

Prepared for
Gardena Farms Irrigation District #13 and
Washington Department of Ecology

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Results of the 2009 Shallow Aquifer Recharge Season at the Locher Road Site, Walla Walla County, Washington

Prepared for

Gardena Farms Irrigation District #13

And

Washington Department of Ecology

Ву

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Executive Summary

This report presents data collected during the 2009 shallow aquifer recharge (SAR) season at the Locher Road Site (the Site) and preliminary interpretations of that data. Work continued to evaluate the feasibility of using SAR to help restore depleted shallow alluvial aquifer groundwater levels and improve flow in spring-fed creeks and streams. SAR at the Site is permitted under a temporary water use authorization granted by the Washington Department of Ecology to Gardena Farms Irrigation District #13 (GFID). This permit authorizes SAR activity for a single season and specifies operating and monitoring conditions for that season. The work done during the 2009 SAR season was done under the third successive permit granted to date.

Given GFID canal operational constraints, observed flow conditions in the Walla Walla River, and the conditions associated with the temporary water use authorization, the 2009 SAR season began on 06 March 2009 and ended on 01 June 2009. Because of flow requirements on the Walla Walla River (both minimum and maximum), SAR activities in the 2009 SAR season occurred in six separate events. A total of approximately 175 acre-feet of water was delivered to the Site and recharged to the alluvial aquifer system during the 56 operational days of the 2009 SAR season.

The alluvial aquifer in the vicinity of the Site does respond to SAR activities, rising and falling as the recharge rate increased and decreased. However, it is not known how far away from the Site the water table response extended during the 2009 SAR season. Water level data collected from off-site wells showed the influence of groundwater pumping which masked the effects of SAR, if any, on the water table more distant from the Site. Following the end of the SAR season, alluvial aquifer water levels at the Site began to fall within one day of the end of operations.

Based on the field and basic water quality parameters collected during the 2009 SAR season, SAR activities at the Site are interpreted to have not degraded groundwater quality in the area of the Site. This data does suggest a high degree of hydraulic continuity between local surface and groundwater, with groundwater at the Site likely influenced by surface water up-gradient of the Site. A few SOC's were detected intermittently. However, the timing of these detections are interpreted to indicate they were not caused by SAR operations and the measured concentrations represent background concentrations related to off-site activities.

It is recommended that SAR at the Site be continued as it does provide recharge to the alluvial aquifer system. Future work should include increasing the size and volume of the recharge basin(s) and conducting currently planned aquifer testing. In addition, every opportunity to deliver water to the Site for SAR should be exploited. This would include several actions being recommended in a draft water plan that GFID will be submitting to the Walla Walla Watershed Partnership in the near future. Finally, given that the alluvial aquifer beneath the Site already receives significant recharge from the Burlingame Canal, and that SOC's are only present rarely and intermittently at extremely low concentrations, it is recommended that SOC sampling be discontinued.

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1.0 - INTRODUCTION

The Locher Road shallow aquifer recharge pilot project (the Project) is one of several projects in the Walla Walla Basin (the Basin) being done to evaluate shallow aquifer recharge (SAR) methods and effects. The objectives of the Project are to gain experience and information on SAR operations, collect information and data to use in evaluating the possible effects of SAR on shallow alluvial aquifer water levels and to contribute recharge water to the alluvial aquifer system. SAR is being explored by water resource stakeholders, in conjunction with other activities in the Basin, to help address water supply, stream flow, water table level, and habitat issues. Data collected from the Locher Road site (the Site) will be used to address these issues in the immediate area of the Site and contribute to basin-wide planning and water resource management efforts.

It is anticipated that pilot SAR activities at the Site will be carried out for several more years, or SAR seasons. Consequently, this report is an interim, or annual, report for 2009 and it does not present final conclusions relative to the project. It is the third annual report produced for the Project to-date.

This report describes the results of the 2009 SAR season, summarizes some basic comparisons between previous seasons and the 2009 SAR season, and presents recommendations for future SAR activities at the Site to be implemented as funding is secured. Site location and basic layout are shown in Figures 1 and 2. As the project progresses over the next several years future project reports will build on the data and recommendations presented in this, previous (GSI, 2007, 2008), and subsequent annual reports.

Descriptions of basic Site physical conditions and infrastructure are found in previous reports (Kennedy/Jenks, 2006; GSI, 2007, 2008). Kennedy/Jenks (2006) also describes: (1) background conditions interpreted for the Site prior to the start of SAR in early 2007 and (2) regulatory constraints under which SAR can be done at the Site. That information will not be repeated in this report.

The data and information described in this report was collected by GSI Water Solutions, Inc. (GSI) as the Project hydrogeologist under contract to Gardena Farms Irrigation District #13 (GFID), by GFID as the Site operator, and by the Walla Walla Basin Watershed Council (WWBWC) as part of its basin-wide groundwater and surface water

data collection and monitoring efforts. This Project is funded by the Washington Department of Ecology (Ecology). The permit authority to operate the Site for the Project was granted by Ecology under a temporary water use authorization issued to GFID.

Topics and information presented in this report include the following:

- A timeline listing the major events associated with the 2009 SAR season.
- Site modifications and changes relative to the 2008 SAR season.
- Rates and volumes of water delivered to the Site in the 2009 SAR season.
- Alluvial aquifer water levels and Mud Creek staff gauge measurements before, during, and after the 2009 SAR season.
- Results of groundwater and surface water quality monitoring for the 2009 SAR season.
- Conclusions and recommendations.

In addition, this report is accompanied by appendices that contain data and information collected during the course of the 2009 SAR season. These appendices are as follows:

• Appendix A. Water quality data, including laboratory QA/QC records.

The Site, including turnouts, control gates, and water distribution was manually operated by the GFID personnel during the 2009 SAR season. GFID staff adjusted flows diverted onto the Site as needed to prevent overtopping of the recharge basins and to meet the conditions and provisions of the temporary water use authorization. GFID staff recorded the time and date of specific actions in field notes, and provided those notes to GSI for use in the preparation of this report. Staff gauge readings also were periodically taken by GFID staff and recorded in project field notes.

WWBWC staff was responsible for water quality sampling before, during, and after the 2009 SAR season, submitting these samples to an analytical laboratory for analysis, and providing analytical results to GSI. WWBWC staff also provided GSI staff with water level data from wells in WWBWC's alluvial aquifer monitoring network that are near the Site.

GSI staff primarily was responsible for data analysis and reporting. In addition, GSI staff attempted to coordinate the activities of GSI, WWBWC, and GFID on the Project.

For the work described herein, the project team included the following people:

- Stuart Durfee GFID Manager, project manager for the Project, and Site operator.
- Kevin Lindsey, Ph.D., L.HG. GSI project manager and hydrogeologist.
- John Fazio, P.E. Fazio Engineering, project engineer, working under subcontract to GSI.
- Jon Travis GSI, technical support.
- Terry Tolan, L.HG. GSI, technical review, hydrogeologist.
- Troy Baker WWBWC, water quality sampling.
- Nella Parks WWBWC, technical support.

The work conducted for this project could not have been done without the cooperation of the Site landowner, Mrs. Patricia Case, the efforts of GFID staff, and the endorsement of the GFID Board of Directors. We thank these people for their support.

2.0 - 2009 TIMELINE

The primary actions associated with the Project during the 2009 SAR season were to operate it and deliver water to the Site (GFID staff primarily responsible), collect monitoring data (WWBWC primarily responsible), and evaluate and interpret monitoring and performance data (GSI staff primarily responsible). Below is a chronological list of basic project activities conducted before, during, and after the 2009 SAR season.

- 28 October 2008 Samples for basic water quality parameters and carbamates collected by WWBWC staff from wells L-1, L-2 and L-3. Source water not collected because Burlingame Canal was not operating.
- 16 December 2008 Samples for basic water quality parameters and SOC's collected by WWBWC staff from wells L-1, L-2 and L-3. Carbamates not analyzed for in well L-1 sample. Source water not collected because Burlingame Canal was not operating.
- 12 February 2009 Samples for basic water quality collected by WWBWC staff from all three wells and source water. Samples for SOC analysis also collected

- by WWBWC staff from all three wells. Source water not collected because Burlingame Canal was not operating.
- 03 March 2009 GSI staff install transducer in flume.
- 05 March 2009 Burlingame Canal operation resumes for spring with opening of head gate on the Walla Walla River near Beet Road by GFID staff.
- 06 March 2009 2009 recharge season begins at 1100 hours. Initial measurement on flume gauge is 0.12 feet or 0.360 cfs.
- 12 March 2009 Samples for basic water quality parameters, carbamates, and herbicides collected by WWBWC staff from all three wells and source water.
 SOC samples not collected.
- 13 March 2009 GSI staff shutdown recharge at 0740 to install transducer in lower basin. Site operations resume at 1135.
- 15 March 2009 GFID shuts down SAR operations.
- 19 March 2009 GFID resumes SAR operations.
- 23 March 2009 GFID shuts down SAR operations.
- 25 March 2009 GFID resumes SAR operations.
- 29 March 2009 GFID shuts down SAR operations.
- 17 April 2009 GFID resumes SAR operations.
- 22 April 2009 GFID shuts down SAR operations.
- 23 April 2009 Samples for basic water quality collected by WWBWC staff from wells L-1 and L-2, source water, and Mud Creek.
- 26 April 2009 GFID resumes SAR operations.
- 6 May 2009 GFID shuts down SAR operations.
- 13 May 2009 GFID resumes SAR operations.
- 28 May 2009 Samples for basic water quality collected by WWBWC staff from wells L-1, L-2 and L-3, source water, and Mud Creek sampling points.
- 01 June 2009 2009 recharge season ends.

3.0 - ON-SITE WORK

No physical modifications were made to the Site for the 2009 SAR season. Work done late in the 2008 SAR season, which focused primarily on increasing the size of the lower basin and deepening the connecting trench between the upper and lower basin, increased the lower basin size approximately four times over its original size (GSI, 2007). For the 2009 SAR season the lower basin had a maximum volume of approximately 50,000 cubic-feet and an approximate wetted surface area of 10,400 square-feet.

4.0 - WATER VOLUME USED IN 2009 SAR SEASON

As was done in the preceding two SAR seasons, the volume of water delivered to the Site during the 2009 SAR season was measured at a ramp flume constructed at the turn-out from Burlingame Canal. A staff gauge for manual readings and a digital pressure transducer for electronic readings were installed in the ramp flume. A conversion chart for the staff gauge (see GSI, 2007) was prepared that allowed the direct conversion of staff gauge readings, in feet, to flow in cfs. Direct staff gauge readings were recorded in field notebooks by GSI and GFID staff during Site visits. The digital transducer was set to record pressure data hourly. The pressure data was converted to the depth of water (height) crossing the measurement sill. These data were subsequently converted to flow estimates using the equation:

 $Q = 10.18 \times h^{1.576}$

Where:

h = height of water above flume measurement sill

Q = flow in cfs through flume

Based on this equation, instantaneous flow onto the Site ranged from approximately 1.0 to 2.5 cfs. The highest measured flow was approximately 5.6 cfs on 12 March 2009. Converting these flow calculations to volume, for the 2009 SAR season approximately 175.8 acre-feet of water (Figure 3) was delivered to the Site during the 56 operating days of the season, for a daily average of approximately 3.14 acre-feet/day.

The 2009 SAR season at the Site is subdivided into six discrete, separate events. The timing and duration of these events was controlled by flow in the Walla Walla River as

measured at the Detour Road gauge. Table 1 summarizes the start and end date, length, average instantaneous flow, and total water volume delivered to the Site for each of these events.

Based on the amount of time it took to drain the lower basin following each of the six events, infiltration rate estimates averaged over the entire lower basin were made. Given that the total surface area of the lower basin is approximately 10,400 square-feet and the water volume and drain time data summarized in Table 2 for each event, infiltration rate on a velocity and unit area basis ranges from 3.8 to 5.3 feet/day and 28 to 39 gallons/square-foot/day, respectively (Table 2). This variability likely reflects such factors as source water temperature (as it relates to viscosity), source water turbidity (as it may influence pore plugging), and water depth in the lower basin (as it may influence driving, pressure, or head).

Comparing water level data discussed in the next section to the known time each SAR event started suggests actual infiltration rates are more rapid than suggested by the average estimates summarized in Table 2. Water levels in the Site monitoring wells generally begin to rise within approximately 12 to 24 hours of the start of each recharge event. Given a depth to water beneath the lower basin of 15 feet, or slightly less, this response time suggests vertical flow velocities through the vadose zone in excess of 30 feet/day. This indicates actual infiltration is not evenly distributed through the substrate underlying the lower basin and that vertical flow along preferential pathways contributes significantly to speeding up infiltration from the basin to the underlying water table.

5.0 - WATER LEVELS DATA

5.1 Water Levels in Wells

As was done in previous SAR seasons, water levels were tracked in on-site monitoring wells L-1, L-2, and L-3. In addition, Ecology staff provided data from a well owned by the agency (referred to in the remainder of this report as the WWGRVL well). Water level data from all four wells was collected using digital transducers.

Although individual water levels differ in each of the Site wells and the WWGRVL well, all four wells display similar trends before, during, and in the case of the Site wells, after the 2009 SAR season (Figure 4). For example:

- In the two months prior to the start of the 2009 SAR season (06 March 2009),
 water levels in the four wells generally rose approximately 1 foot.
- The six separate recharge events comprising the 2009 SAR season are clearly seen in the hydrographs for these wells. In each, water levels began to rise within a few hours of the start of each event (less than 24 hours), and fall within a few hours of the end of each event (less than 24 hours).
- The three Site wells all show water level declines commencing at the end of the 2009 SAR season (01 June 2009) and continuing to the end of the data record for this season (late August 2009). The WWGRVL well data is incomplete and does not show the end of the SAR season.

There is an interesting anomaly in the water level data set. The WWGRVL well commonly displays higher water level elevation than the Site up-gradient well, L-1. The cause of this is unknown, but could be related to such things as an error in the elevation survey, error in the datum correction used to convert depth to water to water table elevation, and/or groundwater mounding related to seepage from the Burlingame Canal.

Off-site water level data was available for the 2009 SAR season from two wells in the Project area that WWBWC personnel had installed transducers in for their basin-wide monitoring program. These two wells are designated by WWBWC as GW 103 and GW 104. These wells are located to the northwest of the site near Frog Hollow Road (Figure 2). Hydrographs for these wells (Figure 5) show water level in GW 103 starting on 15 January 2009 and in GW 104 starting on 26 March 2009. Before testing began the water level at GW 103 was slowly rising and continued to rise after the start of testing on 06 March. GW 104 water level shows a slight rise in water level in early April after the start of testing. Unfortunately, groundwater pumping, in or near both of these wells which began in mid-April, impacts the static level in them. This readily can be seen in the rapid, repeated water level drops on the hydrographs for these wells (Figure 5). These pumping effects make it difficult to assess the impact of SAR operations on those more distant wells after the start of this pumping.

5.2 Mud Creek Stage

Water level, or stage, was also tracked in Mud Creek at one location for the 2009 recharge season; the culvert where Frog Hollow Road crosses the stream. The Frog

Hollow Road site is approximately two miles northwest, and down-gradient, of the Site and it is used to monitor stream levels that may possibly be influenced by Locher Road SAR. Figure 6 shows the stage data recorded by the transducer installed at this location. This data has not been converted to flow volume as a rating curve has not been constructed for this location. In addition, changes in stage have not been checked against such factors as vegetation growth, siltation, and human factors impacting water level. Given these limitations, this stage data is used to provide a general guide to estimate potential relative changes in stream flow before, during, and after the 2009 SAR season.

Mud Creek stage at Frog Hollow Road is variable (Figure 6). Generally it is increasing before the first event of the 2009 SAR season. During the first through fourth events, stage varies, rising and falling. Events 1 and 4 seem to be associated with the greatest stage drops, while events 2 and 3 occur during rises in stage. During and after the fourth event stream stage continues to decline until the start of the fifth event. Stage then begins to increase near the start of the fifth SAR event and continues to rise until SAR shutdown on 01June 2009. Approximately the same time the 2009 SAR season ends, stage also decreases. In addition, the stage decrease seen on 18 April 2009 is coincident with increased groundwater pumping recorded in off-site well GW 104. Given these observations, it is possible, although not certain, that stage changes at Frog Hollow Road might be tied to Locher Road SAR.

6.0 - WATER QUALITY

This section summarizes the results of water quality sampling and analysis done during the 2009 SAR season. Water quality samples were collected from the three Site monitoring wells (L-1, L-2, and L-3), Burlingame Canal at the turn out to the Site, and Mud Creek at Locher Road and at Frog Hollow Road.

6.1 Field and Basic Water Quality

Water quality data for the 2009, 2008, and 2007 SAR seasons are listed in Table 3 and 2009 data is summarized in this section. Complete laboratory results, including laboratory QA records, are reproduced in Appendix A.

For this summary variation in individual parameters during the 2009 SAR season are reviewed individually, with the review focusing on how measured concentrations varied

during the 2009 SAR season. It should be noted, prior to continuing with this review of water quality, that several problems encountered during the SAR season are manifest in the data. Pre-season source water sampling was not done because water was not flowing in the Burlingame Canal on scheduled sampling days. Temperature data was not reported because of problems with field equipment. A post season sampling event was not done because the sampler was unavailable and did not report this fact.

General observations with respect to basic and field water quality during the 2009 SAR season are summarized as follows:

- Pre-season pH for source water was higher than 7.0, while groundwater was less than 7.0 (Figure 7). Mud Creek surface was not sampled during the 2009 SAR season. Groundwater pH was consistently less than 7.0, and showed little variation. This highest groundwater pH during the 2009 SAR season generally was in Well L-1, up-gradient of the Site. The 11.94 pH measured in source water in March 2009 is assumed to be an error, not corrected or rechecked by the sampler while in the field. Source water pH measured later in the 2009 SAR season was similar to groundwater pH, less than 7.0.
- Electrical conductivity (EC) in pre-season groundwater, both up- and down-gradient, was between 125 and 418 micro Siemens per centimeter (mS/cm), with L-1 generally higher than L-2 and L-3 (Figure 8). Source water EC for the 2009 SAR season shows the most variation, although the high measurement on 12 March is, like that for the corresponding pH measurement assumed to be an error. Through the 2009 SAR season wells L-1 and L-2, up-gradient and down-gradient respectively, show the highest groundwater measurements.
- Pre-season hardness values range from a low of 53.7 mg/L in down-gradient well L-3, to a high of 161 mg/L in up-gradient well L-1 (Figure 9). Source water hardness in the pre-season was slightly less than measured in well L-1. During the 2009 SAR the highest hardness seen was in well L-1, which increased from 150 mg/L to 172 mg/L. The lowest hardness seen during the season was in source water, between 20.8 and 37.5 mg/L, and well L-3 between 44.0 and 50.5 mg/L.
- Nitrate-N concentrations in source water and groundwater prior to the start of the
 2009 SAR season was less than 6.08 mg/L with up-gradient well L-1 displaying

the highest concentration (Figure 10). Source water nitrate-N during this period was less than 2.57. During the 2009 SAR season source water and groundwater nitrate-N concentrations in well L-3 were less than 2.05 mg/L. Wells L-1 and L-2 differ from this, displaying a marked increase in nitrate-N concentrations in the April and May samples.

- Total dissolved solids (TDS) in pre-season source and groundwater varied from lows of 145 to 169 mg/L in down-gradient groundwater (well L-3) up to highs of 257 to 259 mg/L in up-gradient groundwater (well L-1) (Figure 11). During the 2009 SAR season well L-3 continued to display the lowest groundwater concentrations while the highest values were shared by wells L-1 and L-2. With the exception of the March 2009 sample, source water hardness of 52 and 61 mg/L were the lowest measured for the season. Although we are not certain, we suspect the March 2009 data point an error, like we suspect for pH and EC.
- The trend seen in the chloride data (Figure 12) closely mimics those seen in the TDS data. Groundwater chloride is highest in well L-1 and L-2 through the preseason and during the season, and lowest in well L-3. Source water is low during the season except for the March 2009 sample which, like it is in pH, EC, and TDS, appears to be anomalously elevated.
- Soluble reactive phosphorus (SRP) concentrations in groundwater and source water ranged between 0.17 and 0.23 mg/L prior of the 2009 SAR season, and between 0.06 and 0.27 mg/L during the season. Before and during the 2009 SAR season source water concentrations generally are less than groundwater concentrations.

Basic water quality parameters summarized above are interpreted to show that SAR activities at the Site did not degrade groundwater quality during the 2009 SAR season. This data, especially the fact that up-gradient groundwater concentrations in most parameters are higher than down-gradient groundwater concentrations and source water, suggests SAR operations at the Site may lead to reductions in parameter concentrations as recharge water is added to the aguifer underlying the Site.

6.2 SOC Water Quality

Three groups of chemicals comprise the SOC analytes for this project: carbonates, SOC's, and herbicides (Table 4). Sampling for these chemicals was done three times prior to the start of the 2009 SAR season, and once following the start of the season. The pre-season sampling events were on 28 October 2008, 19 December 2008, and 12 February 2009. The one sampling event during the season was on 12 March 2009. Sampling and analysis for the three SOC chemical groups on these events were as follows:

- Only carbonates were sampled and analyzed on 28 October 2008 from groundwater.
- All chemical categories were sampled and analyzed for on 16 December 2008 and 12 February 2009 in groundwater. Carbonates in well L-1 were not done.
- For the final sampling event, 12 March 2009, carbonates were sampled for in wells L-2 and L-3 and source water. No sampling was done for SOC's, and herbicides were sampled in all wells and source water.

We have no explanation for this consistency in sampling. Analytical results for this 2009 SAR season sampling is presented in Table 4 and summarized as follows:

- Bromacil was detected, but at extremely low concentrations, in the up-gradient well L-1 in the 16 December 2008 and 12 February 2009 samples. It was not detected in the other wells.
- Atrazine also was detected in well L-1 in the 12 February 2009 sample. It was not detected in the other two wells.
- Dacthal (DCPA) was detected in source water in the 12 March 2009 sample, but not in the wells.

A few SOC and related compounds were detected in up-gradient groundwater and source water before and during the 2009 SAR season. Their inconsistent occurrence, both temporally and spatially, and low concentrations are interpreted to suggest these detections represent very intermittent background conditions and that SAR at the Site is not contributing them to groundwater.

7.0 - COMPARISONS BETWEEN 2007, 2008 AND 2009 SAR SEASONS

This section presents preliminary qualitative comparisons between data collected and observations made during the 2007 and 2008 SAR seasons and the recently completed, 2009 SAR season. In particular:

- The 2009 SAR season was longer than the 2007 and 2008 seasons. Like the 2008 SAR season, it was periodically interrupted because of flow conditions in the Walla Walla River.
- During the 2007 SAR season the recharging capacity of the Site was approximately 1 acre-foot/day. For the 2008 SAR season following the expansion of the lower basin late in the season, average daily infiltration rates increased to approximately 2.4 acre-feet/day. During the 2009 SAR season average daily infiltration rates were 3.3 acre-feet/day. This three-fold increase in recharge capacity corresponded to an approximately four-fold increase in the size of the lower basin between the 2007 and 2009 SAR seasons.
- Water levels in down-gradient wells L-2 and L-3 rose almost 5 feet during the 2009 SAR season, several feet more than the 1.5 to 2.0 foot rise seen during the 2008 and 2007 SAR seasons.
- Field, basic water, and SOC water quality parameters for source water and groundwater during the 2009 SAR season appear to be similar to the 2008 and 2007 seasons. There were concentration fluctuations in basic and field parameters constituents, but no significant changes readily attributed to Site SAR activities.

8.0 - SUMMARY AND RECOMMENDATIONS

8.1 Summary

This report presents the results of the Locher Road 2009 SAR season and preliminary interpretations of some of the data collected to-date. This project is being done in a

continuing effort to evaluate the feasibility of using SAR to help restore depleted shallow alluvial aquifer groundwater levels, improve flow into spring-fed creeks, and increase base flow to streams. SAR at the Site was permitted under a temporary seasonal permit granted by the ECOLOGY. This temporary permit authorizes SAR activity for a single season and specifies operating and monitoring conditions.

The 2009 SAR season began on 06 March 2009 and ended on 01 June 2009. Because of operational restrictions tied to Walla Walla River flow levels specified in the temporary permit, SAR activities in the 2009 SAR season were broken into six separate events. A total of approximately 175.83 acre-feet of water was recharged to the alluvial aquifer during the 56 operational days of the 2009 SAR season.

The alluvial aquifer beneath the Site did respond to SAR activities, rising and falling as the recharge rate increased and decreased. It is not known exactly how far the water table response to 2009 SAR season activities extended away from the Site. Off site wells from which water level data was collected were pumping, or being influenced by pumping wells late in the 2009 SAR season, making an assessment of water table impact distally from the Site problematic. Following the end of the SAR season, water levels in the Site wells began to fall within one to two days of the end of operations.

Based on the field, basic water quality, and SOC parameters measured to-date, SAR activities at the Site are interpreted to have not degraded groundwater quality in the area. This data does suggest a high degree of hydraulic continuity between local surface and groundwater. Much of the impact to groundwater at the Site by surface water appears to be occurring up-gradient of the Site. A few SOC's have been detected intermittently. However, the timing of these detections suggests that they were not caused by SAR operations and the measured concentrations represent background concentrations related to off-site activities.

8.2 Recommendations

Based on the results of the 2009 SAR season described in this report, there are several recommendations for changes to Site operation and monitoring for the 2009/2010 SAR season. These include:

Conduct an aquifer test in a proposed new, purpose-built test well, and use this
data to further constrain aquifer properties and SAR impact on the aquifer in the

- project area. It is our understanding that such a well will be drilled late in 2009 and tested in early 2010.
- Over the past few years the WWBWC, working cooperatively with ECOLOGY, has installed several groundwater monitoring wells in the vicinity of the Site.
 Data from these wells, as was done with GW-103 and GW-104, should be incorporated into the Locher Road SAR project for the 2009/2010 SAR season.
 If done, to the extent possible given operation of these wells by their owners, it would be beneficial to minimize pumping in and near such wells during the SAR season.
- The coordinates and reference point elevation of the WWGRVL well should be done so that an accurate water level elevation may be calculated to determine if it does indeed have an alluvial aquifer water level higher than the Site wells.
- Better coordinate the activities of the consultant with the WWBWC and GFID to prevent data collection oversights, and if they occur, to identify them and quickly rectify them.
- Due to the very low levels of SOC's occurring in source water and in groundwater up-gradient of the site before, during, and after SAR done to-date, it is recommended that water quality monitoring for SOC's be discontinued. The low levels detected in relatively few SOC analytes, suggests that these results reflect background concentrations in both surface water and groundwater unassociated with SAR.

Finally, GFID is nearing completion of a Local Water Plan (LWP) for consideration by the Walla Walla Watershed Partnership. One component of this LWP, the objective of which is to give GFID more flexibility in operation while enhancing in-stream flows, focuses on increased opportunity to operate the Locher Road SAR Site. Acceptance of this portion of the LWP, not to mention the rest of it, would significantly increase the potential for doing SAR at Locher Road as it will give GFID a greater ability to deliver water to the Site.

9.0 - REFERENCES CITED

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GSI, 2007, Results of the first season of shallow aquifer SAR testing at the Locher Road Site, Walla Walla County, Washington. Consultants report prepared for Gardena Farms Irrigation District #13 and Washington Department of Ecology, 23 pgs, 5 tables, 17 figures, 5 Appendices.

GSI, 2008, Results of the 2008 shallow aquifer recharge season at the Locher Road Site, Walla Walla County, Washington. Consultants report prepared for Gardena Farms Irrigation District #13 and Washington Department of Ecology, 10 pgs, 2 tables, 8 figures, 2 Appendices.

Event #	Start Date	End Date	Number of Days	Average Instantaneous Flow (cfs)	Total Volume Delivered (acre-feet)	Reason for Shutdown
1	3/6/2009	3/15/2009	10	1.45	26.85	Walla Walla River over 1000 cfs
2	3/19/2009	3/23/2009	5	1.93	17.40	Walla Walla River over 1000 cfs
3	3/25/2009	3/29/2009	5	1.59	12.89	Walla Walla River over 1000 cfs
4	4/17/2009	4/22/2009	6	1.51	15.85	Walla Walla River over 1000 cfs
5	4/26/2009	5/6/2009	11	1.69	34.37	Walla Walla River over 1000 cfs
6	5/13/2009	6/1/2009	19	1.79	68.47	End of Recharge Season

Table 1. Summary of the six recharge events during the 2009 SAR season.

Recharge Event #	Estimated Volume of Water in Basin in cu-ft	Estimated Time to Drain	Infiltration Rate in ft/day	Infiltration rate by unit area (gal/sq-
	and gal.	Basin in min		ft/day)
1	29,010/216,995	790	5.08	37.4
2	44,083/329,741	1330	4.4	34.6
3	45,632/341,327	1380	4.58	34.6
4	23,022/172,204	840	3.79	28.8
5	32,727/244,798	1110	4.08	30.2
6	47,180/352,906	1250	5.22	38.9

Table 2. Estimated infiltration volumes and rates at the conclusion of each of the six recharge events during the 2009 SAR season.

Sample ID	Date	Lab No.	рН	Town C	Electrical Conductivity (mS/cm)	Turbidity (NTU)	NO ₃ -N	Hardness	TDS	CI (mg/l)	Soluble Reactive Phosphorous	S	COD (mg/L)	Total Coliform (per 100ml)	E-Coli (per 100ml or Absent/Pres
L-1	10/12/2006	85232	рп	Temp. C	(1113/6111)	(NTO)	(mg/L) 6.23	(mg/L) 205.0	(mg/L) 262.0	CI (mg/L) 7.8	(mg/L) 0.100		8.0	(per room)	ent) 0
L-1	1/15/2007	86451	6.77	12.8	432	0.15	6.50	205.0	238.0	1.2	0.100	<	8.0	0	
L-1	4/4/2007	87538	7.24	13.8	401	0.13	5.68	217.0	253.0		< 0.043	<	8.0	0	0
L-1	4/12/2007	87725	7.24	13.5	393	1.92	5.19	213.0	248.0	6.5	0.043	_ <	8.0	0	0
L-1	4/23/2007	87918	7.17	13.9	428	0.42	4.20	624.0	247.0	7.5	0.090	_ <	8.0	0	0
L-1	1/23/2007	2123	7.07	10.0	388	48.6	5.28	135.0	235.0	5.8	0.360		8.0	0	Absent
L-1	2/13/2008	4099	7.13		384	44.4	5.55	154.0	236.0	5.6	0.360		8.0	0	Absent
L-1	5/27/2008	15124	7		401	0.79	5.86	156.0	262.0	6.7	0.270		ND		71000111
 L-1	10/28/2008	32780	6.81		403	0.29	6.38	158.1	265.0	3.5	0.240		ND	absent	Absent
L-1	12/16/2008	37230	6.91		396	0.7	5.85	150.0	257.0	6.1	0.190		ND	absent	Absent
 L-1	2/12/2009	4470	6.78		418	1.78	6.08	161.0	259.0	6.7	0.230		ND	absent	Absent
L-1	3/12/2009	7335	6.72		335	8.12	5.21	150.0	225.0	5.9	0.220		ND	absent	Absent
L-1	4/23/2009	11906	6.71		434	15.5	16.20	167.0	279.0	9.7	0.210		ND	absent	Absent
L-1	5/28/2009	16095	6.89		431	1.3	7.07	172.0	266.0	7.3	0.270		ND	absent	Absent
L-2	10/12/2006	85233					3.27	132.0	184.0	6.4	0.140	<	8.0		0
L-2	1/15/2007	86452	7.05	11.9	281	0.67	3.63	117.0	154.0	0.8	0.130	<	4.0	0	
L-2	4/4/2007	87539	7.19	13.0	284	0.39	4.12	145.0	190.0		< 0.043	<	8.0	0	0
L-2	4/12/2007	87726	7.17	13.0	284	0.65	3.62	148.0	148.0	0.3	0.080	<	8.0	0	0
L-2	4/23/2007	87919	7.17	13.2	288	0.64	1.34	134.0	180.0	4.5	0.140	<	8.0	0	0
L-2	1/23/2008	2124	7.04		287	8.06	3.47	103.0	181.0	5.6	0.330	<	8.0	0	Present
L-2	2/13/2008	4100	7.07		284	8.65	3.50	111.7	196.0	5.5	0.330	<	8.0	0	Present
L-2	5/27/2008	15126	6.93		313	7.13	5.96	117.0	205.0	6.5	0.270		10.0		
L-2	10/28/2008	32781	6.73		291	1.79	4.25	116.0	201.0	6.2	0.240		ND	absent	absent
L-2	12/16/2008	37231	6.87		284	2.03	3.56	105.0	184.0	5.2	0.190		ND	present	absent
L-2	2/12/2009	4471	6.77		301	0.76	4.25	109.0	193.0	5.8	0.230		ND	absent	absent
L-2	3/12/2009	7336	6.59		295	2.02	5.42	118.0	196.0	6.2	0.220		ND	absent	absent
L-2	4/23/2009	11905	6.45		440	10.9	17.30	118.0	302.0	9.8	0.200		ND	absent	absent
L-2	5/28/2009	16096	6.69		422	9.13	15.40	164.0	286.0	8.9	0.260		ND	present	absent
L-3	10/12/2006			-	<u> </u>	·	·	<u> </u>						·	
L-3	1/15/2007	86453	6.88	10.1	202	25.00	2.86	83.4	118.0	< 0.3	0.130	<	8.0	0	
L-3	4/4/2007	87540	7.47	9.6	104	2.51	0.81	54.5	92.5		< 0.043	<	8.0	0	0
L-3	4/12/2007	87727	7.36	9.4	126	2.57	0.90	61.8	92.5	< 0.3	0.070	<	8.0	present	0
L-3	4/23/2007	87920	7.35	9.8	135	1.97	0.47	54.7	86.7	< 0.3	0.100	<	8.0	0	0
L-3	1/23/2008	2125	7.14		187	13.90	2.86	61.9	124.0	3.2	0.300		18.0	0	present
L-3	2/13/2008	4101	7.15		197	29.40	4.51	75.3	148.0	5.5	0.320	<	8.8	0	present
L-3	5/27/2008	15127	6.98		129	7.48	2.11	45.7	98.0	1.8	0.220		16.0		
L-3	10/28/2008	32782	6.75		122	3.61	0.43	46.1	94.0	1.9	0.190		ND	present	absent
L-3	12/16/2008	37232	6.91		125	0.11	1.75	53.7	145.0	2.5	0.170		ND	absent	absent
L-3	2/12/2009	4472	6.81		231	1.62	5.63	84.4	169.0	4.2	0.210		ND	absent	absent
L-3	3/12/2009	7337	6.61		136	2.01	2.05	50.5	91.0	2.2	0.170		ND	absent	absent
L-3	5/28/2009	16097	6.8		121	6.88	1.84	44.0	96.0	1.4	ND		ND	present	present

Table 3. Field and basic water quality results for 2006-2009

Sample ID	Date	Lab No.	рН	Temp. C	Electrical Conductivity (mS/cm)	Turbidity (NTU)	NO ₃ -N (mg/L)	Hardness (mg/L)	TDS (mg/L)	CI (mg/L)	Soluble Reactive Phosphorous (mg/L)	s	COD (mg/L)	Total Coliform (per 100ml)	E-Coli (per 100ml or Absent/Pres ent)
Mud Ck - L	10/12/2006															
Mud Ck - L	1/15/2007	86454	6.21	1.8	262	1.39	2.18	112.0	144.0	2.0		0.060	<	8.0	0	
Mud Ck - L	4/4/2007	87542	8.05	11.2	242	3.48	1.30	132.0	158.0	10.0	<	0.043		8.0	present	present
Mud Ck - L	4/12/2007	87729	7.70	9.6	173	1.49	0.40	95.0	118.0	< 0.3		0.100		8.0	present	present
Mud Ck - L	4/23/2007	87922	8.01	15.3	181	1.26 <		83.1	117.0	< 0.3		0.060	<	8.0	present	present
Mud Ck - L	5/27/2008	15129	7.24	.0.0	162	4.95	0.57	65.0	112.0	2.9		0.180		11.0	р.осол.	ртоссти
Mud Ck - L	3/12/2009	7340	7.62		373	2.68	2.59	142.0	216.0	7.5		0.150		ND	present	present
Mud Ck - L	4/23/2009	11903	6.94		275	0.5	0.98	114.0	163.0	5.3		0.130		ND	present	present
Mud Ck - L	5/28/2009	16099	7.25		310	0.96	1.38	126.0	198.0	6.2		0.660		ND	present	present
Mud Ck - FH	10/12/2006 1/15/2007 4/4/2007 4/12/2007 4/12/2007 4/23/2007 5/27/2008 3/12/2009 4/23/2009 5/28/2009	86455 87541 87728 87921 15130 7339 11902 16098	6.13 8.28 7.89 8.16 7.50 7.69 7.07 7.18	2.8 12.6 10.9 16.5	268 248 175 180 298 272 363 374	3.56 1.81 1.89 1.89 4.11 0.64 1.04 3.97	2.17 1.39 0.52 0.47 0.87 1.82 1.56 1.37	113.0 130.0 95.5 82.4 112.0 112.0 149.0	146.0 165.0 123.0 113.0 188.0 179.0 220.0 226.0	0.8 11.0 0.7 5.0 4.6 5.5 7.4 6.8	<	0.050 0.043 0.040 0.080 0.240 0.150 0.160 0.260	<	9.0 8.0 9.0 10.0 21.0 ND ND	o present present present present present present	present present present present present present present
diversion diversion diversion diversion	10/12/2006 1/15/2007 4/4/2007 4/12/2007	87543 87730	8.02 7.77	9.4 8.0	95 90	6.28 4.27	0.38 0.12	45.2 44.1	95.0 65.0	40.0	<	0.043	<	8.0 12.0	absent present	absent present
	4/23/2007	87923	8.17	12.7	94	6.39	4.21	31.4	73.3	39.0	<	0.043	<	8	present	present
diversion	5/27/2008	15128	7.27	12.1	50.3	17.6	0.11	18.7	75.5 54	0.6		0.12		19	prosont	prosent
diversion diversion					378	0.83	2.57	141.0	224	7.2		0.18		ND	present	absent
diversion		4473	7.63													
diversion diversion	2/12/2009	7338	7.63 11.94					37.5	826	22.0						present
diversion		4473 7338 11904	7.63 11.94 6.52		2560 52.9	21.7 61.8	0.58 0.22	37.5 20.8	826 52	22.0 0.7		0.06 0.12		ND 10	present	present present

Table 3. Field and basic water quality results for 2006-2009 (cont.)

Date	1/15/2007	1/15/2007	1/15/2007
Well ID Chemical	L-1	L-2	L-3
Carbamates i	in Drinking w	ater	
Carbofuran	ND	ND	ND
Oxymal	ND	ND	ND
3-Hydroxycabofuran	ND	ND	ND
Aldicarb Aldicarb sulfone	ND ND	ND ND	ND ND
Aldicarb sulfoxide	ND	ND	ND
Carbaryl	ND	ND	ND
Methomyl	ND	ND	ND
Propoxur (Baygon)	ND	ND	ND
Methiocarb	ND	ND	ND
Synthetic Org Endrin	yanic Compoi ND	unas ND	ND
Lindane (BHC-Gamma)	ND ND	ND	ND ND
Methoxychlor	ND	ND	ND
Alachlor	ND	ND	ND
Atrazine	ND	ND	ND
Benzo(a)pyrene	ND	ND	ND
Chlordane Technical	ND	ND	ND
Di(ethylhexyl)-Adipate	ND ND	ND ND	ND
Di(ethylhexyl)-phthalate Heptachlor	ND ND	ND	ND ND
Heptachlor Epoxide A&B	ND	ND	ND
Hexachlorobenzene	ND	ND	ND
Hexachlorocyclo-Pentadiene	ND	ND	ND
Simazine	ND	ND	ND
Aldrin	ND	ND	ND
Butachlor	ND ND	ND ND	ND ND
Dieldrin Metolachlor	ND ND	ND ND	ND ND
Metribuzin	ND	ND	ND
Propachlor	ND	ND	ND
Bromacil	0.74	ND	ND
Prometon	ND	ND	ND
Terbacil	ND	ND	ND
Diazinon EPTC	ND ND	ND ND	ND ND
4,4-DDD	ND ND	ND	ND ND
4,4-DDE	ND	ND	ND
4,4-DDT	ND	ND	ND
Cyanazine	ND	ND	ND
Malathion	ND	ND	ND
Trifluralin	ND	ND	ND
Napthalene	ND ND	ND	ND ND
Fluorene Acenaphthylene	ND ND	ND ND	ND
Acenaphthene	ND	ND	ND
Anthracene	ND	ND	ND
Benz(A)anthracene	ND	ND	ND
Benzo(B)fluoranthene	ND	ND	ND
Benzo(G,H,I)peryene	ND	ND	ND
Benzo(K)fluoranthene Chrysene	ND ND	ND ND	ND ND
Dibenzo(A,H)anthracene	ND	ND	ND
Fluoranthene	ND	ND	ND
Indeno(1,2,3-CD)pyrene	ND	ND	ND
Phenanthrene	ND	ND	ND
Pyrene	ND	ND	ND
Benzyl Butyl Phthalate	ND 0.05	ND	ND
Di-N-Butyl Phthalate Diethyl Phthalate	0.95 ND	ND ND	ND ND
Dietnyl Phthalate	ND ND	ND ND	ND ND
Toxaphene	ND	ND	ND
Aroclor 1221	ND	ND	ND
Aroclor 1232	ND	ND	ND
Aroclor 1242	ND	ND	ND
Aroclor 1248	ND	ND	ND
Aroclor 1254 Aroclor 1260	ND ND	ND ND	ND ND
Aroclor 1016	ND ND	ND	ND ND
	n Drinking Wa		.10
2,4-D	ND	ND	ND
2,4,5-TP (Silvex)	ND	ND	ND
Pentachlorophenol	ND	ND	ND
Dalapon	ND	ND	ND
Dinoseb	ND	ND	ND
Picloram Dicamba	ND	ND	ND
	ND	ND ND	ND ND
	NID		ND
2,4 DB	ND ND		ΝD
2,4 DB 2,4,5 T	ND	ND	ND ND
2,4 DB			ND ND ND
2,4 DB 2,4,5 T Bentazon Dichlorprop Actiflorfin	ND ND ND ND	ND ND ND ND	ND ND ND
2,4 DB 2,4,5 T Bentazon Dichlorprop	ND ND ND	ND ND ND	ND ND

Date	4/4/2007	4/4/2007	4/4/2007	4/4/2007
Well ID Chemical	Diversion	L-1	L-2	L-3
	nates in Drink	cing water		
Carbofuran	ND	ND	ND	ND
Oxymal	ND	ND	ND	ND
3-Hydroxycabofuran	ND	ND	ND	ND
Aldicarb	ND	ND	ND	ND
Aldicarb sulfone	ND	ND	ND	ND
Aldicarb sulfoxide	ND	ND	ND	ND
Carbaryl	ND	ND	ND	ND
Methomyl	ND	ND	ND	ND
Propoxur (Baygon)	ND	ND	ND	ND
Methiocarb	ND	ND	ND	ND
	tic Organic Co		ND	ND
Endrin Lindane (BHC-Gamma)	ND ND	ND ND	ND ND	ND ND
	ND	ND	ND	ND
Methoxychlor Alachlor	ND	ND	ND	ND
Atrazine	ND	ND	ND	ND
Benzo(a)pyrene	ND	ND	ND	ND
Chlordane Technical	ND	ND	ND	ND
Di(ethylhexyl)-Adipate	ND	ND	ND	ND
Di(ethylhexyl)-phthalate	ND	ND	ND	ND
Heptachlor	ND	ND	ND	ND
Heptachlor Epoxide A&B	ND	ND	ND	ND
Hexachlorobenzene	ND	ND	ND	ND
Hexachlorocyclo-Pentadiene	ND	ND	ND	ND
Simazine	ND	ND	ND	ND
Aldrin	ND	ND	ND	ND
Butachlor	ND	ND	ND	ND
Dieldrin	ND	ND	ND	ND
Metolachlor	ND	ND	ND	ND
Metribuzin	ND	ND	ND	ND
Propachlor	ND	ND	ND	ND
Bromacil	ND	0.2	ND	ND
Prometon	ND	ND	ND	ND
Terbacil	ND	ND	ND	ND
Diazinon	ND	ND	ND	ND
EPTC	ND	ND	ND	ND
4,4-DDD	ND	ND	ND	ND
4,4-DDE	ND	ND	ND	ND
4,4-DDT	ND	ND	ND	ND
Cyanazine	ND	ND	ND	ND
Malathion	ND	0.4	0.5	0.3
Trifluralin	ND	ND	ND	ND
Napthalene	ND ND	ND	ND ND	ND ND
Fluorene Acenaphthylene	ND ND	ND ND	ND ND	ND ND
	ND	ND	ND	ND
Acenaphthene Anthracene	ND	ND	ND	ND
Benz(A)anthracene	ND	ND	ND	ND
Benzo(B)fluoranthene	ND	ND	ND	ND
Benzo(G,H,I)peryene	ND	ND	ND	ND
Benzo(K)fluoranthene	ND	ND	ND	ND
Chrysene	ND	ND	ND	ND
Dibenzo(A,H)anthracene	ND	ND	ND	ND
Fluoranthene	ND	ND	ND	ND
Indeno(1,2,3-CD)pyrene	ND	ND	ND	ND
Phenanthrene				
Pyrene	ND	ND	ND	ND
Benzyl Butyl Phthalate	ND	ND	ND	ND
Di-N-Butyl Phthalate	ND	0.7	0.7	0.5
Diethyl Phthalate	ND	ND	ND	ND
Dimethyl Phthalate	ND	ND	ND	ND
Toxaphene	ND	ND	ND	ND
Aroclor 1221	ND	ND	ND	ND
Aroclor 1232	ND	ND	ND	ND
Aroclor 1242	ND	ND	ND	ND
Aroclor 1248	ND	ND	ND	ND
Aroclor 1254	ND	ND	ND	ND
Aroclor 1260	ND	ND	ND	ND
Aroclor 1016	ND	ND	ND	ND
	ides in Drink	ing Water		
2,4-D				
2,4,5-TP (Silvex)	ND	ND	ND	ND
Pentachlorophenol	ND	ND	ND	ND
	ND	ND	ND	ND
Dalapon	ND	ND	ND	ND
Dinoseb	ND		ND	ND
Dinoseb Picloram	ND	ND		
Dinoseb Picloram Dicamba	ND ND	ND	ND	ND
Dinoseb Picloram Dicamba 2,4 DB	ND ND ND	ND ND	ND ND	ND
Dinoseb Picloram Dicamba 2,4 DB 2,4,5 T	ND ND ND ND	ND ND ND	ND ND ND	ND ND
Dinoseb Picloram Dicamba 2,4 DB 2,4,5 T Bentazon	ND ND ND ND	ND ND ND ND	ND ND ND ND	ND ND ND
Dinoseb Picloram Dicamba 2,4 DB 2,4,5 T Bentazon Dichlorprop	ND ND ND ND ND	ND ND ND ND	ND ND ND ND	ND ND ND ND
Dinoseb Picloram Dicamba 2,4 DB 2,4,5 T Bentazon	ND ND ND ND	ND ND ND ND	ND ND ND ND	ND ND ND

Date	2/13/2008	2/13/2008	2/13/2008
Well ID Chemical	L-1	L-2	L-3
	in Drinking w	ater	
Carbofuran	ND	ND	ND
Oxymal 2 Hydroxygabafyran	ND ND	ND ND	ND ND
3-Hydroxycabofuran Aldicarb	ND	ND	ND
Aldicarb sulfone	ND	ND	ND
Aldicarb sulfoxide Carbaryl	ND ND	ND ND	ND ND
Methomyl	ND	ND	ND
Propoxur (Baygon)	ND	ND	ND
Methiocarb Synthetic Org	ND panic Compo	ND unds	ND
Endrin	ND	ND	ND
Lindane (BHC-Gamma)	ND	ND	ND
Methoxychlor Alachlor	ND ND	ND ND	ND ND
Atrazine	ND	ND	ND
Benzo(a)pyrene	ND	ND	ND
Chlordane Technical Di(ethylhexyl)-Adipate	ND ND	ND ND	ND ND
Di(ethylhexyl)-phthalate	ND	ND	ND
Heptachlor	ND	ND	ND
Heptachlor Epoxide A&B Hexachlorobenzene	ND ND	ND ND	ND ND
Hexachlorocyclo-Pentadiene	ND	ND	ND
Simazine	ND	ND	ND
Aldrin Butachlor	ND ND	ND ND	ND ND
Dieldrin	ND	ND	ND
Metolachlor	ND	ND	ND
Metribuzin	ND	ND	ND
Propachlor Bromacil	ND 0.32	ND ND	ND ND
Prometon	ND	ND	ND
Terbacil	ND	ND	ND
Diazinon EPTC	ND ND	ND ND	ND ND
4,4-DDD	ND ND	ND ND	ND ND
4,4-DDE	ND	ND	ND
4,4-DDT	ND	ND	ND
Cyanazine Malathion	ND ND	ND ND	ND ND
Trifluralin	ND	ND	ND
Napthalene	ND	ND	ND
Fluorene Acenaphthylene	ND ND	ND ND	ND ND
Acenaphthene	ND	ND	ND
Anthracene	ND	ND	ND
Benz(A)anthracene Benzo(B)fluoranthene	ND ND	ND ND	ND ND
Benzo(G,H,I)peryene	ND	ND	ND
Benzo(K)fluoranthene	ND	ND	ND
Chrysene	ND	ND	ND
Dibenzo(A,H)anthracene Fluoranthene	ND ND	ND ND	ND ND
Indeno(1,2,3-CD)pyrene	ND	ND	ND
Phenanthrene	ND	ND	ND
Pyrene Benzyl Butyl Phthalate	ND ND	ND ND	ND ND
Di-N-Butyl Phthalate	ND	ND	ND
Diethyl Phthalate	ND	ND	ND
Dimethyl Phthalate Toxaphene	ND ND	ND ND	ND ND
Aroclor 1221	ND	ND ND	ND
Aroclor 1232	ND	ND	ND
Aroclor 1242	ND	ND	ND
Aroclor 1248 Aroclor 1254	ND ND	ND ND	ND ND
Aroclor 1260	ND	ND	ND
Aroclor 1016	ND	ND	ND
2,4-D	n Drinking Wa	ND ND	ND
2,4,5-TP (Silvex)	ND	ND	ND
Pentachlorophenol	ND	ND	ND
Dalapon Dinoseb	ND ND	ND	ND
Picloram	ND ND	ND ND	ND ND
Dicamba	ND	ND	ND
2,4 DB	ND	ND	ND
2,4,5 T Bentazon	ND ND	ND ND	ND ND
Dichlorprop	ND	ND	ND
Actiflorfin	ND	ND	ND
Dacthal (DCPA)	ND	ND	ND
3,5-Dichlorobenzoic Acid	ND	ND	ND

Date			5/27/2008	5/27/2008	5/27/2008	5/27/200
Well ID	Diversion	L-1	L-2	L-3	MC-L	MC-SL
Chemical Carbar	mates in Drinki	ng water			-	
Carbofuran	ND	ND	ND	ND	ND	ND
Oxymal	ND	ND	ND	ND	ND	ND
3-Hydroxycabofuran	ND	ND	ND	ND	ND	ND
Aldicarb	ND	ND	ND	ND	ND	ND
Aldicarb sulfone	ND	ND	ND	ND	ND	ND
Aldicarb sulfoxide Carbaryl	ND ND	ND ND	ND ND	ND ND	ND ND	ND ND
Methomyl	ND ND	ND	ND	ND	ND ND	ND
Propoxur (Baygon)	ND	ND	ND	ND	ND	ND
Methiocarb	ND	ND	ND	ND	ND	ND
	tic Organic Co					
Endrin	ND	ND	ND	ND	ND	ND
Lindane (BHC-Gamma)	ND	ND	ND	ND	ND	ND
Methoxychlor Alachlor	ND ND	ND ND	ND ND	ND ND	ND ND	ND ND
Atrazine	ND	ND	ND	ND	ND	ND
Benzo(a)pyrene	ND	ND	ND	ND	ND	ND
Chlordane Technical	ND	ND	ND	ND	ND	ND
Di(ethylhexyl)-Adipate	ND	ND	ND	ND	ND	ND
Di(ethylhexyl)-phthalate	ND	ND	ND	ND	ND	ND
Heptachlor	ND	ND	ND	ND	ND	ND
Heptachlor Epoxide A&B Hexachlorobenzene	ND ND	ND ND	ND ND	ND ND	ND ND	ND ND
Hexachlorocyclo-Pentadiene	ND ND	ND	ND ND	ND ND	ND	ND ND
Simazine	ND	ND	ND	ND	ND	ND
Aldrin	ND	ND	ND	ND	ND	ND
Butachlor	ND	ND	ND	ND	ND	ND
Dieldrin	ND	ND	ND	ND	ND	ND
Metolachlor	ND	ND	ND	ND	ND	ND
Metribuzin Propachlor	ND ND	ND ND	ND ND	ND ND	ND ND	ND ND
Bromacil	ND	ND	ND	ND	ND	ND
Prometon	ND	ND	ND	ND	ND	ND
Terbacil	ND	ND	ND	ND	ND	ND
Diazinon	ND	ND	ND	ND	ND	ND
EPTC	ND	ND	ND	ND	ND	ND
4,4-DDD	ND	ND	ND	ND	ND	ND
4,4-DDE 4,4-DDT	ND	ND	ND ND	ND ND	ND ND	ND ND
Cyanazine	ND ND	ND ND	ND ND	ND	ND ND	ND
Malathion	ND	ND	ND	ND	ND	ND
Trifluralin	ND	ND	ND	ND	ND	ND
Napthalene	ND	ND	ND	ND	ND	ND
Fluorene	ND	ND	ND	ND	ND	ND
Acenaphthylene	ND	ND	ND	ND	ND	ND
Acenaphthene Anthracene	ND ND	ND ND	ND ND	ND ND	ND ND	ND ND
Benz(A)anthracene	ND ND	ND	ND	ND	ND	ND
Benzo(B)fluoranthene	ND	ND	ND	ND	ND	ND
Benzo(G,H,I)peryene	ND	ND	ND	ND	ND	ND
Benzo(K)fluoranthene	ND	ND	ND	ND	ND	ND
Chrysene	ND	ND	ND	ND	ND	ND
Dibenzo(A,H)anthracene	ND	ND	ND	ND	ND	ND
Fluoranthene	ND ND	ND ND	ND ND	ND ND	ND ND	ND
Indeno(1,2,3-CD)pyrene Phenanthrene	ND ND	ND ND	ND ND	ND ND	ND ND	ND ND
Pyrene	ND	ND	ND	ND	ND	ND
Benzyl Butyl Phthalate	ND	ND	ND	ND	ND	ND
Di-N-Butyl Phthalate	ND	ND	ND	ND	ND	ND
Diethyl Phthalate	ND	ND	ND	ND	ND	ND
Dimethyl Phthalate	ND	ND	ND	ND	ND	ND
Toxaphene Aroclor 1221	ND ND	ND ND	ND ND	ND ND	ND ND	ND ND
Aroclor 1232	ND ND	ND	ND	ND ND	ND	ND
Aroclor 1242	ND	ND	ND	ND	ND	ND
Aroclor 1248	ND	ND	ND	ND	ND	ND
Aroclor 1254	ND	ND	ND	ND	ND	ND
Aroclor 1260	ND	ND	ND	ND	ND	ND
Aroclor 1016	ND	ND	ND	ND	ND	ND
	cides in Drinkin	ig Water				
2,4-D 2,4,5-TP (Silvex)	ND	ND	ND	ND	NID	NID
Pentachlorophenol	ND ND	ND ND	ND ND	ND ND	ND ND	ND ND
Dalapon	ND	ND	ND	ND	ND	ND
Dinoseb	ND	ND	ND	ND	ND	ND
Picloram	ND	ND	ND	ND	ND	ND
Dicamba	ND	ND	ND	ND	ND	0.18
2,4 DB	ND	ND	ND	ND	ND	ND
2,4,5 T	ND	ND	ND	ND	ND	ND
		ND	ND	ND	ND	ND
Bentazon	ND				N/PS	LIP.
Bentazon Dichlorprop	ND	ND	ND	ND	ND	ND
Bentazon					ND ND 0.4	ND ND 0.2

Date		10/28/2008	
Well ID	L-1	L-2	L-3
Chemical	in Drinking w	otor	
Carbofuran	in Drinking wa	ND ND	ND
Oxymal	ND	ND ND	ND
3-Hydroxycabofuran	ND	ND	ND
Aldicarb	ND	ND	ND
Aldicarb sulfone	ND	ND	ND
Aldicarb sulfoxide	ND	ND	ND
Carbaryl	ND	ND	ND
Methomyl	ND	ND	ND
Propoxur (Baygon)	ND	ND	ND
Methiocarb	ND	ND	ND
Endrin	ganic Compou NS	NS	NS
Lindane (BHC-Gamma)	NS	NS	NS
Methoxychlor	NS	NS	NS
Alachlor	NS	NS	NS
Atrazine	NS	NS	NS
Benzo(a)pyrene	NS	NS	NS
Chlordane Technical	NS	NS	NS
Di(ethylhexyl)-Adipate	NS	NS	NS
Di(ethylhexyl)-phthalate	NS	NS	NS
Heptachlor	NS	NS	NS
Heptachlor Epoxide A&B	NS	NS	NS
Hexachlorobenzene	NS	NS	NS
Hexachlorocyclo-Pentadiene	NS	NS NS	NS
Simazine Aldrin	NS NS	NS NS	NS NS
Butachlor	NS	NS NS	NS
Dieldrin	NS	NS NS	NS
Metolachlor	NS	NS	NS
Metribuzin	NS	NS	NS
Propachlor	NS	NS	NS
Bromacil	NS	NS	NS
Prometon	NS	NS	NS
Terbacil	NS	NS	NS
Diazinon	NS	NS	NS
EPTC	NS	NS	NS
4,4-DDD	NS	NS	NS
4,4-DDE	NS	NS	NS
4,4-DDT	NS	NS	NS
Cyanazine	NS	NS NS	NS
Malathion Trifluralin	NS NS	NS NS	NS NS
Napthalene	NS	NS	NS
Fluorene	NS	NS	NS
Acenaphthylene	NS	NS	NS
Acenaphthene	NS	NS	NS
Anthracene	NS	NS	NS
Benz(A)anthracene	NS	NS	NS
Benzo(B)fluoranthene	NS	NS	NS
Benzo(G,H,I)peryene	NS	NS	NS
Benzo(K)fluoranthene	NS	NS	NS
Chrysene	NS	NS	NS
Dibenzo(A,H)anthracene	NS	NS	NS
Fluoranthene	NS	NS	NS
Indeno(1,2,3-CD)pyrene	NS NS	NS NS	NS
Phenanthrene Pyrene	NS NS	NS NS	NS NS
Benzyl Butyl Phthalate	NS	NS NS	NS
Di-N-Butyl Phthalate	NS	NS	NS
Diethyl Phthalate	NS	NS	NS
Dimethyl Phthalate	NS	NS	NS
Toxaphene	NS	NS	NS
Aroclor 1221	NS	NS	NS
Aroclor 1232	NS	NS	NS
Aroclor 1242	NS	NS	NS
Aroclor 1248	NS	NS	NS
Aroclor 1254	NS	NS	NS
Aroclor 1260	NS	NS	NS
Aroclor 1016	NS n Drinking Wa	NS	NS
	n Drinking Wa		NO
2,4-D 2,4,5-TP (Silvex)	NS NS	NS NS	NS NS
	NS NS	NS NS	NS NS
Pentachlorophenol Dalapon	NS NS	NS NS	NS NS
Dinoseb	NS	NS	NS
Picloram	NS	NS NS	NS
Dicamba	NS	NS	NS
2,4 DB	NS	NS	NS
2,4,5 T	NS	NS	NS
Bentazon	NS	NS	NS
Dichlorprop	NS	NS	NS
Actiflorfin	NS	NS	NS
Dacthal (DCPA)	NS	NS	NS

	10/10/0000		10/10/0000
Date Well ID	12/16/2008 ·	12/16/2008 L-2	12/16/2008 L-3
Chemical			
Carbamates	in Drinking wa	ter	
Carbofuran	NS	ND	ND
Oxymal	NS NS	ND	ND
3-Hydroxycabofuran Aldicarb	NS NS	ND ND	ND ND
Aldicarb sulfone	NS	ND	ND
Aldicarb sulfoxide	NS	ND	ND
Carbaryl	NS	ND	ND
Methomyl Propoxur (Baygon)	NS NS	ND ND	ND ND
Methiocarb	NS	ND	ND
Synthetic Or	ganic Compou	nds	
Endrin	ND	ND	ND
Lindane (BHC-Gamma)	ND	ND	ND
Methoxychlor Alachlor	ND ND	ND ND	ND ND
Atrazine	ND	ND	ND
Benzo(a)pyrene	ND	ND	ND
Chlordane Technical	ND	ND	ND
Di(ethylhexyl)-Adipate Di(ethylhexyl)-phthalate	ND ND	ND ND	ND ND
Heptachlor	ND	ND	ND
Heptachlor Epoxide A&B	ND	ND	ND
Hexachlorobenzene	ND	ND	ND
Hexachlorocyclo-Pentadiene Simazine	ND	ND	ND ND
Aldrin	ND ND	ND ND	ND ND
Butachlor	ND	ND	ND
Dieldrin	ND	ND	ND
Metolachlor	ND	ND	ND
Metribuzin Propachlor	ND ND	ND ND	ND ND
Bromacil	0.09 ug/L J	ND	ND
Prometon	ND	ND	ND
Terbacil	ND	ND	ND
Diazinon	ND	ND	ND
EPTC 4.4-DDD	ND	ND	ND
4,4-DDD 4,4-DDE	ND ND	ND ND	ND ND
4,4-DDT	ND	ND	ND
Cyanazine	ND	ND	ND
Malathion	ND	ND	ND
Trifluralin	ND	ND	ND
Napthalene Fluorene	ND ND	ND ND	ND ND
Acenaphthylene	ND	ND	ND
Acenaphthene	ND	ND	ND
Anthracene	ND	ND	ND
Benz(A)anthracene	ND ND	ND ND	ND ND
Benzo(B)fluoranthene Benzo(G,H,I)peryene	ND	ND	ND
Benzo(K)fluoranthene	ND	ND	ND
Chrysene	ND	ND	ND
Dibenzo(A,H)anthracene	ND	ND	ND
Fluoranthene	ND ND	ND ND	ND ND
Indeno(1,2,3-CD)pyrene Phenanthrene	ND ND	ND	ND ND
Pyrene	ND	ND	ND
Benzyl Butyl Phthalate	ND	ND	ND
Di-N-Butyl Phthalate	ND	ND	ND
Diethyl Phthalate	ND ND	ND ND	ND ND
Dimethyl Phthalate Toxaphene	ND ND	ND ND	ND ND
Aroclor 1221	ND	ND	ND
Aroclor 1232	ND	ND	ND
Aroclor 1242	ND	ND	ND
Aroclor 1248 Aroclor 1254	ND ND	ND ND	ND ND
Aroclor 1260	ND ND	ND	ND
Aroclor 1016	ND	ND	ND
	in Drinking Wat		
2,4-D	ND	ND	ND
2,4,5-TP (Silvex)	ND ND	ND ND	ND
Pentachlorophenol Dalapon	ND ND	ND ND	ND ND
Dinoseb	ND	ND	ND
Picloram	ND	ND	ND
Dicamba	ND	ND	ND
2,4 DB	ND	ND	ND
2,4,5 T Bentazon	ND ND	ND ND	ND ND
Dichlorprop	ND ND	ND	ND ND
	ND	ND	ND
Actiflorfin	ND	IND	140
Actiflorfin Dacthal (DCPA) 3,5-Dichlorobenzoic Acid	ND ND	ND ND	ND ND

3,5-Dichlorobenzoic Acid ND ND NI

J=Indicates an estimated concentration. This occurs when an analyte
concentration is below the calibration curve but is above the method detection limit.

Date	2/12/2009	2/12/2009	2/12/2009
Well ID Chemical	L-1	L-2	L-3
	in Drinking wa	ter	
Carbofuran	NS	ND	ND
Oxymal	NS	ND	ND
3-Hydroxycabofuran	NS.	ND	ND
Aldicarb Aldicarb sulfone	NS NS	ND ND	ND ND
Aldicarb sulfoxide	NS NS	ND	ND
Carbaryl	NS	ND	ND
Methomyl	NS	ND	ND
Propoxur (Baygon)	NS NS	ND	ND
Methiocarb Synthetic Or	NS rganic Compou	ND nde	ND
Endrin	ND	ND	ND
Lindane (BHC-Gamma)	ND	ND	ND
Methoxychlor	ND	ND	ND
Alachlor	ND ND	ND	ND
Atrazine Benzo(a)pyrene	0.04 ug/L J ND	ND ND	ND ND
Chlordane Technical	ND	ND	ND
Di(ethylhexyl)-Adipate	ND	ND	ND
Di(ethylhexyl)-phthalate	ND	ND	ND
Heptachlor	ND	ND	ND
Heyachlorobenzene	ND ND	ND	ND ND
Hexachlorobenzene Hexachlorocyclo-Pentadiene	ND ND	ND ND	ND ND
Simazine	ND	ND	ND
Aldrin	ND	ND	ND
Butachlor	ND	ND	ND
Dieldrin	ND ND	ND	ND
Metolachlor Metribuzin	ND ND	ND ND	ND ND
Propachlor	ND ND	ND	ND ND
Bromacil	0.21 ug/L JJ	ND	ND
Prometon	ND	ND	ND
Terbacil	ND	ND	ND
Diazinon	ND	ND	ND
EPTC 4,4-DDD	ND ND	ND ND	ND ND
4,4-DDE	ND ND	ND	ND
4,4-DDT	ND	ND	ND
Cyanazine	ND	ND	ND
Malathion	ND	ND	ND
Trifluralin Napthalene	ND ND	ND ND	ND ND
Fluorene	ND ND	ND	ND
Acenaphthylene	ND	ND	ND
Acenaphthene	ND	ND	ND
Anthracene	ND	ND	ND
Benz(A)anthracene Benzo(B)fluoranthene	ND ND	ND ND	ND ND
Benzo(G,H,I)peryene	ND	ND	ND
Benzo(K)fluoranthene	ND	ND	ND
Chrysene	ND	ND	ND
Dibenzo(A,H)anthracene	ND	ND	ND
Fluoranthene	ND	ND	ND
Indeno(1,2,3-CD)pyrene Phenanthrene	ND ND	ND ND	ND ND
Pyrene	ND	ND	ND
Benzyl Butyl Phthalate	ND	ND	ND
Di-N-Butyl Phthalate	ND	ND	ND
Diethyl Phthalate	ND	ND	ND
Dimethyl Phthalate	ND ND	ND ND	ND ND
Toxaphene Aroclor 1221	ND ND	ND ND	ND ND
Aroclor 1221	ND	ND	ND
Aroclor 1242	ND	ND	ND
Aroclor 1248	ND	ND	ND
Aroclor 1254	ND	ND	ND
Aroclor 1260 Aroclor 1016	ND ND	ND ND	ND ND
	in Drinking Wa		140
2,4-D	ND ND	ND	ND
2,4,5-TP (Silvex)	ND	ND	ND
Pentachlorophenol	ND	ND	ND
Dalapon	ND	ND	ND
Dinoseb Picloram	ND ND	ND ND	ND ND
Dicamba	ND ND	ND ND	ND ND
2,4 DB	ND	ND	ND
2,4,5 T	ND	ND	ND
Bentazon	ND	ND	ND
Dichlorprop	ND	ND	ND
Actiflorfin Dacthal (DCPA)	ND ND	ND ND	ND ND
3,5-Dichlorobenzoic Acid	ND ND	ND ND	ND ND
.,			

Date Well ID	3/12/2009 L-1	3/12/2009 L-2	3/12/2009 L-3	3/12/2009 Source
Chemical				Cource
Carbama	tes in Drink	ing water		
Carbofuran	NS	ND	ND	ND
Oxymal 3-Hydroxycabofuran	NS NS	ND ND	ND ND	ND ND
Aldicarb	NS	ND	ND	ND
Aldicarb sulfone	NS	ND	ND	ND
Aldicarb sulfoxide	NS	ND	ND	NS
Carbaryl	NS NS	ND ND	ND ND	NS NS
Methomyl Propoxur (Baygon)	NS	ND	ND	NS NS
Methiocarb	NS	ND	ND	NS
Synthetic	Organic C	ompounds		
Endrin	NS	NS	NS	NS
Lindane (BHC-Gamma)	NS NS	NS NS	NS NS	NS NS
Methoxychlor Alachlor	NS	NS	NS	NS
Atrazine	NS	NS	NS	NS
Benzo(a)pyrene	NS	NS	NS	NS
Chlordane Technical	NS	NS	NS	NS
Di(ethylhexyl)-Adipate Di(ethylhexyl)-phthalate	NS NS	NS NS	NS NS	NS NS
Heptachlor	NS	NS	NS	NS
Heptachlor Epoxide A&B	NS	NS	NS	NS
Hexachlorobenzene	NS	NS	NS	NS
Hexachlorocyclo-Pentadiene Simazine	NS NS	NS NS	NS NS	NS NS
Aldrin	NS	NS	NS	NS
Butachlor	NS	NS	NS	NS
Dieldrin	NS	NS	NS	NS
Metolachlor Metribuzin	NS NS	NS NS	NS NS	NS NS
Propachlor	NS	NS	NS	NS
Bromacil	NS	NS	NS	NS
Prometon	NS	NS	NS	NS
Terbacil	NS NS	NS	NS	NS
Diazinon EPTC	NS	NS NS	NS NS	NS NS
4,4-DDD	NS	NS	NS	NS
4,4-DDE	NS	NS	NS	NS
4,4-DDT	NS	NS	NS	NS
Cyanazine Malathion	NS NS	NS NS	NS NS	NS NS
Trifluralin	NS	NS	NS	NS
Napthalene	NS	NS	NS	NS
Fluorene	NS	NS	NS	NS
Acenaphthylene	NS NS	NS NS	NS NS	NS NS
Acenaphthene Anthracene	NS	NS	NS	NS
Benz(A)anthracene	NS	NS	NS	NS
Benzo(B)fluoranthene	NS	NS	NS	NS
Benzo(G,H,I)peryene	NS	NS	NS	NS
Benzo(K)fluoranthene Chrysene	NS NS	NS NS	NS NS	NS NS
Dibenzo(A,H)anthracene	NS	NS	NS	NS
Fluoranthene	NS	NS	NS	NS
Indeno(1,2,3-CD)pyrene	NS	NS	NS	NS
Phenanthrene Pyrene	NS NS	NS NS	NS NS	NS NS
Benzyl Butyl Phthalate	NS	NS	NS	NS
Di-N-Butyl Phthalate	NS	NS	NS	NS
Diethyl Phthalate	NS	NS	NS	NS
Dimethyl Phthalate	NS NS	NS NS	NS NS	NS NS
Toxaphene Aroclor 1221	NS	NS NS	NS	NS NS
Aroclor 1232	NS	NS	NS	NS
Aroclor 1242	NS	NS	NS	NS
Aroclor 1248	NS	NS	NS	NS
Aroclor 1254 Aroclor 1260	NS NS	NS NS	NS NS	NS NS
Aroclor 1200	NS	NS	NS	NS
	es in Drink			
2,4-D	ND	ND	ND	ND
2,4,5-TP (Silvex)	ND	ND	ND	ND
Pentachlorophenol Dalapon	ND ND	ND ND	ND ND	ND ND
Dinoseb	ND	ND ND	ND	ND
Picloram	ND	ND	ND	ND
Dicamba	ND	ND	ND	ND
2,4 DB	ND	ND	ND	ND
2,4,5 T Rentazon	ND ND	ND ND	ND ND	ND ND
Bentazon Dichlorprop	ND ND	ND ND	ND ND	ND ND
Actiflorfin	ND	ND	ND	ND
/ totalion in i				

Jactinal (IJCFA) ND ND N

3,5-Dichlorobenzoic Acid ND ND N

J=Indicates an estimated concentration. This occurs when an analyte concentration is below the calibration curve but is above the method detection limit.

JJ=The amount detected is below the Method's Reporting Level but greater than the lab's Practical Quantitation Level

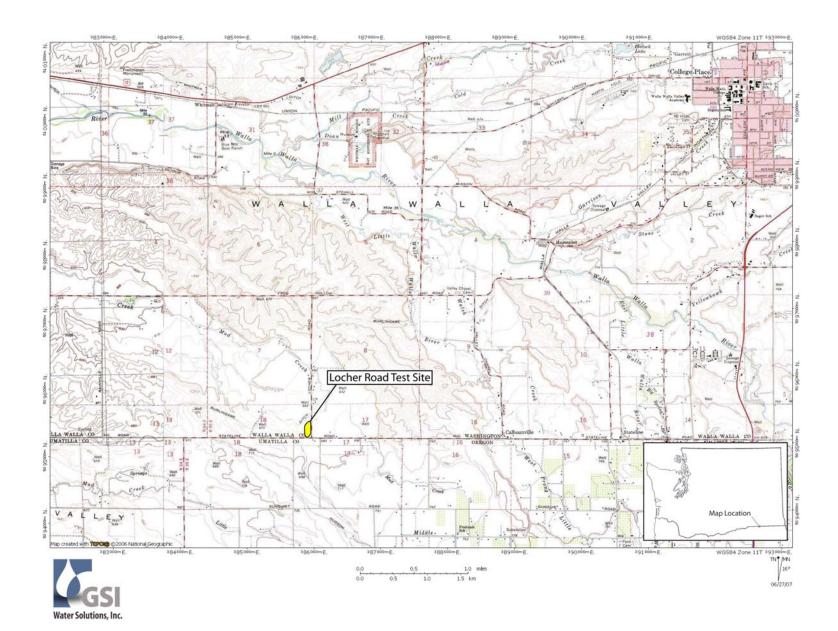


Figure 1. Area and regional setting.



Figure 2. Local setting, site location, and layout.



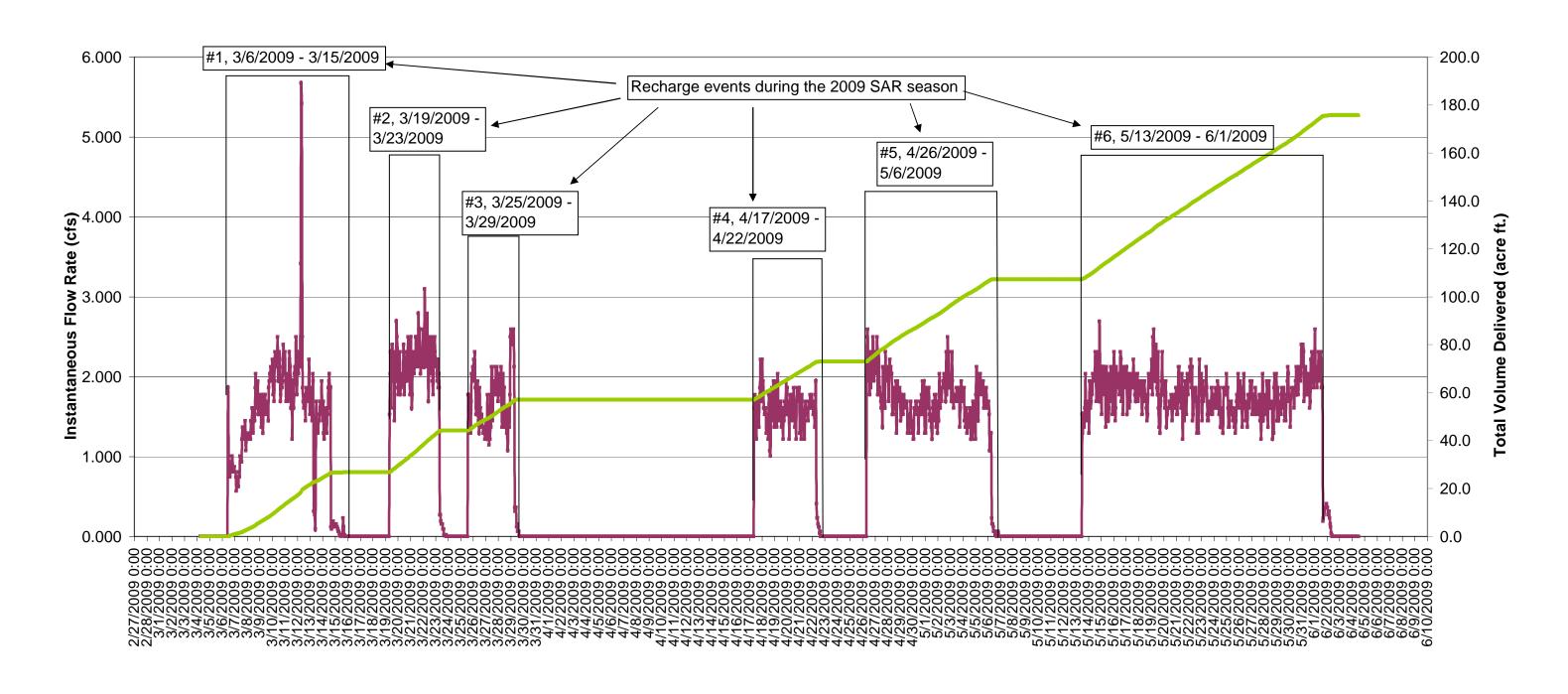
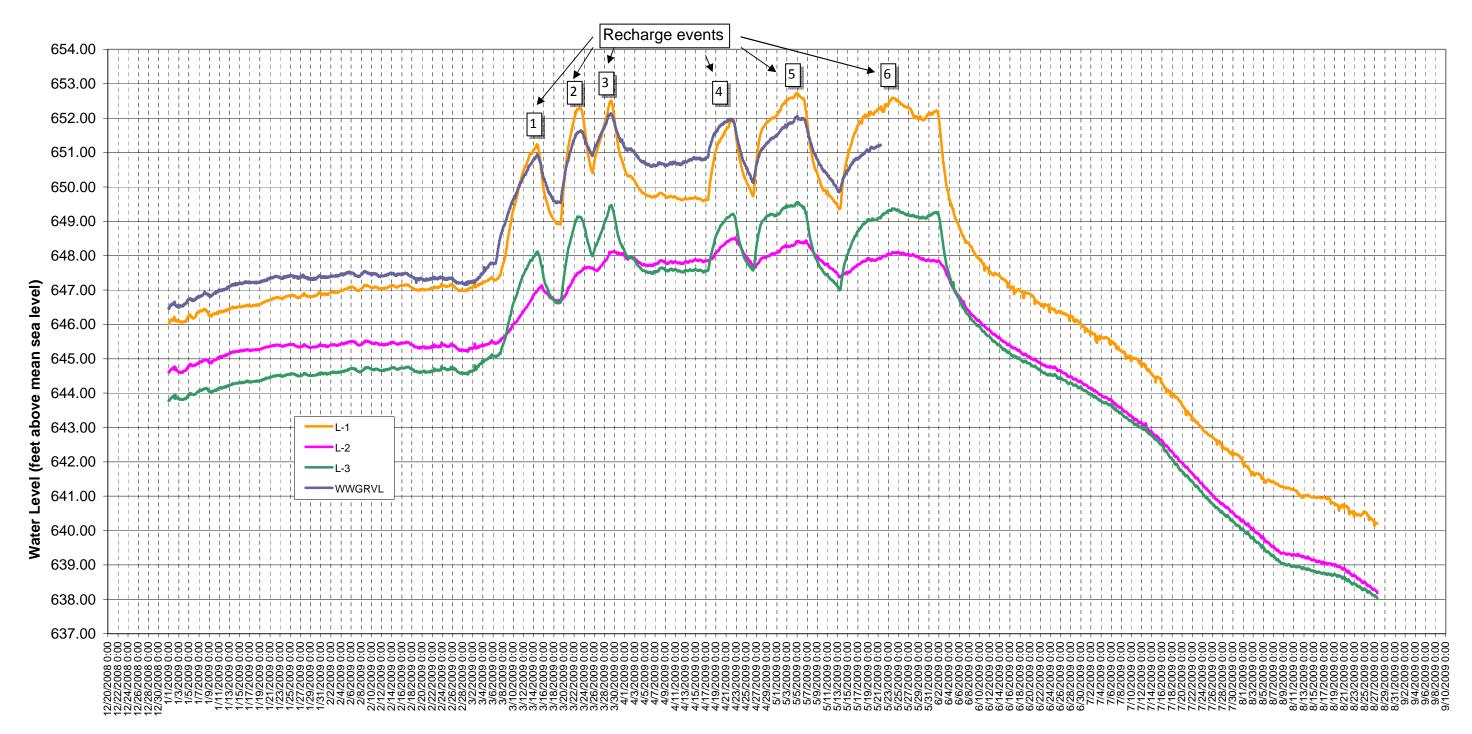


Figure 3. Instantaneous and total flow into the Locher Road SAR site during the 2009 season.

Date and Time

Instantaneous Flow Rate

Total Volume Delivered



Date and Time

Figure 4. Water levels in Locher Road SAR site wells before, during, and after the 2009 SAR season.

Locher Road SAR - Off Site Wells

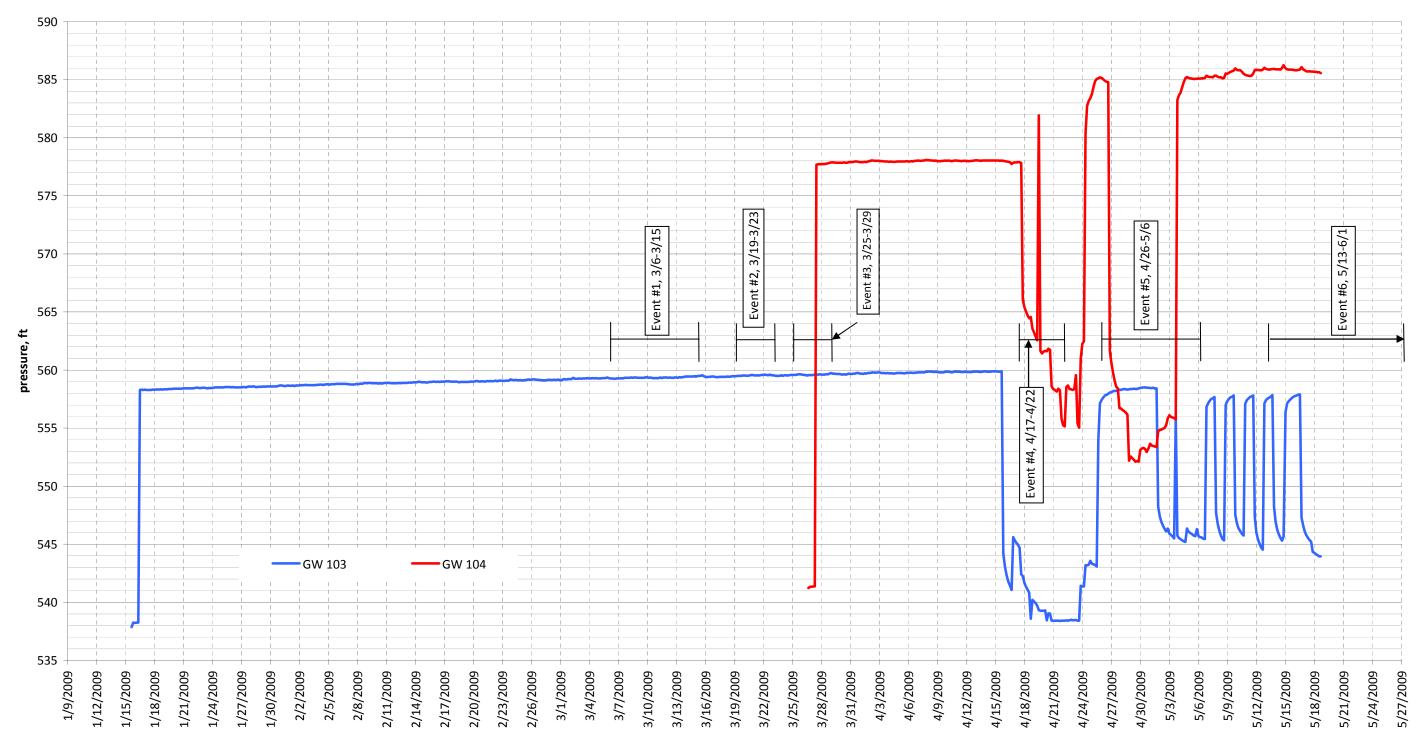
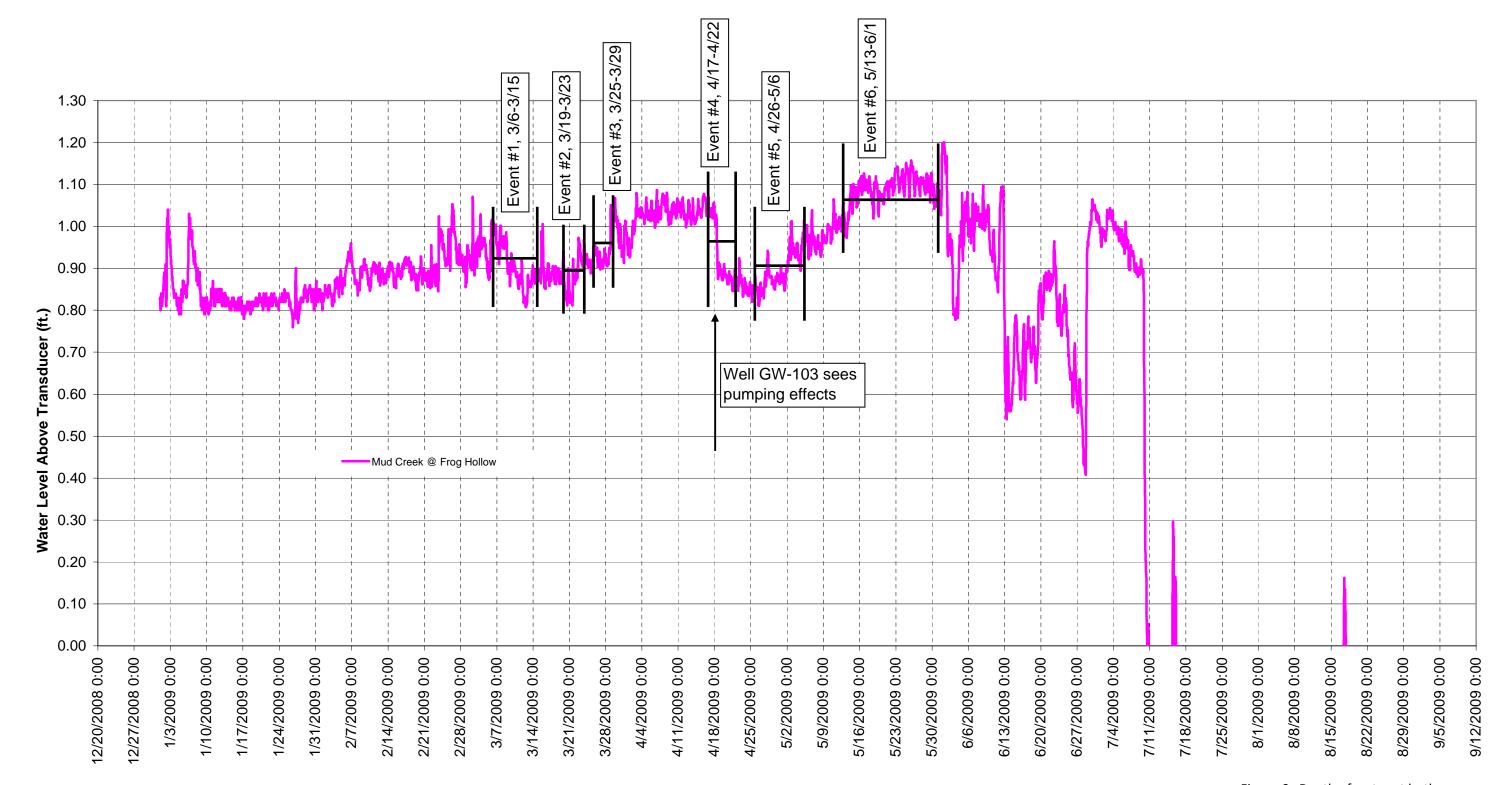


Figure 5. Water levels in Locher Road SAR offsite wells before and during the beginning of the 2009 season.



Date and Time

Figure 6. Depth of water at both surface water monitoring sites on Mud Creek.

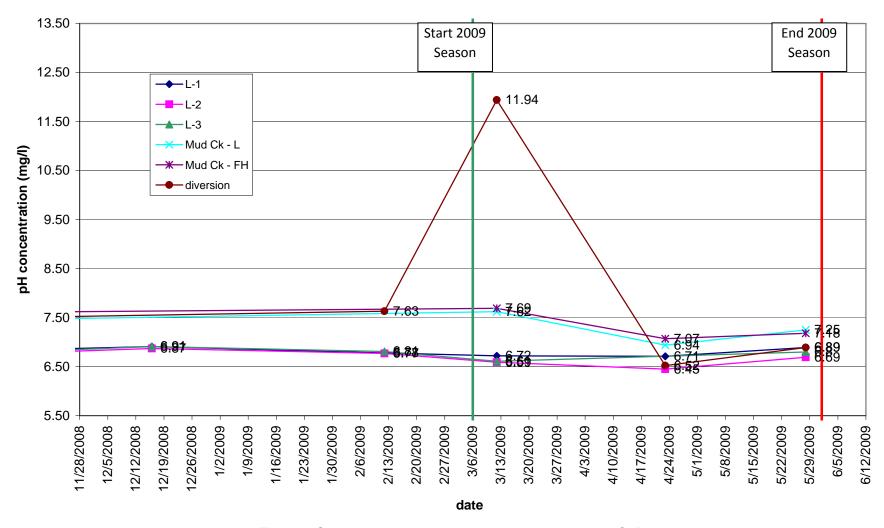


Figure 7. Groundwater and source water pH during the 2009 SAR season.

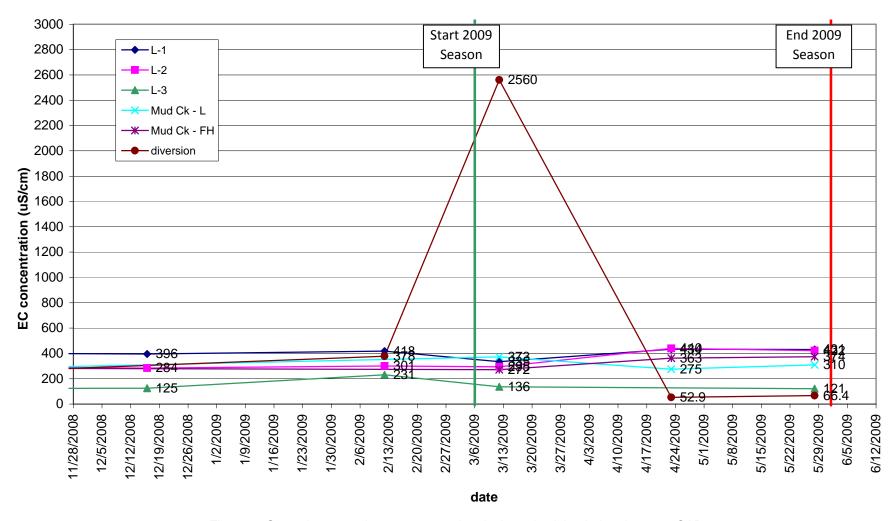


Figure 8. Groundwater and source water electrical conductivity during the 2009 SAR season.

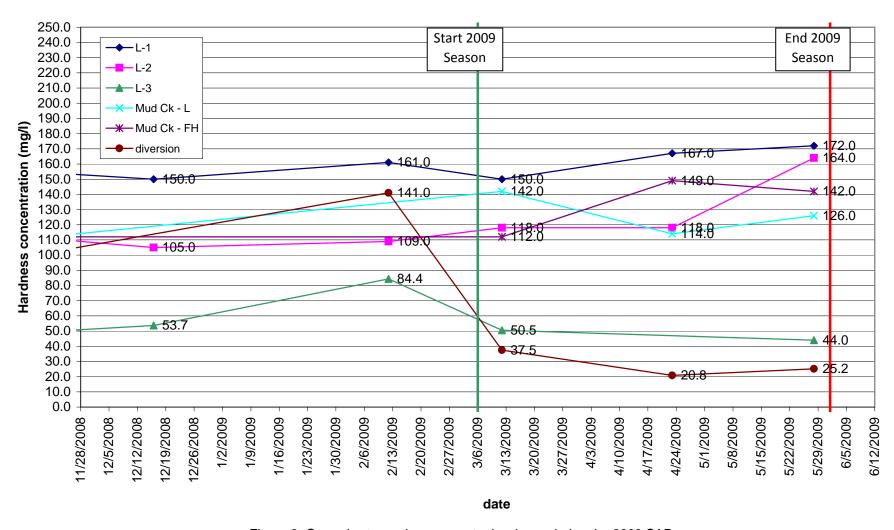


Figure 9. Groundwater and source water hardness during the 2009 SAR season.

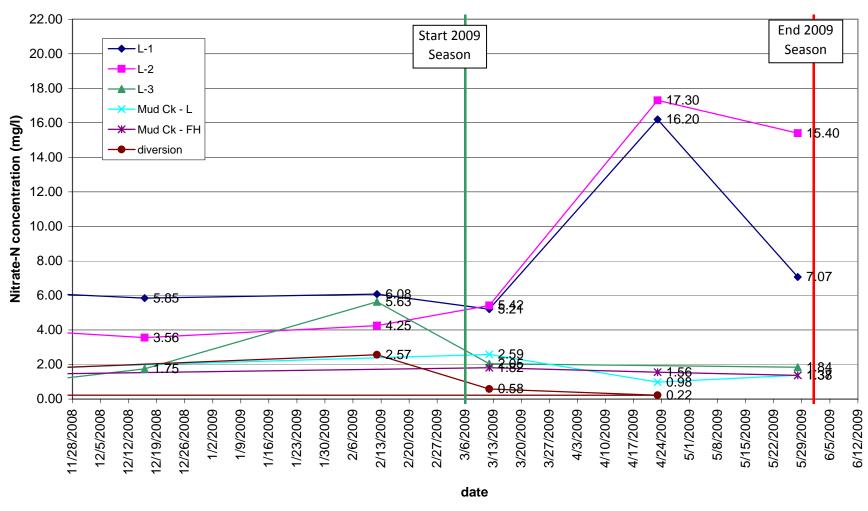


Figure 10. Groundwater and source water nitrate-N during thee 2009 SAR season.

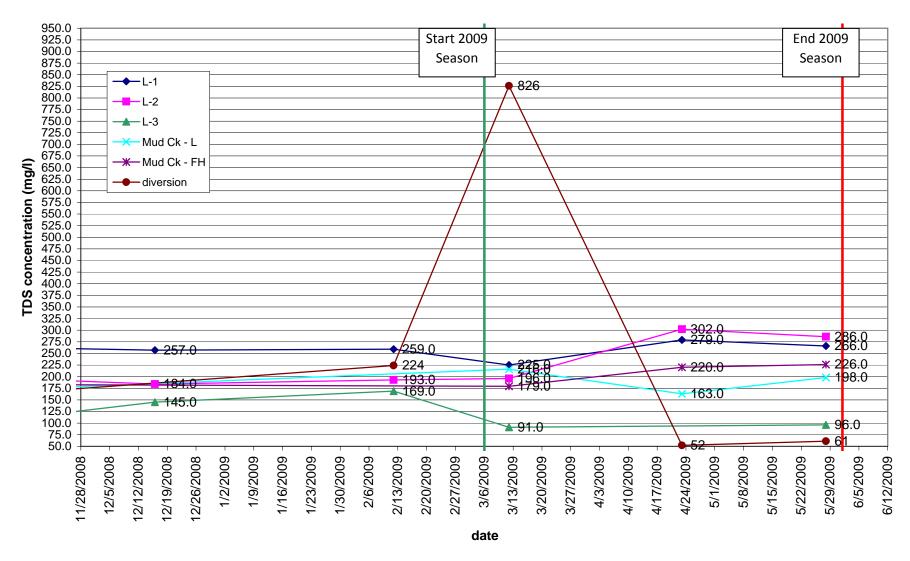


Figure 11. Groundwater and source water total dissolved solids during the 2009 SAR season.

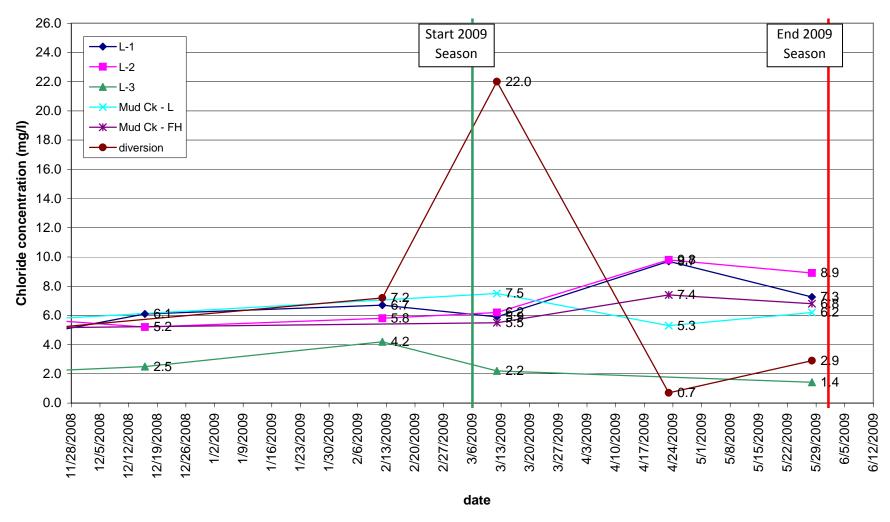


Figure 12. Groundwater and source water chloride during the 2009 SAR season.

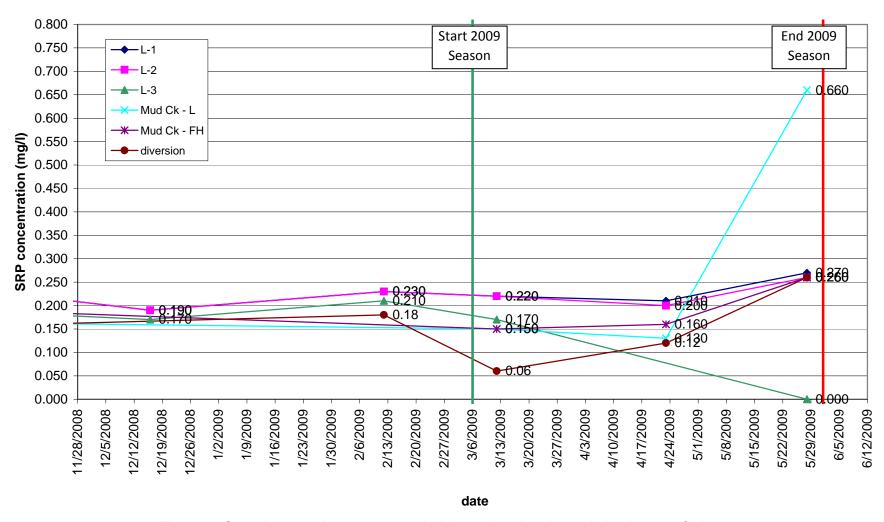


Figure 13. Groundwater and source water soluable reactive phosphorus during the 2009 SAR season.

Appendix A

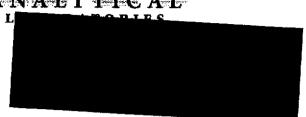


Corporate Office 800

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Bellingham WA

360.671.0688 • 360.671.1577(ax



December 10, 2008

Page 1 of 1

Mr. Troy Baker Walla Walla Basin Watershed Council 810 S Main Street Milton-Freewater, OR 97862

RE: 08-15517 - Locker/Hall-Wetland

Dear Mr. Troy Baker,

Your project: Locker/Hall-Wetland, was received on Wednesday October 29, 2008.

The following comments are reported for your project:

Samples 32783 (H-3) and 32784 (H-2) were also analyzed for PCBs and Toxaphene under method 508.1. Both sample were non-detect for these compounds. There was no charge for this analysis.

If you have questions phone me at 800 755-9295.

Respectfully Submitted,

Lawrence J Henderson, PhD Director of Laboratories

Enclosures Data Report

QC Reports

Chain of Custody



Page 1 of 2

Data Report

Client Name: Walla Walla Basin Watershed Council

810 S Main Street

Milton-Freewater, OR 97862

Report Date: 11/7/2008 Reference Number: 08-15517

Project: Locker/Hall-Wetland

Collected By: Baker

Date Received: 10/29/2008
Peer Review:

Lab Num	nber: 32780 Sam	ple Descriptio	on: L-1 - L	ocker obs#	1			Sample	Date:	10/28/2008	
CAS ID#	Analyte	Result	PQL	MDL	Units	DF	Method	Analyzed	Analys	t Batch	Commen
E-10139	HYDROGEN ION (pH)	6.81	_		pH Units	1.0	SM4500-H+ B	1D/29/2008	MAK	PH_D81029	
14797-55-8	NITRATE-N	6.38	0.100	0.015	mg/L	1.0	300.0	10/29/2008	BJ	1081029A	
16887-00-6	CHLORIDE	3.5	0.1	0.012	mg/L	1.0	300.0	10/29/2008	BJ	1061029A	
E-10173	TOTAL DISSOLVED SOLIDS	265	10	6	mg/L	1.0	SM2540 C	10/20/2008	CCN	TDS_081029	
14265-44-2	ORTHO-PHOSPHATE	0.24	0.01	0.002	mg/L	1.0	SM4500-P F	10/20/2008	60	OPHOS-081029	
E-10184	ELECTRICAL CONDUCTIVITY	403	10		uS/cm	1.0	SM2510 B	11/3/2008	CCN	EC_081103	
E-10617	TURBIDITY	0.29	0.05	0.02	NTU	1.0	180.1	10/29/2008	MAK	TURB_081029	
15541-45-4	BROMATE	ND	0.005	0.0013	mg/L	1.0	300.1	11/4/2008	MAb	D061164A	
E-11778	HARDNESS	158.1	3.30	0.055	mg CaCo	1.0	200.7	11/3/2008	BJ	200.7-081103A	
E-10117	CHEMICAL OXYGEN DEMAND	ND	8	2	mg/L	1.0	SM5220 D	11/4/2008	MAK	COD_081104	
Lab Nun	nber: 32781 Sam	ple Description	on: L-2 - L	ocker obst	‡ 2			Sample	Date:	10/28/2008	
CAS ID#	Analyte	Result	PQL	MDL	Units	DF	Method	Analyzed	Analys	t Batch	Commer
E-10139	HYDROGEN ION (pH)	6.73			pH Units	1.0	SM4500-H+ B	10/29/2006	MAK	PH_081029	
14797-55-8	NITRATE-N	4.25	0.100	0.015	mg/L	1.0	300.0	10/20/2008	81	1081029A	
16887-00-6	CHLORIDE	6.2	0.1	0.012	mg/L	1.0	300.0	10/20/2008	BJ	1081029A	
E-10173	TOTAL DISSOLVED SOLIDS	201	10	6	mg/L	1.0	SM2540 C	10/29/2008	CCN	TCS_081029	
14265-44-2	ORTHO-PHOSPHATE	0.24	0.01	0.002	mg/L	1,0	SM4500-P F	10/20/2008	SD	OPHOS-081029	
E-10184	ELECTRICAL CONDUCTIVITY	291	10		ยS/cm	1.0	SM2510 B	11/3/2008	CCN	EC_061103	
E-10617	TURBIDITY	1.79	0.05	0.02	NTU	1.0	180.1	10/29/2006	MAK	TURB_081029	
15541-45-4	BROMATE	ND	0.005	0.0013	mg/L	1.0	300,1	11/4/2008	MVP	D081104A	
E-11778	HARDNESS	116.0	3.30	0.055	mg CaC	1.0	200.7	11/3/2008	돲	200.7-081103A	
E-10117	CHEMICAL OXYGEN DEMAND	ND	8	2	mg/L	1.0	SM5220 D	11/4/2000	MAK	COD_081184	
Lab Nun	nber: 32782 Sam	nple Descripti	оп: L-3 - I	ocker obs	#3			Sampl	e Date:	10/28/2008	3
CAS ID#	Analyte	Result	PQL	MDL	Units	DF	Method	Analyzed	Analy	st Batch	Comme
E-10139	HYDROGEN ION (pH)	6.75			pH Units	5 1,0	SM4500-H+ B	10/29/2008	MAK	PH_081029	
14797-55-8	NITRATE-N	0.43	0.100	0.015	mg/L	1.0	300.0	10/29/2008	8.1	1081029A	
16887-00-6	CHLORIDE	1.9	0.1	0.012	mg/L	1.G	300.0	10/29/2008	BJ	(081029A	
E-10173	TOTAL DISSOLVED SOLIDS	94	10	6	mg/L	1.0	SM2540 C	10/29/2008	CCN	TDS_981029	
14265-44-2	ORTHO-PHOSPHATE	0.19	0.01	0.002	mg/L	1.0	SM4500-P F	16/29/2006	50	OPHOS-081029	
		400	10		uS/cm	1.D	SM2510 B	11/3/2008	CCN	EC_081103	
E-10184	ELECTRICAL CONDUCTIVITY	122	10		44,011	-,-	Onder D			-	

PQL = Practical Quantitation Limit is the lowest level that can be acheived within specified limits of precision and accuracy during routine laboratory operating conditions.

ND = Not detected above the listed practical quantitation limit (PQL) or not above the Mathod Detection Limit (MDL), if requested.

D.F. - Dilution Factor

WSDOE Lab C1251 WSDOH Lab 046



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CARBAMATES IN DRINKING WATER

Client Name: Walla Walla Basin Watershed Council

810 S Main Street

Milton-Freewater, OR 97862

Reference Number: 08-15517

Project: Locker/Hall-Wetland

Field ID: L-1

Sample Description: Locker obs#1

Sampled By: Baker Sample Date: 10/28/2008

Source Type: Sampler Phone:

Lab Number: 04632780 Report Date: 11/4/2008 Date Analyzed: 10/31/2008 Extraction Date: 531_081031

Analyst: CO Supervisor: Analytical Method: 531.2

Carbamates

CAS	COMPOUND	RESULTS	Units	PQL	MDL	MCL	COMMENT	
	EPA Regulated							
23135-22-0	OXYMAL	ND	ug/L	1.0	0.3	200		
1563-66-2	CARBOFURAN	ND	ug/L	1.0	0.2	40		
	EPA Unregulated							
1646-87-3	ALDICARB SULFOXIDE	ND	rđ/F	1.0	0.3			
1646-88-4	ALDICARB SULFONE	[®] ND	u g/L	1.0	0.3			
16752-77-5	METHOMYL	ND	ug/L	1.0	0.3			
16655-82-6	3-HYDROXYCARBOFURAN	ND	ug/L	1.0	0.3			
116-06-3	ALDICARE	ND	ug/L	1.0	0.3			
63-25-2	CARBARYL	ND	ug/L	1.0	0.2			
	State Unregulated - Other				-			
114-26-1	PROPOXUR (BAYGON)	ND	ug/L	1.0	0.4			
2032-65-7	METHIOCARB	ND	ug/L	1.0	0.3 .			

MCL-Maximum Contaminant Level, maximum permissible level of a contaminant in water established by EPA, NPDWR. State Advisory Level (SAL) for Unrequirted compounds.

A blank MCL or SAL value indicates a level is not currently established.

PQL - Practical Quantitation Limit is the concentration of the standard analyzed during the initial calibration.

MDL - Method Detection Limit is the lab's minimum concentration a compound can be measured and reported with 99% confidence that the compound concentration is greater than zero.



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Page 1 of 1

CARBAMATES IN DRINKING WATER

Client Name: Walla Walla Basin Watershed Council

810 S Main Street

Milton-Freewater, OR 97862

Reference Number: 08-15517

Project: Locker/Hall-Wetland

Field ID: L-2

Sample Description: Locker obs#2

Sampled By: Baker

Sample Date: 10/28/2008

Source Type:

Sampler Phone:

Lab Number: 04632781 Report Date: 11/4/2008 Date Analyzed: 10/31/2008

Extraction Date: 531_081031

Analyst: /C]

Supervisor:

Analytical Method: 531.2

Carbamates

CAS	COMPOUND	RESULTS	Units	PQL	MDL	MCL	COMMENT
	EPA Regulated						
23135-22-0	OXYMAL	ND	υg/L	1.0	0.3	200	
1563-66-2	CARBOFURAN	ND	ug/L	1.0	0.2	40	
	EPA Unregulated						
1646-87-3	ALDICARB SULFOXIDE	ND	ug/L	1.0	0.3		
1646-88-4	ALDICARB SULFONE	ND	ug/L	1.0	0.3		
16752-77-5	METHOMYL	ND	ug/L	1.0	0.3		
16655-82-6	3-HYDROXYCARBOFURAN	ND	ug/L	1.0	0.3		
116-06-3	ALDICARB	ND	ug/L	1.0	0.3		
63-25-2	CARBARYL	ND	ug/L	1.0	0.2		
	State Unregulated - Other						
114-26-1	PROPOXUR (BAYGON)	ND	ug/L	1.0	0.4		
2032-65-7	METHICCARB	ND	ug/L	1.0	0.3		

An amount of "ND" indicates that the compound was not detected above the Lab's Method Defection Limit - MOL MCL Minimum Contaminant Level, maximum permissible level of a contaminant in water established by EPA, NPDWR. State Advisory Level (SAL) for Unregulated compa
A blank MCL or SAL value indicates a level is not currently established.

POL - Practical Quantitation Limit is the concentration of the standard analyzed during the initial cathication.

Mall - Method Detection Limit is the lat's minimum concentration a compound can be measured and reported with 99% confidence that the compound concentration is greater than zero.

These test results meet all the requirements of NELAC, unless otherwise stated in writing, and relate only to these samples. FORM: SOC_GEN



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CARBAMATES IN DRINKING WATER

Client Name: Walla Walla Basin Watershed Council

810 S Main Street

Milton-Freewater, OR 97862

Reference Number: 08-15517

Project: Locker/Hall-Wetland

Field ID: L-3

Sample Description: Locker obs#3

Sampled By: Baker Sample Date: 10/28/2008

Source Type: Sampler Phone: Date Analyzed: 10/31/2008

Extraction Date: 531_081031

Analyst: Supervisor

Analytical Method: 531.2

Lab Number: 04632782

Report Date: 11/4/2008

Carbamates

CAS	COMPOUND	RESULTS	Units	PQL	MDL	MCL	COMMENT	
	EPA Regulated							
23135-22-0	OXYMAL	ND	ug/L	1.0	0.3	200		
1563-66-2	CARBOFURAN	ND	ug/L	1.0	0.2	40		
	EPA Unregulated							
1646-87-3	ALDICARB SULFOXIDE	, ND	ug/L	1.0	0.3			
1646-88-4	ALDICARB SULFONE	ND	ug/L	1.0	0.3			
16752-77-5	METHOMYL	ND	ug/L	1.0	0.3			
16655-82-6	3-HYDROXYCARBOFURAN	ND	ug/L	1.0	0.3			
116-06-3	ALDICARB	ND	ug/L	1.0	0.3			
63-25-2	CARBARYL	ND	ug/L	1.0	0.2			
	State Unregulated - Other	×						
114-26-1	PROPOXUR (BAYGON)	ND	vg/l	1.0	0.4			
2032-65-7	METHIOCARB	ND	ug/L	1.0	0.3			

An emount of "NO" indicates that the compound was not detected above the Lab's Method Detection Limit - MOL.

MCL. Maximum Contaminant Level, maximum permissible level of a contaminant in water established by EPA, NPDWR. State Advisory Level (SAL) for Unregulated compounds.

A blank MCL or SAL value indicates a level is not currently established.

PQL - Practical Quantitation Limit is the concentration of the standard analyzed during the initial calibration.

MDL, - Method Detection Limit is the lab's minimum concentration a compound can be measured and reported with 99% confidence that the compound concentration is greater than zero.

J - Estimated value.

These test results meet all the requirements of NELAC, unless otherwise stated in writing, and relate only to these samples. FORM: SOC_GEN



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SAMPLE INDEPENDENT QUALITY CONTROL REPORT

Laboratory Fortified Blank

Reference Number: 08-15517

		15	True			%		QC	
Batch	Analyte	Result	Value	Units	Method	Recovery	Limits	Qualifier Type*	Comment
200.7-081103A	HARDNESS	73.3	69.5	mg/L	200.7	105	80-120	LFB	
515 <u>0</u> 81031	2,4 - D		_						
313_081031	-	1.87	2	ug/L	515.1	94	70-130	LFB	
	2,4 - DCAA (SURR)	101	_	%	515.1		70-130		
	2,4 DB	- 7.69	8	ug/L	51 5.1	96	70-130		
	2,4,5 - TP (SILVEX)	0.92	1	ug/L	51 5. 1	92	70-130		
	2,4,5 T	0.91	1	ug/L	515.1	91	70-130		
	ACIFLUORFEN	1.04	1	ug/L	515.1	104	70-130		
	BENTAZON	1.79	2	ug/t.	515.1	80	70-130		
	CHLORAMBEN	0.73	1	ug/L	515.1	73	70-130		
	DALAPON	10.1	13	ug/L	515.1	78	70-130	•	
	DICAMBA	0.89	1	ug/L	515.1	89	70-130		
	DICHLORPROP	2.63	3	ug/L	515.1	8B	70-130		
	DINOSEB	1.74	2	ug/L	515.1	87	70-130		
	PENTACHLOROPHENOL	88.0	1	ug/L	515.1	88	70-130		
	PICLORAM	0.91	1	ug/L	515.1	91	70-130		
	TOTAL (DCPA & Metabolites)	0.87	1	ug/L	515.1	87	70-130		
525_081105	1,3-DIMETHYL-2-NITROBENZENE (Suit)	79		A.	tor 0		70 400		
	4,4-DDD	1.2		%	525.2	400	70-130	LFB	
	4,4-DDE		1	ug/L	525.2	120	70-130		
	4,4-DDT	1.12	1	ug/L	525.2	112	70-130		
	BISPHENOL-A	1.09	1	ug/L	525.2	109	70-130		
	DIAZINON	4.1	5	t/g/L	525.2	82	85-115		
		0.9	1	ug/L	525.2	90	7 0- 130		
	LINDANE (BHC - GAMMA)	1.01	1	ug/L	525.2	101	70-130		
	PERYLENE-D12 (Suπ)	94		%	525.2		70-130		
	PYRENE-D10 (Sum)	89		%	525.2		70-130		
	SIMAZINE	0.97	1	ug/L	525.2	97	70-130		
	TRIPHENYLPHOSPHATE (Surr)	120		%	525,2		70-130		
525X_081105	1-NAPHTHALENEACETAMIDE	2.07	2	ug/L	525.2	104	70-130	LFB	
	CHLORPYRIFOS	1.05	1	ug/L	525.2	105	70-130		
	DICOFOL	2.57	2	ug/L	525.2	129	70-130		

[%] Recovery = (Result of Analysis)/(True Value) * 100

NA = Indicates % Recovery could not be calculated.

QCS: Quality Control Sample, a solution containing known concentrations of method analytes which is used to fortify an aliquot of reagent metrix. The QCS is obtained from an external source and is used to check lab performance.

LFB: Laboratory Fortified Blank, an aliquot of reagent matrix to which known quantities of method analytes are added in the leb. The LFB is analyzed exactly like a sample, and its purpose is to determine whether method performance is within accepted control limits.

MB or LRB: Method Blank or Laboratory Reagent Blank, an alliquol of reagent matrix is analyzed exactly like a sample, and its purpose is to determine if there is background contamination.



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SAMPLE INDEPENDENT QUALITY CONTROL REPORT

Laboratory Fortified Blank

Reference Number: 08-15517 Report Date: 12/10/08

		96	True		•	%		QC	
Batch	Analyte	Result	Value	Units	Method	Recovery	v Limits	Qualifier Type"	Comment
525X_081105	FENARIMOL	1.29	1	ng/L	525,2	129	70-130	LFB	- CONTRACT IL
	HEXAZINONE	1.26	1	ug/L	525.2	126	70-130		
	MALATHION	1.04	1	ug/L	525.2	104	70-130		
	METALAXYL	2.26	2	ug/L	525.2	113	70-130		
	METHIDATHINON *	2.92	2	ug/L	52 5.2	146	85-115	HR	
	MEVINPHOS	1.12	1	ug/L	525.2	112	70-130	****	
	MYCLOBUTANIL	2.43	2	ug/L	525.2	122	85-115		
	NAPROPAMIDE	1.16	1	ug/L	525.2	116	70-130		
	PARATHION-ETHYL	1.05	1	ug/L	525.2	105	70-130		
	PHOSMET	2.68	2	ug/L	525.2	134	70-130	HR	
	PROPARGITE	2.23	2	ug/L	525.2	112	85-115	****	
	TRIADIMEFON	1.05	1	ug/L	525.2	105	70-130		
	TRIFLUMIZOLE	1.35	2	ug/L	525.2	68	85-115	M1	
			_		020.2	00	00-115	N I	
531 _08 1031	3-HYDROXYCARBOFURAN	8	10	ug/L	531.2	80	70-130	LFB	
	ALDICARB	7.9	10	ug/L	531.2	79	70-130		
	ALDICARB SULFONE	8.1	10	ug/L	531.2	81	70-130		
	ALDICARB SULFOXIDE	~ 8.4	10	ug/L	531.2	B4	70-130		
	BDMC (SURR)	85		%	531.2	- -	70-130		
	CARBARYL	10	10	ug/L	531.2	100	70-130		
	CARBOFURAN	8	10	ug/L	531.2	80	70-130		
	METHIOCARB	7.9	10	ug/L	531,2	79	70-130		
	METHOMYL	9.8	10	ug/L	531.2	98	70-130		
	OXYMAL	. 8	10	ug/L	531.2	80	70-130		
	PROPOXUR (BAYGON)	7.3	10	ug/L	531.2	73	70-130		
				-5-	001.2		10-100		
COD_081104	CHEMICAL OXYGEN DEMAND	49	50	mg/L	SM5220 D	98	80-120	LFB	
OPHOS-081029	ORTHO-PHOSPHATE	1.02	1.00	mg/L	SM4500-P F	102	70-130	LFB	
tds_081029	TOTAL DISSOLVED SOLIDS	494	500	mg/L	SM2540 C	99	80-120	LFB	
tds_081029	TOTAL DISSOLVED SOLIDS	496	500	mall	C) 12540 C	00	60 40C	. —	
		400	aru.	mg/L	SM2540 C	99	80-120	LFB	

^{*}Notation:

[%] Recovery = (Result of Analysis)/(True Value) * 100

NA = Indicates % Recovery could not be calculated.

QCS: Quality Control Sample, a solution containing known concentrations of method analytes which is used to fortify an alkquot of reagent matrix. The QCS is obtained from an external source and is used to check lab performance.

LFB: Laboratory Fortified Blank, an aliquot of reagant matrix to which known quantities of method analytes are added in the lab. The LFB is analyzed exactly like a sample, and its purpose is to determine whether method performance is within accepted control limits.

MB or LRB; Method Blank or Laboratory Reagent Blank, an aliquot of reagent matrix is analyzed exactly like a sample, and its purpose is to determine if there is background contamination.



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SAMPLE INDEPENDENT QUALITY CONTROL REPORT

Laboratory Fortified Blank

Reference Number: 08-15517

			True			%		QC	
Batch	Analyte	Result	Value	Units	Method	Recove	ry Limits	Qualifier Type*	Comment
tds_081029	TOTAL DISSOLVED SOLIDS	498	500	mg/L	SM2540 C	100	80-120	LFB	
lds_081029	TOTAL DISSOLVED SOLIDS	500	500	mg/L	SM2540 C	100	80-120	LFB	

^{*}Notation:

[%] Recovery = (Result of Analysis)/(True Value) * 100

NA = Indicates % Recovery could not be calculated.

QCS: Quality Control Sample, a solution containing known concentrations of method analytes which is used to fortify an aliquot of reagent matrix. The QCS is obtained from an external source and is used to check tab performance.

LFB: Laboratory Fortified Blank, an aliquot of reagent matrix to which known quantities of method analytes are added in the lab. The LFB is analyzed exactly like a sample, and its purpose is to determine whether method performance is within accepted control limits.

MB or LRB: Method Blank or Laboratory Reagent Blank, an aliquot of reagent matrix is analyzed exactly like a sample, and its purpose is to determine if there is background contamination.



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SAMPLE INDEPENDENT QUALITY CONTROL REPORT

Low Level Laboratory Fortified Blank

Reference Number: 08-15517

			True			%		QC	
Betch	Analyte	Result	Value	Units	Method	Recovery	Limits	Qualifier Type*	Comment
531_081031	3-HYDROXYCARBOFURAN	0.6	1	ug/L	531.2	60	50-150	LFBD	
	ALDICARB	0.8	1	ug/L	531.2	80	50-150		
	ALDICARE SULFONE	0.6	1	ug/L.	531.2	60	50-150		
	ALDICARB SULFOXIDE	8.0	1	ug/L	531.2	80	50-150		
	BDMC (SURR)	79		%	531.2		50-150		
	CARBARYL	. 0.8	1	ug/L	531.2	80	50-150		
	CARBOFURAN	0.6	1	ug/L	531.2	60	50-150		
	METHIOCARB	0.6	1	ug/L	531,2	60	50-150		
	METHOMYL	0.7	1	ug/L	531.2	70	50-150		
OX	OXYMAL	1	1	ug/L	531.2		50-150		
	PROPOXUR (BAYGON)	0.6	1	ug/L	531.2	60	50-150		

^{*}Notation:

[%] Recovery = (Result of Analysis)/(True Value) * 100

NA = Indicates % Recovery could not be calculated.

QCS: Quality Control Sample, a solution containing known concentrations of method analytes which is used to fortify an aliquot of reagent matrix. The QCS is obtained from an external source and is used to check lab performance.

LFB: Laboratory Fortified Blank, an aliquot of reagent matrix to which known quantities of method analytes are added in the lab. The LFB is analyzed exactly like a sample, and its purpose is to determine whether method performance is within accepted control limits.

MB or LRB: Method Blank or Laboratory Reagent Blank, an aliquot of reagent matrix is analyzed exactly like a sample, and its purpose is to determine if there is background contamination.



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SAMPLE INDEPENDENT QUALITY CONTROL REPORT

Laboratory Reagent Blank

Reference Number: 08-15517

			True			%	QC	
Batch	Analyte	Result	Value	Units	Method	Recovery Limits	Qualifier Type*	Comment
200.7-081103A	HARDNESS	ND		mg/L	200.7	10.000	X LRB	
COD_081104	CHEMICAL OXYGEN DEMAND	ND		mg/L	SM5220 D	4.0000) LRB	
D081104A	BROMATE	` ND		mg/L	300.1	0.00500) LRB	
1081029A	CHLORIDE NITRATE-N	ND ND		mg/L mg/L	300.0 300.0	0.10000 0.10000		
OPHOS-081029	ORTHO-PHOSPHATE	ND		mg/L	SM4500-P F	0.10000) LRB	

[%] Recovery = (Result of Analysis)/(True Value) * 100

NA = Indicates % Recovery could not be calculated.

QCS: Quality Control Sample, a solution containing known concentrations of method analytes which is used to fortify an aliquot of reagent matrix. The QCS is obtained from an external source and is used to check lab performance.

LFB: Laboratory Fortified Blank, an aliquot of reagent matrix to which known quantities of method analytes are added in the lab. The LFB is analyzed exactly like a sample, and its purpose is to determine whether method performance is within accepted control limits.

MB or LRB. Method Blank or Laboratory Reagent Blank, an assount of reagent matrix is analyzed exactly like a sample, and its purpose is to determine if there is background contamination.



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SAMPLE INDEPENDENT QUALITY CONTROL REPORT

Method Blank

Reference Number: 08-15517

			True			%		QC	
Batch	Analyte	Result	Value	Units	Method	Recovery Li	imits	Qualifier Type*	Comment
200.7-081103A	HARDNESS	ND		mg/L	200.7		82000	MB	
515_081031	2,4 - D	ND		ug/L	515.1	0.0	05000	MB	
	2,4 - DCAA (SURR)	96		%	515.1	0.	00000	PAL	
	2,4 DB	ND.		ug/L	515.1	0.5	25000		
	2,4,5 - TP (SILVEX)	ND		ug/L	515.1		10000		
	2,4,5 T	ND		ug/L	515.1		100D0		
	ACIFLUORFEN	ND		ug/L	515.1		50000		
	BENTAZON	ND		ug/L	515.1		12000		
	CHLORAMBEN	ND	,	ug/L	515.1		20000		
	DALAPON	ND		ug/L	515.1		50000		
	DCPA (ACID METABOLITES)	ND		ug/L	515.1		10000		
	DICAMBA	ND		ug/L	515.1		05000		
	DICHLORPROP	ND		ug/L	515.1		12000		
	DINOSEB	ND		ug/L	515.1		10000		
	PENTACHLOROPHENOL	ND		ug/L	515.1		02000		
	PICLORAM	ND		ug/L	515.1		35000		
	TOTAL (DCPA & Metabolites)	- ND		ug/L	515.1		02000		
525_081105	1.3-DIMETHYL-2-NITROBENZENE (Sum)	80		%	525.2			MB	
	4,4-DDD	ND		ug/L	525.2	0.0	05000		
	4,4-DDE	ND		ug/L	525.2	0.0	05000		
	4,4-DDT	ND		ug/L	525.2	0.0	05000		
	BISPHENOL-A	" ND		ug/L	525.2	1.0	00000		
	DIAZINON	ND		ug/L	525.2		05000		
	LINDANE (BHC - GAMMA)	ND		ug/L	525.2	0.0	2000		
	PERYLENE-D12 (Surr)	85		%	525.2				
	PYRENE-D10 (Surr)	91		%	525.2				
	SIMAZINE	ND		ug/L	525.2	0.0	2000		
4	TRIPHENYLPHOSPHATE (Surr)	, 117		%	525.2				
525X_081105	1-NAPHTHALENEACETAMIDE	NĐ		ug/L	525.2	0.1	10000	MB	
	AZINPHOS-METHYL	ND		ug/L	525.2		00000	•	
				•					

^{*}Notation:

[%] Recovery = (Result of Analysis)/(True Value) * 100

NA = Indicates % Recovery could not be calculated.

QCS: Quality Control Sample, a solution containing known concentrations of method analytes which is used to fortify an aliquot of reagent matrix. The QCS is obtained from an external source and is used to check lab performance.

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SAMPLE INDEPENDENT QUALITY CONTROL REPORT

Method Blank

Reference Number: 08-15517

		*	True			%		QC	
Baich	Алаlyte	Result	Value	Units	Method	Recovery L	imils	Qualifier Type*	Comment
525X_0B1105	CHLORPYRIFOS	ND		ug/L	525.2	O	0.00000	МВ	
	DICOFOL	ND		ug/L	525.2	D	0.0000		
	DIMETHOATE	ND		ug/L	525.2	0	0.00000.0		
	FENARIMOL	ND		ug/L	525.2	O	000000		
	HEXAZINONE	ND		ug/L	525.2	0	000000		
	MALATHION	` ND		ug/L	525.2	0	.05000		
	METALAXYL	ND		ug/L	525.2	0	.10000		
	METHIDATHINON	ND		ug/L	525.2	o	.50000		
	METHYL PARATHION	ND		ug/L	525.2	o	.00000		
	MEVINPHOS	ND		ug/L	525.2		.00000		
	MYCLOBUTANIL	ND		ug/L	525.2		.50000		
	NAPROPAMIDE	. ND		ug/L	525.2		.00000		
	PARATHION-ETHYL	ND		ug/L	525.2		.05000		
	PHOSMET	ND		ug/L	525.2		.10000		
	PROPARGITE	ND		ug/L	525.2		.00000		
	TRIADIMEFON	ND		ug/L	525.2		.00000		
	TRIFLUMIZOLE	ND		ug/L	525.2		.00000		
531_081031	3-HYDROXYCARBOFURAN	ND		ug/L	531.2	0.	.50000	MB	
	ALDICARB	ND		ug/L	531.2	0.	.25000		
	ALDICARB SULFONE	ND		ug/L	531.2	0.	.40000		
	ALDICARB SULFOXIDE	ND		ug/L	531.2	0.	.25000		
	BDMC (SURR)	110		%	531.2	0.	.00000		
	CARBARYL	ND		ug/L	531.2	0.	.50000		
	CARBOFURAN	ND		ug/L	531.2	0.	45000		
	METHIOCARB	ND		ug/L	531.2	1.	.00000		
	METHOMYL	ND		ug/L	531.2	0.	25000		
	OXYMAL	ND		ug/L	531.2	1.	00000		
	PROPOXUR (BAYGON)	ND		ug/L	531.2	0.	25000		
004400									
ec_081103	ELECTRICAL CONDUCTIVITY	ND		uS/cm	SM2510 B	2.	50000	MB	
ec_081103	ELECTRICAL CONDUCTIVITY	ND		uS/cm	SM2510 B	2.	50000	МВ	
ec_ 08 1103	ELECTRICAL CONDUCTIVITY	ND		uS/cm	SM2510 B	2.	50000	MB	

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SAMPLE INDEPENDENT QUALITY CONTROL REPORT

Method Blank

Reference Number: 08-15517

		14	True			%	QC	
Batch	Analyte	Result	Value	Units	Method	Recovery Limits	Qualifier Type*	Comment
ec_081103	ELECTRICAL CONDUCTIVITY	ND		uS/cm	SM2510 B	2.50000) MB	4/1/4
OPHOS-081029	ORTHO-PHOSPHATE	ND		mg/L	SM4500-P F	0.10000) MB	
tds_081029	TOTAL DISSOLVED SOLIDS	ND		mg/L	SM2540 C	2.50000) MB	
lds_081029	TOTAL DISSOLVED SOLIDS	ND		mg/L	SM2540 C	2.50000	мв	
tds_081029	TOTAL DISSOLVED SOLIDS	ND		mg/L	SM2540 C	2.50000) MB	
tds_081029	TOTAL DISSOLVED SOLIDS	. ND		mg/L	SM2540 C	2.50000	МВ	
turb 081029	TURBIDITY							
IUID_00 1029	IURDIUIT	ND		NTU	180.1	0.02000	MB	

^{*}Notation:

[%] Recovery = (Result of Analysis)/(True Value) * 100

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SAMPLE INDEPENDENT **QUALITY CONTROL REPORT**

Quality Control Sample

Reference Number: 08-15517

		**							
			True			%		QC	
Batch	Analyte	Result	Value	Units	Method	Recovery	Limits	Qualifier Type*	Comment
200.7-0811D3A	HARDNESS	133	132.3	mg/L	200.7	101	80-120	QCS	
COD_081104	CHEMICAL OXYGEN DEMAND	125	133	mg/L	SM5220 D	94	80-120	QCS	
D081104A	BROMATE	0.0181	0.0182	mg/L	300.1	99	75-125	qcs	
ec_081103	ELECTRICAL CONDUCTIVITY	158	150.5	uS/cm	SM2510 B	105	80-120	qcs	
ec_081103	ELECTRICAL CONDUCTIVITY	158	150.5	uS/cm	SM2510 B	105	80-120	qcs	
ec_081103	ELECTRICAL CONDUCTIVITY	158	150.5	uS/cm	SM2510 B	105	80-120	qcs	
ec_081103	ELECTRICAL CONDUCTIVITY	158	150.5	uS/cm	SM2510 B	105	80-120	QCS	
I081029A	CHLORIDE	30.4	30.0	mg/L	300.0	101	80-120	QCS	
	NITRATE-N	2.58	2.50	mg/L	300.0	103	80-120		
OPHOS-081029	ORTHO-PHOSPHATE	0.47	0.49	mg/L	SM4500-P F	96	70-130	qcs	
ph_081029	HYDROGEN ION (pH)	7.97	8.00	pH Units	SM4500-H+ B	100	80-120	qcs	
ph_081029	HYDROGEN ION (pH)	8.05	8.00	pH Units	SM4500-H+ B	101	80-120	qcs	
turb_081029	TUDDIDATA								
mm_001058	TURBIDITY	0.95	1.00	UTM	180.1	95	70-130	QCS	

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Report Date: 12/10/2008 Reference Number: 08-15517 Page 1 of 5 Comments 밁 뎔 큠 출출 필 급 몱 ם 밁 밁 불물 를 Qualifier မ္ပ %RPD Limits 0-45 0.45 35 4 3 0.45 5 54 8 45 f 9-45 5 5 0-45 4.3 6. 4. 0.0 ۳. 0.0 0.0 3 3 0.0 4 9.0 A.1 4.1 4.0 Duplicate and Matrix Spike/Matrix Spike Duplicate Report QUALITY CONTROL REPORT Corporate Office 800.755.9295 • 360.757.1400 • 360.757.14021ax Bellingham WA 805 Orchard Dr Sulfe 4 • 98225 Merobelepy 350.671.0588 • 360.671.1577[ax тв СаСОЗ/L PH Units uS/cm uS/cm ES/CIII 턞 mg/L mg/L πgγ mg/L тgЛ щg/L 뺿 mg/L ě mg/L Duplicate 0.0126 Resul 7380 89.8 0.8 0.24 6.42 122 512 519 99 552 58 7 ţ 8 0.0130 Result 514 122 7440 161 89.4 10.8 0.25 6.38 95 554 516 7.1 98 5 33163 CHEMICAL OXYGEN DEMAND 33136 CHEMICAL OXYGEN DEMAND 32782 ELECTRICAL CONDUCTIVITY 33010 ELECTRICAL CONDUCTIVITY 32571 ELECTRICAL CONDUCTIVITY 32817 TOTAL DISSOLVED SOLIDS 32571 TOTAL DISSOLVED SOLIDS 32762 TOTAL DISSOLVED SOLIDS 32784 ORTHO-PHOSPHATE 32784 HYDROGEN ION (pH) 32892 HARDNESS HARDNESS 32775 NITRATE-N CHLORIDE CHLORIDE 32775 CHLORIDE 32565 BROMATE Sample Analyte 32853 32853 32817 **Duplicate OPHOS-081029** 200.7-081103A COD_081104

525_081105

Batch

EC_081103

1081029A

D081104A

TURB_081029

TDS_081029

PH_081029

%RPD = Relative Percent Difference

NA = indicates %RPD could not be calculated

Matrix Spike (MS)/Matrix Spike Duplicate (MSD) analyses are used to determine the accuracy (MS) and precision (MSD) of a analytical method in a given sample matrix. Therefore, the usertuiness of this report is limited to samples of similar matrices analyzed in the same analytical batch.



Duplicate

Batch

Comments Quelifier ဗ %RPD Limits 95 0 9.5 ä Duplicate Result 5 Result 32853 TURBIDITY Sample Analyte

Page 2 of 5 Reference Number: 08-15517 Report Date: 12/10/2008

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Reference Number: 08-15517 Report Date: 12/10/2008 Page 3 of 5

Matrix Spike

Matrix Spike	pike			Duplicate										
			Spika	Spike	Spike		Percent	Percent Recovery				8		
Batch	Sample Analyte	Residt	Result	Result	Conc	Units	MS	MSD	Limits	%RPD	Limits	Qualifier	Comments	
200.7-081103A						- Carallel							77.	
	32853 HARDNESS	161	230	229	69.5	mg CaCO3/L 99	66	86	80-120	1.5	08-9	5	LFM	
	32892 HARDNESS	89.4	158	162	69.5	mg/L	66	104	80-120	5.7	09-0	=	FW	
515_081031														
	29825 2,4-D	9	1.69		2	ug/L	88	¥2	65-135	٧	0-60	ב	LFM	
	29825 2,4,5-TP (SILVEX)	2	0.88		_	ug/L	88	NA A	65-135	¥	09-0	5	LFM	
-	29825 PENTACHLOROPHENOL	S	0.78		-	ug/L	82	NA A	65-135	۷ ۲	09-0	5	LFM	
	29825 DALAPON	皇	10.4		13	ug/L	80	Y.	65-135	Ϋ́	0-60	5	FM	
	28825 DINOSEB	용	1.66		2	ug/L	83	Ą	65-135	¥Z	8	"	LFM	
	29825 PICLORAM	2	1.13		_	√L n ∂ /L	113	¥	65-135	¥	9	٣	LFIZ	
	29825 DICAMBA	Ş	0.83	•	_	ng/L	22	Ą	65-135	¥	9 9	5	LPM	**
	29825 TOTAL (DCPA & Metabolites)	Q	6.0		_	ug/L	8	¥	65-135	Ą	09-0	<u>"</u>	LFM	
	29825 2,4 DB	9	7.9		60	ug/L	8	ΝA	65-135	¥2	09-0	5	LFM	
	29825 2,4,5 T	용	0.93		_	ug/L	89	ΝĀ	65-135	۷ ۷	9 0	<u>ה</u>	LFM	
	29825 BENTAZON	g	1.77		N.	ug/L	69	NA A	65-135	Υ	0-80	5	LFM	
	29825 DICHLORPROP	2	2.52		m	ng/L	84	NA	65-135	Υ	09-0	5	LFM	
	29825 ACIFLUORFEN	2	-		_	ng/L	9	NA A	65-135	Ą	080	5	LFM	
	29825 CHLORAMBEN	2	0.75		_	иgЛ	75	Y.	65-135	Ą	0-20	5	LFM	
	29825 2,4 - DCAA (SURR)	105	99			*		NA	70-130	¥	0-60	5	M#7	
		2	1.75		N	mg/L	88	NA A	65-135	Ϋ́	09-0	5	LFM	
	32909 2,4,5 - TP (SILVEX)	욮	0.82		_	mg/L	82	ΝĀ	65-135	Ϋ́	0-60	*	LFM	
	32909 PENTACHLOROPHENOL	2	0.78		_	ug/L	78	¥	65-135	¥	0-60	5	LFM	
		2	10,4		13	mg/L	80	NA A	65-135	Ϋ́	9-6	5	LFM	
		문	1.59		CNI.	mg/L	89	¥	65-135	Ϋ́	09-0	5	LFM	
		문	0.92		_	mg/L	92	¥.	65-135	Y Y	99	5	LFM	
	32809 DICAMBA	S	98.0		_	ug/L	98	¥	65-135	Š	9	5	LFM	
		9	6:0		_	ug/L	26	NA NA	65-135	Ą	08-0	ב	LFM	
	32909 2,4 DB	용	7.37		.	ug/L	92	NA	65-135	۲	09-0	5	LFM	
	32909 2,4,5T	2	0.83		_	ng/L	83	¥	65-135	ΥN	09-0	5	TEM .	
	32909 BENTAZON	9	1.8		~	ng/L	8	Z.	65-135	¥	09-0	"	LFM	
	32909 DICHLORPROP	S	2.41			ug/L	90	NA	65-135	Ā	09-0	<u>"</u>	LFM	
	32909 ACIFLUORFEN	Q Z	0.92		_	₩.	92	NA A	65-135	¥2	09-0	<u>"</u>	N.	
	32909 CHLORAMBEN	9	0.73		_	ηĝγ	73	¥.	65-135	Š	0-20	5	LFM.	
	32809 2,4 - DCAA (SURR)	2 8	69			%		NA A	70-130	¥	0 9 0	5	-FM	
525_081105														
	32424 BISPHENOL-A	Ş	က		ĽΩ	ug/L	400	NA	70-130	ž	0-20	"	LFM	

%RPD = Refative Percent Difference

NA = Indicates %RPD could not be calculated
Matrix Spike (MSJ/Matrix Spike Duplicate (MSD) and prediston (MSD) of a analytical method in a given sample matrix. Therefore, the usefulness of this raport is limited to samples of samples of samples of



Reference Number: 08-15517 Page 4 of 5

Report Date: 12/10/2008

Matrix Spike	pike				Duplicate									
				Spike	Spike	Spike	~-	Percent Recovery	SCOVERY				90	
Batch	Sample		Result	Result	Result (Солс	Units		MSD	Limits	%RPD	Umits	Qualifier	Comments
	32424	SIMAZINE	S	_			ug/L.	9	NA	70-130	Ą.	0-60	LFM	
	32424	DIAZINON	2	0.93	•	_		93	¥.	70-130	¥	09-0	E.F.	_
	32424	LINDANE (BHC - GAMMA)	Q	0.97	`	_	ng/L	97	ΝĀ	70-130	Ą	09-0	LFM	
	32424	. 4,4-DDD	Q	1.15	•	_	ug/L 1	115	¥	70-130	¥.	09-0	LFM	
	32424	4,4-DDE	Q	1.12	•	_	ug/L 1	112	¥	70-130	¥	09-0	LFM	_
	32424	4,4-DDT	민	1.1	•	_	-	110	¥	70-130	ž	0-90	LFM	_
	32424	MALATHION	무	1.23	•	_		123	Ā	70-130	۲ ۲	98-0	Ę	_
	32424	PARATHION-ETHYL	9	1.3	•	_	-	130	NA A	70-130	¥ X	9-60	LFM	_
	32424	1,3-DIMETHYL-2-NITROBENZENE (Surr	26	79			%	_	NA NA	70-130	¥	09-0	LFM	
	32424	PYRENE-D10 (Surr)	8	06			%	_	NA	70-130	¥	09-0	LFIX	_
	32424	PERYLENE-D12 (Sum)	91	66		٠,	%		NA	70-130	₹	09-0	LFIM	Α.
	32424	TRIPHENYLPHOSPHATE (Sur)	120	121			*		Ą	70-130	ş	08-0	LFM	
525X 081105														
	32424	PROPARGITE	2	2.4	•••	n.	ng/L	120	Y.	70-130	¥	0-20	NE NE	
	32424	METALAXYL	2	2.28		ru.		114	Ā	70-130	¥	0-20	E.	_
	32424	NAPROPAMIDE	9	1,16	•	_	·	116	¥	70-130	¥	0-20	LFI	
	32424	1-NAPHTHALENEACETAMIDE	문	2.26	.,	-		113	AN	70-130	¥	0-50	LFM	
	32424	FENARIMOL	2	1.5B	•	_		158	Y.	70-130	Y Y	0-20	LFM	
	32424	MEVINPHOS	Q	1.16	•	_	ug/L.	116	NA	70-130	¥	0-20	Ę	
	32424	CHLORPYRIFOS	皇	1.08		_		108	¥.	70-130	¥	05-0	LFM	_
	32424	DICOFOL	S	3.06	.,	N.		153	Ā	70-130	¥	0-20	HR LFI	
	32424	MALATHION	2	1.15	•	_		115	NA A	70-130	¥	0-60	LFM	
	32424	PARATHION-ETHYL	9	1.13	,	_		113	¥	70-130	Ą	09-0	LFM	_
	32424	PHOSMET	Q	_{co}	.,	.		150	NA	70-130	¥	0-20	HR LFM	_
	32424	TRIADIMEFON	9	1.18	•	_	ug/L 1	118	¥	70-130	¥	0-20	LFM	•
	32424	TRIFLUMIZOLE	R	0.77	••	N	ug/L s	88	NA	70-130	¥	0-50	N1 LFM	_
	32424	METHIDATHINON	오	3.02	••	ο.		151	Ą	70-130	Ą	0-20		_
	32424	MYCLOBUTANIL	9	2.5	••	N	ng/u.	125	¥	70-130	Ā	0-50	LFM	_
	32424	HEXAZINONE	Q	1.29	•	_		129	AA.	70-130	Ą.	0-20	LFI	
531_081031														
l	31657	OXYMAL	Q	14.3	13	5	mg/L 6	92	87	70-130	8.5	0-20	LFM	
	31657	CARBOFURAN	S S	14	12.7	5		93	85	70-130	5.7	0-20	LFM	
	31657	ALDICARB SULFOXIDE	P	15.6	14.7	9	mg/L 1	104	98	70-130	5.9	0-20	LFM	_
	31657	ALDICARB SULFONE	2	13.9	13.6	5		93	.	70-130	2:5	0-20	LFM	•
	31657	METHOMYL	Q.	13.7		5	ug/L {		87	70-130	5,2	0-30	Ę	•
	31657	3-HYDROXYCARBOFURAN	Ð	13.6	Oi	15		9	£	70-130	10.9	0-20	N±1	_

%RPD = Relative Percent Difference

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Matrix Spike (MSJ/Matrix Spike Duplicate (MSD) analyses are used to determine the accuracy (MS) and precision (MSD) of a analytical method in a given sample metrix. Therefore, the usefulness of this report is limited to samples of similar matrices analyzed in the same analytical batch.



Page 5 of 5 Reference Number: 08-15517

Comments LFM LFM E LFM Ē Ē Ē Ē Ē Ē E Ę Ē Ę, Ę Ę 돌 Ę FE Ē Ξ <u>F</u> Ę ဗ္ဗ Limits 8 9 99 S-20 8 ş 0-50 8 0-50 650 50 55 20 9-50 0.50 0-50 8 9 9 9-0 9-0 ş 8 8 ş ¥ ¥ 3,5 ž ≨ ž ž ¥ ş ≨ ₹ ٤ ¥ ž 6 9 8 7 § § ž 70-130 70-130 70-130 70-130 70-130 80-120 80-120 75-125 80-120 80-120 **B0-120** 80-120 80-120 70-130 70-130 70-130 70-130 70-130 70-130 70-130 70-130 70-130 70-130 80-120 80-120 80-120 70-130 Percent Recovery 106 108 **50** ş Š ž ¥ ¥ ž ž Š ≨ ≨ 100 둳 505 102 113 108 5 5 ¥ 107 30 90 100 100 S 듣 117 107 8 4 100 6 얆 mg/L пgЛ 뼥 퉏 mg/L ng/L 퉏 mg/L mg/L щg/L Ag/ 셤 βľ J₀ 7 υg/L ď 0.010 20.00 20.00 8 8 8 9. 20 20 10 5 5 5 5 5 2 5 5 9 8 Duplicate Result Spike 5,3 14.8 11.7 건 61 8 99 0.010 28.5 9. 1.26 10.5 10.2 1.3 15.1 12.9 13.4 11.7 10.7 30.1 9.2 9.4 9.4 8.5 2.7 69 8 3 87 10.8 0.25 9 88 7. 2 9 용 물 물 9 9 물 물 물 5 5 5 8 5 5 5 8 9 9 5 운 7.1 CHEMICAL OXYGEN DEMAND 3-HYDROXYCARBOFURAN ALDICARB SULFOXIDE PROPOXUR (BAYGON) PROPOXUR (BAYGON) 32784 ORTHO-PHOSPHATE ALDICARB SULFONE 32780 CARBOFURAN METHIOCARB BDMC (SURR) METHIOCARB BDMC (SURR) METHOMYL NITRATE-N **NITRATE-N** CARBARYL CARBARYL CHLORIDE CHLORIDE CHLORIDE ALDICARB ALDICARB 31852 BROMATE OXYMAL Analyte 32853 32853 31657 31657 31857 32780 32780 32780 32780 32780 32780 32780 32780 32782 33163 33171 32775 32775 Sample 31657 32780 32780 32820 33136 32817 31657 Matrix Spike OPHOS-081029 COD_081104 D081104A 1081029A Betch

[%]RPD = Relative Percent Difference

NA = Indicates %RPD could not be calculated

Metrix Spike (MS)/Matrix Spike Duplicate (MSD) analyses are used to determine the accuracy (MS) and precision (MSD) of a analytical method in a given sample matrix. Therefore, the usefulness of this report is limited to samples of similar matrices analyzed in the same analytical batch



N1

Page 1 of 1

Qualifier Definitions

Reference Number: 08-15517

Report Date: 12/10/08

Qualifier	Definition
B1	The source of the contamination has been identified as a contaminate in the lab purified water. Data for this compound is suspect if reported.
HR	High QCS recovery due to increased detector response No sample dectections, therefore, no further action taken for this analysis set.
ME	Matrix spike shows a possible matrix induced bias. The LFB was within acceptance limits, results for this compound are suspect.

Acceptance limits have not been established, the limits listed are for guidance only.







QUALITY CONTROL REPORT SURROGATE REPORT

Reference Number: 08-15517

	·	1 m			
Lab No	Analyte	Result Qualifier	Units	Method	Limit
531_081031 32780	BDMC (SURR)	83	%	531.2	
531_081031 32781	BDMC (SURR)	79	%	531.2	
531_081031 32782	BDMC (SURR)	75	%	531.2	
515_081031 32783 525_081105	2,4 - DCAA (SURR)	96	%	5 15 .1	Acceptance Range is 70 - 130%
32783	1,3-DIMETHYL-2-NITROBENZENE (Surr) PYRENE-D10 (Surr) PERYLENE-D12 (Surr)	79 90 84	% % %	525.2	Acceptance Range is 70% to 130% Acceptance Range is 70% to 130% Acceptance Range is 70% to 130%
531_081031 32783	TRIPHENYLPHOSPHÄTE (Sum) BDMC (SURR)	75	%		Acceptance Range is 70% to 130%
515 081031		75	· %	531.2	
32784 525_081105	2,4 - DCAA (SURR)	94	%	515.1	Acceptance Range is 70 - 130%
32784	1,3-DIMETHYL-2-NITROBENZENE (Surr) PYRENE-D10 (Surr) PERYLENE-D12 (Surr)	80 89 ** 82	% % %	5 2 5.2	Acceptance Range is 70% to 130% Acceptance Range is 70% to 130% Acceptance Range is 70% to 130%
531_081031 32784	TRIPHENYLPHOSPHATE (Surr) BDMC (SURR)	113 81	%	531.2	Acceptance Range is 70% to 130%
	•			~~	

^{*}Notation:

Chain of Custody / Analysis Request Be specific in analysis requests.
 Check off analyses to be performed for Š Sample Receipt Request (Must include FAX or Email) Ship Address: 810 S Main Street Report to: Relinquished by Project Phone: Use one line per sample (Location). Enter number of containers. each sample Location. 1-1 1-3 1-2 Locker/Hall-weithout Walla Walla Basin Watershed Cour 541.938-2170 FAX: BERRY LIVEY Dakar Milton-Freews St. Baker Have - Wentland#2 HALL-WONT BANK # 3 racher OBS # Lockes Obs #2 Locher Obs #3 OR Zip: 97862 Date Phone: 9382170 Turn Around Time Required

X Standard Time 읓 B# 60 Phone: Card#: P.O.#: ∏ Visa Address Quickest (100% surcharge Emergency (Phone Call Req. Half-time (50% surcharge) Received by Melrix 810 S Main Street Walla Walla Basin Watershed Counc Milton-Freewa St. ₹ 6 10/28/8/9:40 10,28/08 10,41 10/28/0 8:50 10/28/82 10:06 Date 2:11 80/87/2 (Please complete all applicable shaded sections) FA Time _ } Nitrate 유 TDS, CI. ZIÞ: 1929/08 O-Phos.pH,Turb,Ec 97862 Date Hardness Analyses Requested ã X Other Check Regulatory Program Time RCRA / CERCLA Clean Water Act Safe Drinking Water Act For Lab Use Only troy, bakes@wwbuck, are Bromate Custody seals intact Sample temp 2 C satisfactory Samples received intact 525(Hexazinone) 515 805 W. Orchard Dr. Suite 4 Bellingham, WA 98225 1620 S. Walnut St. Burlington, WA 98233 1.800.755.9295 Number of Containers ANALYTICAL LABORATORIES Total Containers No 525 Page Special Instructions Conditions on Receipt \es CO004495 4,495 8 ۵, Z

08-15517

D9-15517

Chain of custody & labels agree



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January 28, 2009

Page 1 of 1

Mr. Troy Baker Walla Walla Basin Watershed Council 810 S Main Street Milton-Freewater, OR 97862

RE: 08-17751 - Locher Road Recharge Sites/Hall Wetland

Dear Mr. Troy Baker,

Your project: Locher Road Recharge Sites/Hall Wetland, was received on Wednesday December 17, 2008. The following comments are reported for your project:

Sample 37231 - BisPhenol-A was detected in the EPA Method 525.2 analysis, estimated at 1.7 ug/L.

If you have questions phone me at 800 755-9295.

Respettfully Submitted,

Lawrence J Henderson, PhD Director of Laboratories

Enclosures Data Report



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Page 1 of 3

Data Report

Client Name: Walla Walla Basin Watershed Council

810 S Main Street

Milton-Freewater, OR 97862

Reference Number: 08-17751

Project: Locher Road Recharge Sites/Hall W.

Report Date: 1/28/09 Date Received: 12/17/08

Peer Review:

Sample Description: L-1 - Locher Rd

Sample Date: 12/16/08

Latin	umber: 3/230						Colle	cted By:	Unkno	WN	
CAS ID#	Parameter	Result	PQL	MDL	Units	DF	Method	Analyzed	Anely	st Batch	Comment
E-10139	HYDROGEN ION (pH)	6.91			pH Units	1	SM4500-H+ B	12/17/08	MAK	PH_081217	
14797-55-8	NITRATE-N	5.85	0.100	0.015	mg/L	1	300.0	12/17/06	8.7	ID61217A	
16887-00-6	CHLORIDE	6.1	0.1	0.012	mg/L	1	300.0	12/17/06	LE	1061217A	
E-10173	TOTAL DISSOLVED SOLIDS	257	10	6	mg/L	1	SM2540 C	12/19/08	CCN	TDS_051219	
14265-44-2	ORTHO-PHOSPHATE	0.19	0.01	0.002	mg/L	1	SM4500-P F	12/17/08	50	OPHOS-081217	
E-10184	ELECTRICAL CONDUCTIVITY	396	10		uS/cm	1	SM2510 B	12/23/06	CCN	EC_081223	
E-10617	TURBIDITY	0.70	0.05	0.02	NTU	1	180,1	12/17/08	MAK	TURB_061217	
15541-45-4	BROMATE	ND	0.005	0.0016	mg/L	1	300.1	12/30/08	MVP	D081230A	
E-11778	HARDNESS	150	3.30	0.055	mg CaCO3/L	1	200.7	12/23/08	BJ.	200.7-081223A	
E-10117	CHEMICAL OXYGEN DEMAND	ND	8	2	mg/L	1	SM5220 D	12/29/08	MAK	COD_081229	

			_
Sample i	Description:	L-2 - Locher	Rd

Sample Date: 12/16/08

Lab N	umber: 37231						Colle	cted By:	Unkno	wn	
CAS ID#	Parameter	Result	PQL	MDL	Units	DF	Method	Analyzed	Analy	st Baich	Comment
E-10139	HYDROGEN ION (pH)	6.87			pH Units	1	SM4500-H+ B	12/17/08	MAK	PH_081217	
14797-55-8	NITRATE-N	3.56	0.100	0.015	mg/L	1	300.0	12/17/08	BJ	1981217A	
16887-00-6	CHLORIDE	5.2	0.1	0.012	mg/L	1	300.0	12/17/08	BJ	1081217A	
E-101 7 3	TOTAL DISSOLVED SOLIDS	184	10	6	mg/L	1	SM2540 C	12/10/08	CCN	TDS_081219	
14265-44-2	ORTHO-PHOSPHATE	0.19	0.01	0.002	mg/L	,	SM4500-P F	12/17/08	SQ	OPHOS-081217	
-10184	ELECTRICAL CONDUCTIVITY	284	10		uS/cm	t	SM2510 B	12/23/08	CGN	EC_081223	
-10617	TURBIDITY	2.03	0.05	0.02	NTU	1	180.1	12/17/08	MAK	TURB_081217	
5541-45-4	BROMATE	ND	0.005	0.0016	mg/L	1	300.1	12/30/06	MVP	D081230A	
-11778	HARDNESS	105	3.30	0.055	mg CaCO3/L	1	200.7	12/23/08	BI	200.7-081223A	
-10117	CHEMICAL OXYGEN DÉMAND	ND	В	2	mg/L	1	SM5220 D	12/29/08	MAK	COD_081220	

ND = Not detected above the listed practical quantitation limit (PQL) or not above the Mathod Detection Limit (MDL), if requested.

PQL = Practical Quantitation Limit is the lowest level that can be achieved within specified limits of precision and accuracy during routine laboratory operating conditions.



Sample Description: L-3 - Locher Rd

Page 2 of 3 Reference Number: 08-17751 Report Date: 1/28/09

Data Report

Sample Date: 12/16/08 Lab Number: 37232 Collected By: Unknown CAS ID# Parameter Result POL MDL Units DF Method Analyzed Analyst Batch Comment E-10139 HYDROGEN ION (pH) 6.91 pH Units SM4500-H+ B 12/17/08 PH_081217 14797-55-8 NITRATE-N 2.11 0.100 0.015 mg/L 300.0 12/17/08 1061217A 16887-00-6 CHLORIDE 2.2 0.012 0.1 mg/L 1 300.0 12/17/08 BJ 8081217A E-10173 **TOTAL DISSOLVED SOLIDS** 125 10 mg/L SM2540 C 12/19/06 TDS_081219 14265-44-2 ORTHO-PHOSPHATE 0.15 0.01 0.002 ma/L SM450D-P F 12/17/08 OPHOS-081217 E-10184 **ELECTRICAL CONDUCTIVITY** 149 10 uS/cm SM2510 B 12/23/08 CCN EC_081223 E-10817 TURBIDITY 4.32 0.05 NTU 0.02 180.1 12/17/08 TURB_081217 15541-45-4 BROMATE ND 0.005 0.0016 mg/L 300.1 12/30/08 D081230A E-11778 HARDNESS 53.7 3.30 0.055 mg 200.7 12/23/08 200.7-081223A · CaCO3/L E-10117 **CHEMICAL OXYGEN DEMAND** ND 2 . mg/L SM5220 D 12/29/08 COD_081229



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Page 1 of 2

SYNTHETIC ORGANIC COMPOUNDS (SOC) REPORT

Client Name: Walla Walla Basin Watershed Council

810 S Main Street

Milton-Freewater, OR 97862

System Name:

System ID Number:

DOH Source Number: **Multiple Sources:**

Sample Type:

Sample Purpose: Investigative or Other

Sample Location: Locher Rd

County:

Sampled By: Unknown

Sampler Phone:

Reference Number: 08-17751

Project: Locher Road Recharge Sites/Ha

Field ID: L-1

Lab Number: 046-37230 Date Collected: 12/16/08 10:45 Date Extracted: 525_081222 Date Analyzed: 12/23/08

Report Date: 1/28/09 Analyst: CO

Peer Review:

EPA Method 525.2 For State Drinking Water Compliance

DOH#	COMPOUNDS	RESULTS	UNITS	SRL	Trigger	MCL	COMMENT
	EPA Regulated						
33	ENDRIN	ND	ug/L	0.02	0.02	2	
34	LINDANE (BHC - GAMMA)	ND	ug/L	0.04	0.04	0.2	
35	METHOXYCHLOR	ND	ug/L	0.2	0.2	40	
117	ALACHLOR	ND	ug/L	0.4	0.4	2	
119	ATRAZINE	ND	ug/L	0.2	0.2	3	
120	BENZO(A)PYRENE	ND	ug/L	0.04	0.04	0.2	
122	CHLORDANE, TECHNICAL	ND	ug/L	0.4	0.4	2	
124	DI(ETHYLHEXYL)-ADIPATE	ND	ug/L	1.3	1.3	400	
125	DI(ETHYLHEXYL)-PHTHALATE	ND	ug/L	1.3	1.3	6	
126	HEPTACHLOR	ND	ug/L	0.08	0.08	0.4	
127	HEPTACHLOR EPOXIDE	ND	ug/L	0.04	0.04	0.2	
128	HEXACHLOROBENZENE	ND	ug/L ·	0.2	0.2	1	
129	HEXACHLOROCYCLO-PENTADIENE	ND	ug/L	0.2	0.2	50	
133	SIMAZINE	ND	ug/L	0.15	0.15	4	
134	PENTACHLOROPHENOL	ND	ug/L	0.4	0.08	1	
Ì	EPA Unregulated						
118	ALDRIN	ND	ug/L	0.2	0.2	1	
121	BUTACHLOR	ND	ug/t.	0.4	0.4		
123	DIELDRIN	ND	ug/L	0.2	0.2		
130	METOLACHLOR	ND	ug/L	1.0	1.0		
131	METRIBUZIN	ND	ug/L	0.2	0.2		
132	PROPACHLOR	ND	ug/L	0.2	0.2		
[
	PROMETON	ND	ug/L	0.1			
179	BROMACIL	0.09 J	ug/L	0.2	0.2	j	Fleld dup - 0.09 ug/L

If a compound is detected > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.

MCL (Maximum Conteminant Level) maximum permissible level of a conteminant in water established by EPA; Federal Action Levels are 0.015 mg/L for Lead and 1.3 mg/L for Copper. Sodium has a recommended limit of 20 mg/L. A blank MCL value indicates a level is not currently established.

Trigger Level: DOH Drinking Water Response level. Systems with compounds detected in excess of this level are required to take additional samples. Contact your regional DOH office.

ND (Not Detected); indicates that the parameter was not detected above the State Reporting Limit (SRL).



Lab Number: 046-37230 Report Date: 1/28/09 15:44

SYNTHETIC ORGANIC COMPOUNDS (SOC) REPORT

DOH#	COMPOUNDS	RESULTS	UNITS	SRL	Trigger	MCL	COMMENT
190	TERBACIL	ND	ug/L	0.2	0.2		
202	DIAZINON	ND	ug/L	0.2	0.2		Unstable in Acidified Sample Matrix
208	EPTC	ND	ug/L	0.3	0.3		
232	4,4DDD	ND	ug/L	0.2	0.2		
233	4,4-DDE	ND	ug/L	0.2	0.2		
234	4.4-DDT	ND	ug/L	0.2	0.2		·
236	CYANAZINE	ND	ug/L	0.2	0.2		
239	MALATHION	ND	ug/L	0.2	0.2		
240	PARATHION	NĐ	ug/L	0.2	0.2		
243	TRIFLURALIN	ND	ug/L	0.2	0.2		
	- PAHs			ŀ			
96	NAPTHALENE	ND	ug/L	0.1	0.1		
254	FLUORENE	ND	ug/L	0.2	0.2	1	
245	ACENAPHTHENE	ND	ug/L	0.2	0.2		
246	ANTHRACENE	ND	ug/L	0.2	0.2		
247	BENZ(A)ANTHRACENE	ŃD	ug/L	0.1	0.1		,
249	BENZO(G,H,I)PERYLENE	ND	ug/L	0.2	0.2		
251	CHRYSENE	ND	ug/L	0.2	0.2		
252	DIBENZO(A,H)ANTHRACENE	ND	ug/L	0.2	0.2	•	
253	FLUORANTHENE	ND	ug/L	0.2	0.2		
255	INDENO(1,2,3-CD)PYRENE	ND	ug/L	0.2	0.2		
256	PHENANTHRENE	ND	ug/L	0.2	0.2		,
į	- Phthalates						
258	BENZYL BUTYL PHTHALATE	ND	ug/L	0.6	0.6		·
259	DI-N-BUTYL PHTHALATE	ND	ug/L	0.6	0.6		
260	DIETHYL PHTHALATE	ND	ug/L	0.6	0.6		
261	DIMETHYL PHTHALATE	ND	ug/L	0.6	0.6		
	Other Compounds						
0	HEXAZINONE (Velpar)	ND	ug/L	0.1			
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NOTES:



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SYNTHETIC ORGANIC COMPOUNDS (SOC) REPORT

Client Name: Walla Walla Basin Watershe-

810 S Main Street

Milton-Freewater, OR 97862

System Name:

System ID Number: DOH Source Number:

Multiple Sources:

Sample Type:

Sample Purpose: Investigative or Other

Sample Location: Locher Rd

County:

Sampled By: Unknown

Sampler Phone:

Reference Number: 08-17751

Project: Locher Road Recharge Sites/Ha

Field ID: L-2

Lab Number: 046-37231 Date Collected: 12/16/08 10:45 Date Extracted: 525 081222 Date Analyzed: 12/23/08

Report Date: 1/7/09

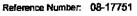
Analyst: Peer Review:

EPA Method 525.2 For State Drinking Water Compliance

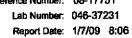
DOH#	COMPOUNDS	RESULTS	UNITS	SRL	Trigger	MCL	COMMENT
	EPA Regulated			1			
33	ENDRIN	ND	ug/L	0.02	0.02	2	
34	LINDANE (BHC - GAMMA)	ND	ug/L	0.04	0.04	0.2	
35	METHOXYCHLOR	ND	ug/L	0.2	0.2	40	
117	ALACHLOR	ND	ug/L	0.4	0.4	2	
119	ATRAZINE	ND	ug/L	0.2	0.2	3	
120	BENZO(A)PYRENE	ND	ug/L	0.04	0.04	0.2	
122	CHLORDANE, TECHNICAL	ND	ug/L	0.4	0.4	2	
124	DI(ETHYLHEXYL)-ADIPATE	ND	ug/L	1.3	1.3	400	
125	DI(ETHYLHEXYL)-PHTHALATE	ND	ug/L	1.3	1.3	6	
126	HEPTACHLOR	ND	ug/L	0.08	0.08	0.4	
127	HEPTACHLOR EPOXIDE	ND	ug/L	0.04	0.04	0.2	İ
128	HEXACHLOROBENZENE	ND	ug/L	0.2	0.2	1	
129	HEXACHLOROCYCLO-PENTADIENE	ND	ug/L	0.2	0.2	50	
133	SIMAZINE	ND	ug/L	0.15	0.15	4	
134	PENTACHLOROPHENOL	ND	ug/L	0.4	0.08	1	screening only / compliance by 515.1
	EPA Unregulated						
118	ALDRIN	ND	ug/L	0.2	0.2		
121	BUTACHLOR	ND	ug/L	0.4	0.4		
123	DIELDRIN	ND	ug/L	0.2	0.2	1	
130	METOLACHLOR	ND	ug/L	1.0	1.0		
131	METRIBUZIN	ND	ug/L	0.2	0.2		
132	PROPACHLOR	ND	ug/L	0.2	0.2		
	State Unregulated - Other						
179	BROMACIL	ND	ug/L	0.2	0.2		
190	TERBACIL	ND	ug/L	0.2	0.2		.]

specified increased monitoring trequencies may occur per DOH.

s contaminant in water established by EPA; Federal Action Levels are 0.015 mg/L for Lead and 1.3 mg/L for Copper. Sodium has a recommended limit of 20 mg/L. A MCL (Maximum Contentions) the state of the state of the state of the contention of the state of



Page 2 of 2



SYNTHETIC ORGANIC COMPOUNDS (SOC) REPORT

DOH#	COMPOUNDS	RESULTS	UNITS	SRL	Trigger	MCL	COMMENT
202	DIAZINON	ND	ug/L	0.2	0.2		Unstable in Acidified Sample Matrix
208	EPTC	ND	ug/L	0.3	0.3		
232	4.4-DDD	ND	ug/L	0.2	0.2		
233	4.4-DDE	ND	ug/L	0.2	0.2		
234	4,4-DDT	ND	ug/L	0.2	0.2	Ì	
236	CYANAZINE	ND	ug/L	0.2	0.2]	Qualitative Analysis Only
239	MALATHION	ND	ug/L	0.2	0.2		
240	PARATHION	ND	ug/L	0.2	0.2		
243	TRIFLURALIN	ND	ug/L	0.2	0.2	1	·
	- PAHs						
96	NAPTHALENE	ND	ug/L	0.1	0.1		
254	FLUORENE	ND	ug/L	0.1	0.1		
244	ACENAPHTHYLENE	ND	1 -	0.2	0.2		
245	ACENAPHTHENE	ND .	ug/L	0.2	0.2		
245	ANTHRACENE	ND .	ug/L	0.2	0.2		
247	BENZ(A)ANTHRACENE	ND	ug/L	0.2	0.2		•
248	BENZO(B)FLUORANTHENE	ND	ug/t.	0.1	0.1	j	
249	BENZO(G,H,I)PERYLENE	ND	ug/L	0.2	1		
250	BENZO(K)FLUORANTHENE	1	ug/L		0.2		
250 251	CHRYSENE	ND	ug/L	0.2	0.2		
252		ND	ug/L	0.2	0.2		
	DIBENZO(A,H)ANTHRACENE	ND	ug/L	0.2	0.2		
253	FLUORANTHENE	ND	ug/L	0.2	0.2	-	
255	INDENO(1,2,3-CD)PYRENE	ND	ug/L	0.2	0.2		
256	PHENANTHRENE	ND	ug/L	0.2	0.2	1	
257	PYRENE	ND	ug/L	0.2	0.2		
	- Phthalates						
258	BENZYL BUTYL PHTHALATE	ND	ug/L	0.6	0.6		
259	DI-N-BUTYL PHTHALATE	ND	ug/L	0.6	0.6		
260	DIETHYL PHTHALATE	ND	ug/L	0.6	0.6		
261	DIMETHYL PHTHALATE	ND	ug/L	0.6	0.6		
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NOTES:
If a compound is detected > or = to the State Reporting Level, SRL, apacified increased monitoring frequencies may occur per DOH.

MCL (Maximum Contaminant Level) maximum permissible level of a contaminant in water established by EPA; Federal Action Levels are 0.015 mg/L for Lead and 1.3 mg/L for Copper. Sodium has a recommended limit of 20 mg/L. A blank MCL value indicates a level is not currently established.

Trigger Level: DOH Drinking Water Response level. Systems with compounds detected in excess of this level are required to take additional samples. Contact your regional DOH office.

ND (Not Detected): indicates that the parameter was not detected above the State Reporting Limit (SRL).



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Page 1 of 2

SYNTHETIC ORGANIC COMPOUNDS (SOC) REPORT

Client Name: Walla Walla Basin Watershe-

810 S Main Street

Milton-Freewater, OR 97862

System Name:

System ID Number:

DOH Source Number:

Multiple Sources:

Sample Type:

Sample Purpose: Investigative or Other

Sample Location: Locher Rd

County:

Sampled By: Unknown

Sampler Phone:

Reference Number: 08-17751

Project: Locher Road Recharge Sites/Ha

Field ID: L-3

Lab Number: 046-37232

Date Collected: 12/16/08 10:25

Date Extracted: 525_081222

Date Analyzed: 12/23/08

Report Date: 12/31/08

Analyst: CO-

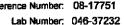
Peer Review:

DOH#	COMPOUNDS	RESULTS	UNITS	SRL	Trigger	MCL	COMMENT
	EPA Regulated		 		1		
33	ENDRIN	ND	ug/L	0.02	0.02	2	
34	LINDANE (BHC - GAMMA)	ND	ug/L	0.04	0.04	0.2	
35	METHOXYCHLOR	ND	ug/L	0.2	0.2	40	
117	ALACHLOR	ND	ug/L	0.4	0.4		
119	ATRAZINE	ND	ug/L	0.2	0.2	2 3	
120	BENZO(A)PYRENE	ND	ug/L	0.04	0.04	0.2	
122	CHLORDANE, TECHNICAL	ND	ug/L	0.4	0.4	2	
124	DI(ETHYLHEXYL)-ADIPATE	ND	ug/L	1.3	1.3	400	
125	DI(ETHYLHEXYL)-PHTHALATE	ND	ug/L	1.3	1.3	6	
126	HEPTACHLOR	ND	ug/L	0.08	0.08	0.4	
127	HEPTACHLOR EPOXIDE	ND	ug/L	0.04	0.04	0.2	·
128	HEXACHLOROBENZENE	ND	ug/L	0.2	0.2	1	
129	HEXACHLOROCYCLO-PENTADIENE	ND	ug/L	0.2	0.2	50	
133	SIMAZINE	ND	ug/L	0.15	0.15	4	
134	PENTACHLOROPHENOL	ND	ug/L	0.4	0.08	1	
	EPA Unregulated						
118	ALDRIN	ND	ug/L	0.2	0.2		
121	BUTACHLOR	ND	ug/L	0.4	0.4		
123	DIELDRIN	ND	ug/L	0.2	0.2		
130	METOLACHLOR	ND	ug/L	1.0	1.0		
131	METRIBUZIN	ND	ug/L	0.2	0.2		
132	PROPACHLOR	ND	ug/L	0.2	0.2		
	State Unregulated - Other						·
179	BROMACIL	ND	ug/L	0.2	0.2		
190	TERBACIL	ND	ug/L	0.2	0.2		

Tripper Level: DOH Drinking Water Response level. Systems with compounds detected in excess of this level are required to take additional samples. Cantact your regional DOH office.

ND (Not Datecled): Indicates that the parameter was not detected above the State Reporting Limit (SRL).





Report Date: 12/31/08 10:41

Page 2 of 2



SYNTHETIC ORGANIC COMPOUNDS (SOC) REPORT

OH#	COMPOUNDS	RESULTS	UNITS	SRL	Trigger	MCL	COMMENT
202	DIAZINON	ND	ug/L	0.2	0.2	1	Unstable in Acidified Sample Matrix
208	EPTC	ND	ug/L	0.3	0.3		
232	4.4-DDD	ND	ug/L	0.2	0.2	1	
233	4,4-DDE	ND	ug/L	0.2	0.2		
234	4,4-DDT	ND	ug/L	0.2	0.2		
236	CYANAZINE	ND	ug/L	0.2	0.2	1	
239	MALATHION	ND	ug/L	0.2	0.2		
240	PARATHION	ND	ug/L	0.2	0.2		
243	TRIFLURALIN	ND	ug/L	0.2	0.2	1	
245	- PAHs	1,5	Ogic	U.E	J	-	
				<u> </u>			
96	NAPTHALENE	ND	ug/L	D.1	0.1		•
254	FLUORENE	ND	ug/L	0.2	0.2	Parameter.	
245	ACENAPHTHENE	ND	ug/L	0.2	0.2		
246	ANTHRACENE	ND	ug/L	0.2	0.2	1	
247	BENZ(A)ANTHRACENE	ND	ug/L	0.1	0.1	1	
249	BENZO(G,H,I)PERYLENE	ND	ug/L	0.2	0.2		
	CHRYSENE	ND	ug/L	0.2	0.2	1	
1	DIBENZO(A,H)ANTHRACENE	ND	ug/L	0.2	0.2		
- 1	FLUORANTHENE	ND	ug/L	0.2	0.2		
	INDENO(1,2,3-CD)PYRENE	ND	ug/L	0.2	0.2		
256	PHENANTHRENE	ND	ug/L	0.2	0.2		·
	- Phthalates						
258	BENZYL BUTYL PHTHALATE	ND	ug/L	0.6	0.6		
	DI-N-BUTYL PHTHALATE	ND	ug/L	0.6	0.6	1	
	DIETHYL PHTHALATE	ND	ug/L	0.6	0.6		
- 1	DIMETHYL PHTHALATE	ND	ug/L	0.6	0.6		
			94.				
					1		
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-						1	
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NOTES:
If a compound is delected > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.

MCL (Maximum Contaminant Level) maximum permissible level of a contaminant in water established by EPA; Federal Action Levels are 0.015 mg/L for Lead and 1.3 mg/L for Copper. Sodium has a recommended limit of 20 mg/L. A blank MCL value indicates a layel is not currently established.

Trigger Level: DOH Drinking Water Response level. Systems with compounds detected in excess of this level are required to take additional samples. Contact your regional DOH office.

ND (Not Detected): Indicates that the parameter was not detected above the State Reporting Limit (SRL).



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WSDOE Lab C1251

DATA REPORT

Page 1 of 1

Client Name: Walla Walla Basin Watershed Council

810 S Main Street

Milton-Freewater, OR 97862

Reference Number: 08-17751

Project: Locher Road Recharge Sites

Lab Number: 37230

Field ID: L-1

Sample Description: Locher Rd

Matrix: Water

Sample Date: 12/16/08

Extraction Date: 12/22/08

Extraction Method: 3535

Report Date: 1/6/09

Date Analyzed: 12/23/08

Analyst: CO

Peer Review:

Analytical Method: 525.2

Batch: 525X_081222

CAS Compound **RESULT**

ND

Flag

UNITS PQL .

MDL

COMMENT D.F.

Other Compounds 51235-04- HEXAZINONE (Velpar)

0.1

0.14

1.00

Flags are data qualifiers. If there are data qualifiers on your report definitions can be found on an accompanying sheet.

NO - indicates the compound was not detected above the POL or MDL.

PQL = Practical Quantification Limit is the lowest level that can be achieved within specified limits of precision and accuracy during routine laboratory operating conditions. D.F. - Dilution Factor.



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WSDOE Lab C1251

DATA REPORT

Page 1 of 1

Client Name: Walla Walla Basin Watershed Council

810 S Main Street

Milton-Freewater, OR 97862

Reference Number: 08-17751

Project: Locher Road Recharge Sites

Lab Number: 37231

Field ID: L-2

Sample Description: Locher Rd

Matrix Water

Sample Date: 12/16/08

Extraction Date: 12/22/08

Extraction Method: 3535

Report Date: 1/6/09

Date Analyzed: 12/23/08

Anatyst: |QO

Peer Review: À

Analytical Method: 525.2

Batch: 525X 081222

COMMENT D.F. MDL UNITS PQL RESULT Flag CAS Compound Other Compounds 0.14 1.00 0.1 ND ug/L 51235-04 HEXAZINONE (Velpar)



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DATA REPORT

Page 1 of 1

Client Name: Walla Walla Basin Watershed Council

810 S Main Street

Milton-Freewater, OR 97862

Reference Number: 08-17751

Project: Locher Road Recharge Sites

Lab Number: 37232

Field ID: L-3

Sample Description: Locher Rd

Matrix: Water

Sample Date: 12/16/08

Extraction Date: 12/22/08

Extraction Method: 3535

Report Date: 1/6/09

Date Analyzed: 12/23/08

Analyst: CO

Peer Review: VML Analytical Method: 525.2

Batch: 525X 081222

CAS	Compound	RESULT	Flag	UNITS	PQL	MDL	D.F.	COMMENT		
E477E 0A	Other Compounds	MD			0.4	0.44			,	
51235-04-	HEXAZINONE (Velpar)	ND		ug/L	0.1	0.1^	1.0	00		



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Page 1 of 1

HERBICIDES IN DRINKING WATER

Client Name: Walla Walla Basin Watershe-

810 S Main Street

Milton-Freewater, OR 97862

Project:

Field ID: L-1

Sample Description: Locher Rd

Sampled By: Unknown

Sample Date: 12/16/08

Source Type:

Sampler Phone:

Reference Number: 08-17751

Project: Locher Road Recharge Sites/Ha

Lab Number: 37230

Report Date: 1/6/09 Date Analyzed: 12/31/08

Date Extracted: 515 081222

Analyst: ←ÇO

Peer Review: VM Analytical Method: 1515.1

Chlorophenoxy Herbicides

CAS	COMPOUND	RESULTS	UNITS	PQL	MDL.	MCL	COMMENT
EPA R	egulated						
94-75-7	2,4 - D	ND	ug/L	0.2	0.11	70	
93-72-1	2,4,5 - TP (SILVEX)	ND	ug/L	0.1	0.02	50	
87-86-5	PENTACHLOROPHENOL	ND	ug/L	0.1	0.044	1	
75-99-0	DALAPON	ND	ug/L	1.3	0.80	200	
88-85-7	DINOSEB ,	ND	ug/L	0.2	0.16	7	
1918-02-1	PICLORAM	ND	ug/L	0.1	0.089	500	
EPA U	nregulated						
1918-00-9	DICAMBA	ND	υ <u>α</u> /L	0.1	0.045		
State U	Inregulated						
1861-32-1	TOTAL (DCPA & Metabolites)	ND	ug/L	0.1	0.089		
E-14028	DCPA (ACID METABOLITES)	ND ·	ug/L	0.1	0.1		
94-82-6	2,4 DB	ND	ug/L	8.0	0.10		
93-76-5	2,4,5 T	ND	ug/L	0.1	0.044		
25057-89-0	BENTAZON	ND	ug/L	0.2	0.067		
120-36-5	DICHLORPROP	ND	ug/L	0.3	0.089		
50594-66-6	ACIFLUORFEN	ND	ug/L	0.1	0.089		
133-90-4	CHLORAMBEN	ND	ug/L	0.2	0.2		
51-36-5	3,5 - DICHLOROBENZOIC ACID	ND	nB\r	0.1	0.044		

NOTES:
If a compound is detected > or = to the State Reporting Level, SRL, specified increased showlong frequencies may occur per DOM.
MCL (Maximum Contaminant Level) maximum permissible level of a contaminant in water established by EPA; Federal Action Levels are 0.015 mg/L for Lead and 1.3 mg/L for Copper. Sodium has a recommended limit of 20 mg/L. A blank MCL value indicates a level is not currently established.
Trigger Level: DOH Orinking Water Response level. Systems with compounds defacted in excess of this level are required to take additional samples. Contact your regional DOH office.



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Page 1 of 1

HERBICIDES IN DRINKING WATER

Client Name: Walla Walla Basin Watershe-

810 S Main Street

Milton-Freewater, OR 97862

Project::

Field ID: L-2

Sample Description: Locher Rd

Sampled By: Unknown

Sample Date: 12/16/08

Source Type:

Sampler Phone:

Reference Number: 08-17751

Project: Locher Road Recharge Sites/Ha

Lab Number: 37231

Report Date: 1/6/09

Date Analyzed: 12/31/08

Date Extracted: 515 081222

Analyst:

Peer Review:

Analytical Method:

Chlorophenoxy Herbicides

CAS	COMPOUND	RESULTS	UNITS	PQL	MDL	MCL	COMMENT
EPA R	egulated						***************************************
14-75-7	2,4 - D	ND	ug/L	0.2	0.11	70	
3-72-1	2,4,5 - TP (SILVEX)	ND	ug/L	0.1	0.02	50	•
7-86-5	PENTACHLOROPHENOL	ND	ug/L	0.1	0.044	1	
5-99-0	DALAPON	ND	ug/L	1.3	0.80	200	
8-85-7	DINOSEB	ND	ug/L	0.2	0.16	7	
918-02-1	PICLORAM	ND	ug/L	0.1	0.089	500	
EPA U	nregulated						·
918-00-9	DICAMBA	ND	ug/L	0.1	0.045		
State l	Inregulated						
861-32-1	TOTAL (DCPA & Metabolites)	ND	ug/L	0.1	0.089		
-14028	DCPA (ACID METABOLITES)	ND %	ug/L	0.1	0.1		
4-82-6	2,4 DB	ND	ug/L	8.0	0.10		
3-76-5	2,4,5 T	ND	ug/L	0.1	0.044		
5057-89-0	BENTAZON	ND	ug/L	0.2	0.067		
20-36-5	DICHLORPROP	ND	ug/L	0.3	0.089		
0594-66-6	ACIFLUORFEN	ND	ug/L	0.1	0.089		
33-90-4	CHLORAMBEN	ND	ug/L	0.2	0.2		
1-36-5	3,5 - DICHLOROBENZOIC ACID	ND	ug/L	0.1	0.044		

ND (Not Detected): Indicates that the parameter was not detected above the State Reporting Limit (SRL).

If a compound is detected > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.

MCL (Maximum Contaminant Level) maximums permissible level of a contaminant in water established by EPA; Federal Action Levels are 0.015 mg/L for Lead and 1.3 mg/L for Copper. Sodium has a recommended limit of 20 mg/L, A blank MCL wake indicates a level is not currently established.

Trigger Level: DOH Drinking Water Response level. Systems with compounds detacted in excess of this level are required to take additional samples. Contact your regional DOH office.



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Page 1 of 1

HERBICIDES IN DRINKING WATER

Client Name: Walla Walla Basin Watershe

810 S Main Street

Milton-Freewater, OR 97862

Project:

Field ID: L-3

Sample Description: Locher Rd

Sampled By: Unknown Sample Date: 12/16/08

Source Type:

Sampler Phone:

Reference Number: 08-17751

Project: Locher Road Recharge Sites/Ha

Lab Number: 37232 Report Date: 1/6/09

Date Analyzed: 12/31/08 Date Extracted: 515 081222

Analyst: 40 Peer Review:

Analytical Method: 515.1

Chlorophenoxy Herbicides

CAS	COMPOUND	RESULTS	UNITS	PQL	- MDL	MCL	COMMENT
EPA R	egulated					·	
94-75-7	2,4 - D	ND	ug/L	0.2	0.11	70	
93-72-1	2,4,5 - TP (SILVEX)	ND	ug/L	0.1	0.02	50	
87-86-5	PENTACHLOROPHENOL	ND	u g/L	0.1	0.044	1	
75- 99 -0	DALAPON	ND	ug/L	1.3	0.80	200	
88-85-7	DINOSEB	ND	u g/L	0.2	0.16	7	
1918-02-1	PICLORAM	ND	ug/L	0.1	0.089	500	
EPA U	nregulated						
1918-00-9	DICAMBA	ND	ug/L	0.1	0.045		
State L	Inregulated						
1861-32-1	TOTAL (DCPA & Metabolites)	0.2	ug/L	0.1	0.089		
94-82-6	2,4 DB	ND 🤭	ug/L	8.0	0.10		
93-76-5	2,4,5 T	ND	u g/L	0.1	0.044		•
25057-89-0	BENTAZON	ND	ug/L	0.2	0.067		
120-36-5	DICHLORPROP	ND	⊔g/L	0.3	0.089		
50594-66-6	ACIFLUORFEN	NĐ	ug/L	0.1	0.089		
133-90-4	CHLORAMBEN	ND	ug/L	0.2	0.2		
51-36-5	3,5 - DICHLOROBENZOIC ACID	ND	υg/L	0.1	0.044		•

NOTES:

NOTES:
If a compound is detected > or a to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.
MCL (Maximum Contaminant Level) maximum permissible level of a contaminant in water established by EPA; Federal Action Levels are 0.015 mg/L for Lead and 1.3 mg/L for Copper. Sodium has a recommended limit of 20 mg/L. A blank MCL value indicates a level is not currently established.
Tripper Level: DOH Drinking Water Response level. Systems with compounds detacted in excess of this level are required to take additional samples. Contact your regional DOH office.
ND (Not Defacted): indicates that the parameter was not detected above the State Reporting Limit (SRL).



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Page 1 of 1

CARBAMATES IN DRINKING WATER

Client Name: Walla Walla Basin Watershe

810 S Main Street

Milton-Freewater, OR 97862

Project::

Field ID: L-2

Sample Description: Locher Rd

Sampled By: Unknown

Sample Date: 12/16/08

Source Type:

Sampler Phone:

Reference Number: 08-17751

Project: Locher Road Recharge Sites/Ha

Lab Number: 37231

Report Date: 12/22/08

Date Analyzed: 12/17/08

Date Extracted: 531_081217

Analyst: CQ

Peer Review:

Analytical Method: \$31.2

Carbamates

CAS	COMPOUND	RESULTS	UNITS	PQL '	MDL	MCL	COMMENT
EPA R	egulated			, ,,			
23135-22-0	OXYMAL	ND	ug/L	1.0	0.3	200	
1563-66-2	CARBOFURAN	NĐ	ug/L	1.0	0.2	40	
EPA U	nregulated						
1646-87-3	ALDICARB SULFOXIDE	ND	ug/L	1.0	0.3		
1646-88-4	ALDICARB SULFONE	ND	ng∕l	1.0	0.3		
16752-77-5	METHOMYL	ND	ug/L	1.0	0.3	-	
16655-82-6	3-HYDROXYCARBOFURAN	ND	ug/L	1.0	0.3		
116-06-3	ALDICARB	ND	ug/L	1.0	0.3		
53-25-2	CARBARYL	ND	ug/L	1.0	0.2		
State U	nregulated - Other						
114-26-1	PROPOXUR (BAYGON)	ND ''	ug/L	1.0	0.4		
2032-65-7	METHIOCARB	ND	ug/L	1.0	0.3		

NOTES;
If a compound is defected > or = to the State Reporting Level, SRL, spatified increased monitoring fraquencies may occur per DOH.

If a compound is defected > or = to the State Reporting Level, SRL, spatified increased monitoring fraquencies may occur per DOH.

MCL (Maximum Contaminant Level) maximum permissible level of a contaminant to water established by EPA; Federal Action Levels are 0.015 mg/L for Copper. Sodium has a recommended limit of 20 mg/L. A blank MCL value indicates a level is not currently established.

Trigger Level; DOH Orinking Water Response level. Systems with compounds detected in excess of this level are required to take additional samples. Contact your regional DOH office.

ND (Not Detected): indicates that the parameter was not detected above the State Reporting Limit (SRL).



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Page 1 of 1

CARBAMATES IN DRINKING WATER

Client Name: Walla Walla Basin Watershe

810 S Main Street

Milton-Freewater, OR 97862

Project::

Field ID: L-3

Sample Description: Locher Rd

Sampled By: Unknown

Sample Date: 12/16/08

Source Type:

Sampler Phone:

Reference Number: 08-17751

Project: Locher Road Recharge Sites/Ha

Lab Number: 37232

Report Date: 12/22/08 Date Analyzed: 12/17/08

Date Extracted: 531, 081217

Analyst: CO Peer Review:

Analytical Method: 531.2

Carbamates

CAS	COMPOUND	RESULTS	UNITS	PQL	MDL	MCL	COMMENT
EPA Re	egulated						
23135-22-0	OXYMAL.	ND	ug/L	1.0	0.3	200	
1563-66-2	CARBOFURAN	ND	nB/r	1.0	0.2	40	
EPA U	nregulated						
646-87-3	ALDICARB SULFOXIDE	ND	ug/L	1.0	0.3		
646-88-4	ALDICARB SULFONE	ND	ug/L	1.0	0.3		
6752-77-5	METHOMYL	ND	цg/L	1.0	0.3		
16655-82-6	3-HYDROXYCARBOFURAN	ND	ug/L	1.0	0.3		
16-06-3	ALDICARB	ND	ug/L	1.0	0.3		
3-25-2	CARBARYL	ND	ug/L	1.0	0.2		
State U	inregulated - Other						
114-26-1	PROPOXUR (BAYGON)	ND **	ug/L .	1.0	0.4		
2032-65-7	METHIOCARB	ND	ug/L	1.0	0.3		

If a compound is detected > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.

MCL (Maximum Contaminant Level) maximum permissible level of a contaminant in water established by EPA; Federal Action Levels are 0.015 mg/L for Copper, Sodium has a recommended limit of 20 mg/L. A blank MCL value indicates a level is not currently established.

Trigger Level: DOH Drinking Water Response level. Systems with compounds detacted in excess of this level are required to take additional samples. Contact your regional DOH office.

ND (Not Delected): indicates that the parameter was not delected above the State Reporting Limit (SRL).







QUALITY CONTROL REPORT SURROGATE REPORT

Reference Number: 08-17751

Lab No	Analyte	Result Qualifier	Units	Method	Limit
515 081222				1	
37230	2,4 - DCAA (SURR)	94	%	515.1	Acceptance Range is 70 - 130%
525 081222	_, , _ , _ , _ , _ , _ , _ ,				
37230	1,3-DIMETHYL-2-NITROBENZENE (Surr)	109	%	525.2	Acceptance Range is 70% to 130%
21200	PYRENE-D10 (Surr)	93	%		Acceptance Range is 70% to 130%
	PERYLENE-D12 (Surr)	109	%		Acceptance Range is 70% to 130%
	TRIPHENYLPHOSPHATE (Suit)	106	%		Acceptance Range is 70% to 130%
	· · · · · · · · · · · · · · · · · · ·				
515 081222					
37231	2,4 - DCAA (SURR)	94	%	515.1	Acceptance Range is 70 - 130%
525 081222	Z,4 - DCMA (SUNN)	34	7 0	0.011	recopedition (and in 1 a 1 a 1 a 1 a 1 a 1 a 1 a 1 a 1 a 1
	4.2 DIMETION O NITRODENZENE (C)	109	%	525.2	Acceptance Range is 70% to 130%
37231	1,3-DIMETHYL-2-NITROBENZENE (Surr)	93	<i>7</i> 6	020.2	Acceptance Range is 70% to 130%
·	PYRENE-D10 (Sum) PERYLENE-D12 (Sum)	102	%		Acceptance Range is 70% to 130%
	TRIPHENYLPHOSPHATE (Surr)	. 107	%		Acceptance Range is 70% to 130%
531 081217	MINICIPAL HOSPINIE (Sul)	, 101	7 0		
	POMC (CLIDD)	118	%	531.2	
37231	BDMC (SURR)	110	<i>7</i> e	001.L	
515_081222					1
37232	2,4 - DCAA (SURR)	88	%	515.1	Acceptance Range is 70 - 130%
525_081222			4.		A
37232	1,3-DIMETHYL-2-NITROBENZENE (Surr)	103	%	5 2 5.2	Acceptance Range is 70% to 130%
	PYRENE-D10 (Surt)	92	%		Acceptance Range is 70% to 130%
	PERYLENE-D12 (Sum)	109	%		Acceptance Range is 70% to 130%
	TRIPHENYLPHOSPHATE (Sum)	110	%		Acceptance Range is 70% to 130%
531_081217					
37232	BDMC (SURR)	110	%	531.2	
515 081222					
37233	2,4 - DCAA (SURR)	97	%	515.1	Acceptance Range is 70 - 130%
525 081222					
37 2 33	1,3-DIMETHYL-2-NITROBENZENE (Surr)	99	%	525.2	Acceptance Range is 70% to 130%
	PYRENE-D10 (Surr)	84	%		Acceptance Range is 70% to 130%
	PERYLENE-D12 (Surr)	102	%		Acceptance Range is 70% to 130%
	TRIPHENYLPHOSPHATE (Sum)	104	%		Acceptance Range is 70% to 130%
531 081217		14.			
37233	BDMC (SURR)	116	%	531.2	
	·				
515 081222					
37234	2,4 - DCAA (SURR)	91	%	515.1	Acceptance Range is 70 - 130%
525 081222	2,17 20.01 (00.01)				•
37234	1,3-DIMETHYL-2-NITROBENZENE (Surr)	102	%	525.2	Acceptance Range is 70% to 130%
01204	PYRENE-D10 (Sur)	92	%		Acceptance Range is 70% to 130%
	PERYLENE-D12 (Surr)	99	%		Acceptance Range is 70% to 130%
	TRIPHENYLPHOSPHATE (Surr)	109	%		Acceptance Range is 70% to 130%
531 090113	THE PERSON NAME OF THE PARTY				
37234	BDMC (SURR)	107	%	531.2	•
01201	DDIRE (CC/A/)				
E4E 004000					
515_081222	CA DOAA (CUDD)	95	%	515.1	Acceptance Range is 70 - 130%
37235	2,4 - DCAA (SURR)	30	78	313.1	Acceptance vande is to a look
525_081222	4.0 DATETING OF LITTLE CO	400	₽/	E95 9	Acceptance Range is 70% to 130%
37235	1,3-DIMETHYL-2-NITROBENZENE (Surr)	106	%	525.2	Acceptance Range is 70% to 130% Acceptance Range is 70% to 130%
	PYRENE-D10 (Surr)	88	% %		Acceptance Range is 70% to 130% Acceptance Range is 70% to 130%
	PERYLENE-D12 (Sur)	103	% %		Acceptance Range is 70% to 130%
504 000440	TRIPHENYLPHOSPHATE (Surr)	106	70		Undeltatiles utilide is un to m 1964)
531_090113	20140 4014001	447	ev.	531.2	
37235	BDMC (SURR)	117	%	031.2	

^{*}Notation:

A surrogate is a pure compound added to a sample in the laboratory just before processing so that the overall efficiency of a method can be determined.



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SAMPLE INDEPENDENT **QUALITY CONTROL REPORT**

Laboratory Fortified Blank

Reference Number: 08-17751

			True			%		QC	
Batch	Analyle	Result	Value	Units	Method	Recovery	Limits	QualifierType*	Comment
200.7-081223A	HARDNESS	71.3	69.5	mg/L	200.7	103	80-120	LFB	
515_081222	2,4 - D	1.61	2	ug/L	515. 1	81	70-130	LFB	
	2,4 - DCAA (SURR)	100		%	515.1		70-130		
	2,4 DB	7.25	8	ug/L	515.1	91	70-130		
	2,4,5 - TP (SILVEX)	0.88	1	ug/L	515.1	88	70-130		•
	2,4,5 T	0.79	1	ug/L	515.1	79	70-130		
	ACIFLUORFEN	0.72	1 .	ug/L	515.1	72	70-130		
	BENTAZON	1.7	2	ug/L	515.1	85	70-130		
	CHLORAMBEN	0.98	1	ug/L	515.1	98	70-130		
	DALAPON	15.3	13	ug/L	515.1	118	70-130		
	DICAMBA	0.86	1	ug/L	515.1	86	70-130		
	DICHLORPROP	2.63	3	ug/L	515.1	88	70-130		
	DINOSEB	1.69	2	ug/L	515. 1	85	70-130		
	PENTACHLOROPHENOL	0.92	1	սց/L	5 15 .1	92	70-130		
	PICLORAM	0.97	1	ug/L	515.1	97	70-130		•
	TOTAL (DCPA & Metabolites)	1.34	1	ug/L	515.1	134	70-130	HQ	
525 <u>0</u> 81222	1,3-DIMETHYL-2-NITROBENZENE (Surr)	100		%	525.2		70-130	LFB	
_	4.4-DDD	1	1	ug/L	525.2	100	70-130		
	4,4-DDD	1	1	ug/L	525.2	100	70-130		
	4,4-DDE	1 **	1	ug/L	525.2	100	70-130		
	4,4-DDE	1	1	ug/L	525.2	100	70-130		
	4,4-DDT	1.12	1	ug/L	525.2	112	70-130		
	4,4-DDT	1.12	1	ug/L	525.2	112	70-130		
	ACENAPHTHYLENE	1.11	1	ug/L	525.2	111	70-130		
	ALACHLOR	2.27	2	ug/L	525.2	114	70-130		
	ALDRIN	0.98	1	ug/L	525.2		70-130		
	ANTHRACENE	1.06	1	ug/L	525.2	106	70-130		
	ATRAZINE	2.35	2	ug/L	525.2	118	70-130		
	BENZ(A)ANTHRACENE	1.09	1	ug/L	525.2	109	70-130		
	BENZO(A)PYRENE	0.96	1	ug/L	525.2	96	70-130		
	BENZO(B)FLUORANTHENE	1.05	1	ug/L	525.2	105	70-130		
	BENZO(G,H,I)PERYLENE	0.82	1	ug/L	525.2	82	70-130		

[&]quot;Notation:

[%] Recovery = (Result of Analysis)/(True Value) * 100

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SAMPLE INDEPENDENT QUALITY CONTROL REPORT

Laboratory Fortified Blank `

Reference Number: 08-17751

			True			%		QC	
Batch	Analyte	Resuit	Value	Units	Method	Recovery	Limits	Qualifier Type*	Comment
525_081222	BENZO(K)FLUORANTHENE	1.04	1	ug/L	525.2	104	70-130	LFB	
	BENZYL BUTYL PHTHALATE	1.14	1	n ∂ √	525.2	114	70-130		
	BROMACIL	1.05	1	ug/L	525.2	105	70-130		
	BUTACHLOR	1.13	1	ug/L	52 5.2	113	70-130		
	CHLORDANE, TECHNICAL	1.04	1	n ∂ /Γ	525.2	104	70-13 0		
	CHRYSENE	1.09	1	ug/L	525.2	109	70-130		
	CYANAZINE	2.13	2	ug/L	525.2	107	70-130		
	DI(ETHYLHEXYL)-ADIPATE	1.09	1	ug/L	525.2	109	70-130		
	DI(ETHYLHEXYL)-PHTHALATE	1.4	1	ug/L	525.2	140	70-130	AC	
	DIAZINON	3.12	3	ug/L	525.2	104	70-130		
	DIAZINON	3.12	3	ug/L	525.2	104	70-130		
	DIBENZO(A,H)ANTHRACENE	0.92	1	ug/L	525.2	92	70-130		
•	DIELDRIN	1. 18	1	ug/L	525.2	118	70- 130		
	DIETHYL PHTHALATE	1.2	1	ug/L	525.2	120	70-130		
	DIMETHYL PHTHALATE	1.11	1	ug/L	525.2	111	70-130		
	DI-N-BUTYL PHTHALATE	1.15	1	ug/L	525.2	115	70-130		
	ENDRIN	1.06	1	ug/L	525.2	106	70-130		
	EPTC	1.14	1	ug/L	525.2	114	70-1 30		
	FLUORENE	1.16	1	ug/L	525.2	116	70-130		
	HEPTACHLOR	1.02	1	ug/L	525.2	102	70-130		
	HEPTACHLOR EPOXIDE	1.11	1	ug/L	525.2	111	70-130		
	HEXACHLOROBENZENE	1.12	1	ug/L	525.2	112	70-130		
	HEXACHLOROCYCLO-PENTADIENE	1.13	1	ug/L	525.2	113	70-130		
	INDENO(1,2,3-CD)PYRENE	0.91	1 .	ug/L	525.2		70-130		
	LINDANE (BHC - GAMMA)	1.01	1	ug/L	525.2	101	70-130		
	LINDANE (BHC - GAMMA)	1.01	1	ug/L	525.2	101	70-130		
	MALATHION	2.15	2	ug/L	525.2	108	70-130		
	MALATHION	2.15	2	ug/L	525,2		70-130		
	METHOXYCHLOR	1.09	1	ug/L	525.2		70-130		
	METOLACHLOR	1.15	1	ug/L	525.2		70-130		
	METRIBUZIN	0.82	1	ug/L	525.2		70-130		•
	PARATHION	2.39	2	ug/L	525.2		70-130		
	PARATHION-ETHYL	2.39	2	ug/L	525.2		70-130		
	PENTACHLOROPHENOL	4.99	4	ug/L	525.2		70-130		
	PERYLENE-D12 (Surr)	95	•	%	525.2	-	70-130		
	PHENANTHRENE	1.11	1	ug/L	525.2		70-130		
		••••	•	-Alir	JEU.E	• • •	10-100		

[&]quot;Notation:

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SAMPLE INDEPENDENT QUALITY CONTROL REPORT

Laboratory Fortified Blank

Reference Number: 08-17751

			True			%		QC	
Batch	Analyte	Result	Value	Units	Method	Recovery	Limits	QualifierType*	Comment
525_0B1222	PROPACHLOR	1.21	1	ug/L	525.2	121	70-130	LFB	
	PYRENE	1.11	1	ug/L	525.2	111	70-130		
	PYRENE-D10 (Surr)	90		%	525.2		70-130		
	SIMAZINE	1.04	1	ug/L	525.2	104	70-130		
	SIMAZINE	1.04	1	ug/L	525.2	.104	70-130		
	TERBACIL	1.07	1	ug/L	525.2	107	70-130		
	TRIFLURALIN	1.15	1	ug/L	525.2	115	70-130		
	TRIPHENYLPHOSPHATE (Surr)	105		%	525.2		70-13 0		
525X_081222	1-NAPHTHALENEACETAMIDE	2.1	2	ug/L	525.2	1 05	70-130	LFB	
	AZINPHOS-METHYL	1.1	1	ug/L	525.2	110	70-130		
	CHLORPYRIFOS	3.6	3	ug/L	525.2	120	70-130		
-	DICOFOL	3.5	3	ug/L	525.2	117	70-130		
	DIMETHOATE	0.7	1	ug/L	525.2	70	70-130		
	FENARIMOL	1.9	2	ug/L	525.2	95	70-130		
	HEXAZINONE	1	1	ug/L	525.2	100	70-130		
	HEXAZINONE (Velpar)	1	1	ug/L	525.2	100	70-130		
	METALAXYL	2.2	2	ug/L	525.2	110	70-130		
	METHIDATHINON	3.4	2	ug/L	525.2	170	85-115		-
	METHYL PARATHION	1.6	2	ug/L	525.2	80	70-130		
	MEVINPHOS	2.5	2	ug/L	525.2	125	70-130		
	MYCLOBUTANIL	2.4	2	ug/L	525.2	120	85-115		
	NAPROPAMIDE	1.08	1 .	ug/L	525.2	108	70-130		
	PHOSMET	3.3	2	ug/L	525.2	165	70-130	N1	
	PROPARGITE	3.5	2	ug/L	525.2	175	85-115	N1	
	TRIADIMEFON	2.45	2	ug/L	525.2	123	70-130		
	TRIFLUMIZOLE	2.3	2	ug/L	525.2		85-115		
525X_081222	1-NAPHTHALENEACETAMIDE	2.5	2	ug/L	525.2	125	70-130	LFB	
	AZINPHOS-METHYL	1.5	1	.ug/L	525.2	150	70-130	НО	
	CHLORPYRIFOS	3.7	3	ug/L	525.2	123	70-130		
	DICOFOL	3.5	3	ug/L	525.2	117	70-130		
	DIMETHOATE	8.0	1	ug/L	525.2	80	70-130		
	FENARIMOL	2.3	2	ug/L	525.2	115	70-130		
	HEXAZINONE	1.2	1	ug/L	525.2		70-130		
	HEXAZINONE (Velpar)	1.2	1	ug/L	525.2		70-130		

^{*}Notation:

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SAMPLE INDEPENDENT QUALITY CONTROL REPORT

Laboratory Fortified Blank

Reference Number: 08-17751

			True			%		QC	
Batch	Analyte	Result	Value	Units	Method	Recovery	Limits	Qualifier Type*	Comment
525X_081222	METALAXYL	2.3	2	ug/L	525.2	115	70-130	LFB	
	METHIDATHINON	3.8	2 .	ug/L	525.2	190	85-115	N1	
	METHYL PARATHION	1.6	2	ug/L	525.2	80	70 -130		
	MEVINPHOS	2.8	2	ug/L	525.2	140	70-130	HQ	
	MYCLOBUTANIL	2.7	2	ug/L	525.2	135	85-115	N1	
	NAPROPAMIDE	1.1	1	ug/L	525.2	110	70-130		
	PHOSMET	3.3	2	ug/L	525.2	165	70-130	N1	
	PROPARGITE	3.5	2	ug/L	525.2	175	85-115	N1	
	TRIADIMEFON	2.3	2	ug/L	525.2	115	70-130		
	TRIFLUMIZOLE	2.3	2	ug/L	525.2	115	85-115		
531_081217	3-HYDROXYCARBOFURAN	11.4	10	ug/L	531,2	114	70-130	LFB	
	ALDICARB	11.3	10	ug/L	531.2	113	70-130	G G	
	ALDICARB SULFONE	11,2	10	ug/L	531.2	112	70-130		
	ALDICARB SULFOXIDE	10.8	10	ug/L	531.2	108	70-130		
	BDMC (SURR)	101		~ <u>~</u> ~	531.2		70-130		
	CARBARYL	11.3	10	ug/L	531.2	113	70-130		•
	CARBOFURAN	10.4	10	ug/L	531.2	104	70-130		
	METHIOCARB	10.4	10	ug/L	531.2	104	70-130		
	METHOMYL	11	10	ug/L	531.2	110	70-130		
	OXYMAL.	10.6	10	ug/L	531.2	106	70-130		
	PROPOXUR (BAYGON)	11.1	10	ug/L	531.2	111	70-130		
531_081217	3-HYDROXYCARBOFURAN	17.4	20	ug/L	531.2	87	70-130	LFB	
	ALDICARB	18	20	ug/L	531.2	90	70-130		
	ALDICARB SULFONE	17.7	20	ug/L	531,2	89	70-130		
	ALDICARB SULFOXIDE	17.6	20	ug/L	531.2	88	70-130		
	BDMC (SURR)	97		%	531.2		70-130		
	CARBARYL	18.5	20	ug/L	531.2	93	70-130		
	CARBOFURAN	16.9	20	ug/L	531.2	85	70-130		
	METHIOCARB	19.5	20	u g/L	531.2	98	70-130		
	METHOMYL	18	20	ug/L	531.2	90	70-130		
	OXYMAL	17.3	20	ug/L	531.2	87	70-130		
	PROPOXUR (BAYGON)	18.1	20	ug/L	531.2	91	70-130		

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SAMPLE INDEPENDENT QUALITY CONTROL REPORT

Laboratory Fortified Blank

Reference Number: 08-17751

			True			%		QC	
Balch	Analyte	Result	Value	Units	Method	Recovery	Limits	Qualifier Type*	Comment
531_090113	BOMC (SURR)	113		%	531,2	·	70-130	LFB	
531_090113	3-HYDROXYCARBOFURAN	21.6	20	ug/L	531.2	108	70-130	LFB	
	ALDICARB	21.6	20	ug/L	531,2	108	70-130		
	ALDICARB SULFONE	21.2	20	ug/L	531.2	106	70-130		
	ALDICARB SULFOXIDE	21.3	20	ug/L	531.2	107	70-130		
	BDMC (SURR)	89		%	531.2		70-130		
	CARBARYL	21.2	20	ug/L	531.2	106	70-130		
	CARBOFURAN	19.6	20	ug/L	531.2	98	70-130		
•	METHIOCARB	17.3	20	ug/L	531.2	87	70-130		
	METHOMYL	21.7	20	ug/L	531.2	109	70-130		
	OXYMAL	20.6	20	ug/L	531.2	103	70-130		
	PROPOXUR (BAYGON)	21.1	20	nð/L	531.2	106	70-130		
COD_081229	CHEMICAL OXYGEN DEMAND	55	50	mg/L	SM5220 D	110	60-120	LFB	
OPHOS-081217	ORTHO-PHOSPHATE	1.01	1.00	mg/L	SM4500-P F	101	70-130	LFB	
tds_081219	TOTAL DISSOLVED SOLIDS	514	500	mg/L	SM2540 C	103	80-120	LFB	
tds_081219	TOTAL DISSOLVED SOLIDS	500	500	mg/L	SM2540 C	100	80 -120	LFB	
tds_081219	TOTAL DISSOLVED SOLIDS	494	500	mg/L	SM2540 C	99	80-120	LFB	
lds_081219	TOTAL DISSOLVED SOLIDS	472	500	mg/L	SM2540 C	94	80-120	LFB	

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SAMPLE INDEPENDENT **QUALITY CONTROL REPORT**

Low Level Laboratory Fortified Blank

Reference Number: 08-17751

			True			%		QC	
Batch	Analyle	Result	Value	Units	Method	Recovery	Limits	Qualifier Type*	Comment
531_061217	3-HYDROXYCARBOFURAN	0.73	1	ug/L	531.2	73	50-150	LFBD	
	ALDICARB	0.99	1	ug/L	531.2	99	50- 150		
	ALDICARB SULFONE	0.95	1	ug/L	531.2	95	50- 150		
	ALDICARB SULFOXIDE	0.86	1	ug/L	531.2	86	50-150		
	BDMC (SURR)	91		%	531.2		50-150		
	CARBARYL	1.1	1	ug/L	531.2	110	50-150		
	CARBOFURAN	0.89	1	ug/L	531.2	89	50-150		
	METHIOCARB	1.1	1	ug/L	531.2	110	50-150		
	METHOMYL	0.81	1	⊔g/L	531.2	B1	50-150		
	OXYMAL	0.96	1	ug/L	531.2	96	50-150		
	PROPOXUR (BAYGON)	0.99	1	ug/L	531.2	99	50-150		
531_090113	3-HYDROXYCARBOFURAN	1.1	1	ug/L	531.2	110	50-150	LFBD	
	ALDICARB	1.15	1	ug/L	531.2	115	50-150		
	ALDICARB SULFONE	0.95	1	ug/L	531.2	95	50-150		
	ALDICARB SULFOXIDE	0.97	1	ug/L	531.2	97	50-150		
	BDMC (SURR)	113		%	531.2		50-150		
	CARBARYL	1.16	1	ug/L	531.2	116	50-150		
	CARBOFURAN	0.96	1	ug/L	531.2	96	50-150		
	METHIOCARB	1.13	1	ug/L	531.2	113	50-150		
	METHOMYL	1	1	ug/L	531.2	100	50-150		
	OXYMAL	0.93	1	ug/L	531.2	93	50-150		
	PROPOXUR (BAYGON)	0.95	1 .	ug/L	531.2	95	50-150		

[&]quot;Notation:

[%] Recovery = (Result of Analysis)/(True Value) * 100

NA = Indicates % Recovery could not be calcufated.

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SAMPLE INDEPENDENT QUALITY CONTROL REPORT

Laboratory Reagent Blank

Reference Number: 08-17751

			True			%	QC	
Batch	Analyte	Result	Value	Units	Method	Recovery Limits	Qualifier Type*	Comment
200.7-081223A	HARDNESS	ND		mg/L	200.7	10.000		
COD_081229	CHEMICAL OXYGEN DEMAND	NĐ		mg/L	SM5220 D	4.0000	0 LRB	
D081230A	BROMATE	ND		mg/L	300.1	0.0050	O LR6	
I081217A	CHLORIDE NITRATE-N	ND ND		mg/L mg/L	300.0 300.0	0.1000 0.1000		
OPHOS-081217	ORTHO-PHOSPHATE	ND		mg/L	SM4500-P F	0.1000	D LRB	·
TURB_081217	TURBIDITY	NĎ		NTU	180.1	0.0200) LRB	

[&]quot;Notation:

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SAMPLE INDEPENDENT **QUALITY CONTROL REPORT**

Method Blank

Reference Number: 08-17751

			True			%		QC	
Batch	Analyte	Result	Value	Units	Method	Recovery	Limits	Qualifier Type*	Commen
200.7-081223A	HARDNESS	ND		mg/L	200.7		0.82000	MB	
515_081222	2,4 - D	ND		ug/L	515.1		0.05000	MB	
	2,4 - DCAA (SURR)	95		%	515.1	•			
	2,4 DB	ND		ug/L	515.1		0.25000		
	2,4,5 - TP (SILVEX)	ND		ug/L	515.1		0.10000		
	2,4,5 T	ND		ug/L	515.1		0.10000		
	ACIFLUORFEN	ND		ug/L	515.1		0.50000		
	BENTAZON	ND		ug/L	515.1		0.12000		
	CHLORAMBEN	ND		ug/L	515.1		0.20000		
	DALAPON	ND		ug/L	515.1		0.50000		
	DCPA (ACID METABOLITES)	ND		ug/L	515.1		0.10000		
	DICAMBA	ND		ug/L	515.1		0.05000		
	DICHLORPROP	ND		ug/L	515.1		0.12000		
	DINOSEB	ND		ug/L	515.1		0.10000		
	PENTACHLOROPHENOL	ND		ug/L	515.1		0.02000		
	PICLORAM	ND		ug/L	515.1		0.05000		
	TOTAL (DCPA & Metabolites)	ND		ug/L	515.1	-	0.02000		
25_081222	1,3-DIMETHYL-2-NITROBENZENE (Surr)	106		0,	ror n				
25_001222	4,4-DDD			% !!	525.2			MB	
•	4,4-DDD	ND ND		ug/L	525.2		0.05000		
	4,4-DDE		·	ug/L. -	525.2		0.05000		
	4,4-DDE	ND		ug/L -	525.2		0.05000		
	4,4-DDT	ND		ug/L	525.2		0.05000		
	4,4-DDT	ND		ug/L	525.2		0.05000		
	ACENAPHTHENE	ND		ug/L	525.2		0.05000		
	ALACHLOR	ND ND		ug/L	525.2		0.05000		
	ALDRIN	ND		ug/L	525.2		0.02000		
	ANTHRACENE	ND		ug/L	525.2		0.05000	-	
		ND		ug/L	525.2		0.05000		
	ATRAZINE	ND		ug/L	525.2		0.02000		
	BENZ(A)ANTHRACENE	ND		ug/L	525.2		0.02000		
	BENZO(A)PYRENE	ND		ug/L	525.2		0.02000		
	BENZO(B)FLUORANTHENE	ND		ug/L	525.2		0.05000		

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SAMPLE INDEPENDENT **QUALITY CONTROL REPORT**

Method Blank

Reference Number: 08-17751

			True			%	QC .	
Batch	Analyte	Result	Value	Units	Method	Recovery Limits	Qualifier Type*	Comment
525_081222	BENZO(G,H,I)PERYLENE	ND		ug/L	525.2	0.0500		
	BENZO(K)FLUORANTHENE	ND		ng/L	525.2	0.0500	0	
	BENZYL BUTYL PHTHALATE	ND		ug/L	525.2	0.6000	0	
	BROMACIL	ND		ug/L	525.2	0.0500	0	
	BUTACHLOR	ND		ug/L	525.2	0.1000	D	
	CHLORDANE, TECHNICAL	ND		ug/L	525.2	0.0200	0	
	CHRYSENE	ND		ug/L	525.2	0.0500	0	
	CYANAZINE	ND		ug/L	525.2	0.0500	0	
	DI(ETHYLHEXYL)-ADIPATE	NĐ		ug/L	525.2	0.0200	0	
	DI(ETHYLHEXYL)-PHTHALATE	ND		ug/L	525.2	0.6000	0	
	DIAZINON	ND		ug/L.	525.2	0.0500	0	
	DIAZINON	ND		ug/L	525.2	0.0500	D	
	DIBENZO(A,H)ANTHRACENE	ND		ug/L	525.2	0.0500	0	
	DIELDRIN	ND		ug/L	525.2	0.0500	0	
	DIETHYL PHTHALATE	ND		ug/L	525.2	0.6000	0	
	DIMETHYL PHTHALATE	ND		ug/L	525.2	0.6000		
	DI-N-BUTYL PHTHALATE	ND		ug/L	525.2	0.6000	0	
	ENDRIN	ND		ug/L	525.2	0.0200		
	EPTC	ND		ug/L	525.2	0.0700	0	
•	FLUORANTHENE	ND		ug/L	525.2	0.0500	0	
	FLUORENE	ND		ug/L	525.2	0.0500)	
	HEPTACHLOR	ND		ug/L	525.2	0.0200)	
	HEPTACHLOR EPOXIDE	ND		ug/L	525.2	0.0200)	
	HEXACHLOROBENZENE	ND		⊔g/L	525.2	0.0200)	
	HEXACHLOROCYCLO-PENTADIENE	ND		ug/L	525.2	0.0200)	
	INDENO(1,2,3-CD)PYRENE	ND		ug/L	525.2	0.0500)	
	LINDANE (BHC - GAMMA)	ND		ug/L	525.2	0.0200		
	LINDANE (BHC - GAMMA)	ND		ug/L	525.2	0.0200		
	MALATHION	. ND		ug/L	525.2	0.0500		
	MALATHION	ND		ug/L	525.2	0.0500		
	METHOXYCHLOR	ND		ug/L.	525.2	0.0200		
	METOLACHLOR	ND		ug/L	525.2	0.2500		
	METRIBUZIN	ND		ug/L	525.2	0.0500		
	NAPTHALENE	ND		ug/L	525.2	0.0200		
	PARATHION	ND		ug/L	525.2	0.0500		
	PARATHION-ETHYL	ND		ug/L	525.2	0.0500		

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SAMPLE INDEPENDENT **QUALITY CONTROL REPORT**

Method Blank

Reference Number: 08-17751

			True			%	QC	
Batch	Analyte	Result	Value	Units	Method	Recovery Limits	Qualifier Type*	Comment
525_081222	PENTACHLOROPHENOL	ND		ug/L	525.2	0.0400		
	PERYLENE-D12 (Surr)	100		%	525.2			
	PHENANTHRENE	ND		ug√L	525.2	0.0500	0	
	PROPACHLOR	ND		⊔ g /L	525.2	0.0500	0	
	PYRENE	ND		ug/L	525.2	0.0500	D	
	PYRENE-D10 (Surr)	94		% .	525.2			
	SIMAZINE	ND		ug/L	525.2	0.0200	0	
	SIMAZINE	ND		ug/L	525.2	0.0200	0	
	TERBACIL	ND		ug/L	525.2	0.0500)	
	TRIFLURALIN	ND.		ug/L	525.2	0.0500)	
	TRIPHENYLPHOSPHATE (Suri)	107		%	525.2			
525X_081222	1-NAPHTHALENEACETAMIDE	ND		ug/L	525.2	0.10000) MB	
	AZINPHOS-METHYL	ND		ug/L	525.2	0.0000		
	CHLORPYRIFOS	ND		ug/L	525.2	0.0000		
	DICOFOL	ND		ug/L	525.2	0.0000		
	DIMETHOATE	ND		ug/L	525.2	0.0000		
	FENARIMOL	ND		ug/L	525.2	0.0000		
	HEXAZINONE	ND		ug/L	525.2	0.00000		·
	HEXAZINONE (Velpar)	ND		ug/L	525.2	0.02000		
	METALAXYL	ND		ug/L	525.2	0.10000		
	METHIDATHINON	ND		ug/L	525.2	0.50000		
	METHYL PARATHION	ND		ug/L	525.2	0.00000		
	MEVINPHOS	ND		ug/L	525.2	0.0000		
	MYCLOBUTANIL	ND		ug/L	525.2	0.50000		
	NAPROPAMIDE	ND		ug/L	525.2	0.00000		
	PHOSMET	ND		ug/L	525.2	0.10000		
	PROPARGITE	ND		ug/L	525.2	0.00000		
	TRIADIMEFON	ND		ug/L	525.2	0.00000		
	TRIFLUMIZOLE	ND		ug/L	525.2	1.00000		
31_081217	3-HYDROXYCARBOFURAN	ND		ug/L	531.2	0.50000	MB	
	ALDICARB	ND		ug/L	531.2	0.25000		
	ALDICARB SULFONE	ND		ug/L	531.2	0.40000		

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SAMPLE INDEPENDENT **QUALITY CONTROL REPORT**

Method Blank

Reference Number: 08-17751

			True		•	%		QC	
Batch	Analyte	Result	Value	Units	Method	Recovery	Limits	Qualifier Type*	Comment
531_081217	ALDICARB SULFOXIDE	ND		ug/L	531.2		0.25000	MB	
	BDMC (SURR)	96		%	531.2		0.00000		
	CARBARYL	ND		ug/L	531.2		0.50000		
	CARBOFURAN	ND		ug/L	531.2		0.45000		
	METHIOCARB	ND		ug/L	531.2		1.00000		
	METHOMYL	ND		ug/L	531.2		0.25000	r -	
	OXYMAL	ND		ug/L	531.2		1.00000		
	PROPOXUR (BAYGON)	ND		u g/L	531.2		0.25000		
531_090113	3-HYDROXYCARBOFURAN	ND		ug/L	531.2		0.50000	MB	
•	ALDICARB	ND		ug/L	531.2		0.25000		
	ALDICARB SULFONE	ND		ug/L	531.2		0.40000		
	ALDICARB SULFOXIDE	ND		⊔g/L	531.2		0.25000		
	BDMC (SURR)	121		%	531.2		0.00000		
	CARBARYL	ND		ug/L	531.2		0.50000		
	CARBOFURAN	ND		ug/L	581.2		0.45000		
	METHICCARB	ND		ug/L	531.2		1.00000		
	METHOMYL	ND		ug/L	531.2		0.25000		
	OXYMAL	ND		ug/L	531.2		1.00000		
	PROPOXUR (BAYGON)	ND		ug/L	531.2		0.25000		
ec_081223	ELECTRICAL CONDUCTIVITY	NÖ		u\$/cm	SM2510 B		2.50000	MB	
ec_081223	ELECTRICAL CONDUCTIVITY	ND		uS/cm	SM2510 B		2.50000	МВ	
ec_081223	ELECTRICAL CONDUCTIVITY	ND		uS/cm	SM2510 B		2.50000	MB	
ec_081223	ELECTRICAL CONDUCTIVITY	ND	-	uS/cm	SM2510 B		2.50000	MB	
							-		
OPHOS-081217	ORTHO-PHOSPHATE	ND		mg/L	SM4500-P F		0.10000	MB	
tds_081219	TOTAL DISSOLVED SOLIDS	ND		mg/L	SM2540 C		2.50000	MB	
tds_081219	TOTAL DISSOLVED SOLIDS	ND		_	CLIDE IO O	-			
20_201210	OTHE SINGULACH SAFINS	ND		mg/L	SM2540 C		2.50000	MB	

[&]quot;Notation:

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SAMPLE INDEPENDENT **QUALITY CONTROL REPORT**

Method Blank

Reference Number: 08-17751

Report Date: 01/28/09

-			True			%	QC	
Batch	Analyte	Result	Value	Units	Method	Recovery Limits	Qualifier Type*	Comment
tds_081219	TOTAL DISSOLVED SOLIDS	ND	······································	mg/L	SM2540 C	2.50000	MB	· · · · · · · · · · · · · · · · · · ·
tds_081219	TOTAL DISSOLVED SOLIDS	ND		mg/L	SM2540 C	2.50000	MB	

*Notation:

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SAMPLE INDEPENDENT QUALITY CONTROL REPORT

Quality Control Sample

Reference Number: 08-17751

			True			%		QC	
Balch	Analyte	Result	Value	Units	Method	Recovery	Limits	Qualifier Type*	Comment
200.7-081223A	HARDNESS	131	132.3	mg/L	200.7	99	80-120	QCS	······································
,		•							
531_061217	3-HYDROXYCARBOFURAN	41.2	40	ug/L	531.2	103	70-130	QCS	
	ALDICARB	37.5	37.3	ug/L	531.2	101	70-130		
	ALDICARB SULFONE	41.2	44.9	ug/L	531.2	92	70-130		•
	ALDICARB SULFOXIDE	43.2	40.2	ug/L	531.2	107	70-130		
	BDMC (SURR)	96		%	531,2		70-130		
	CARBARYL	47	46	ug/L	531.2	102	70-130		
	CARBOFURAN	60.6	60.9	ug/L	531.2	100	70-130		
	METHIOCARB	119.1	121	ug/L	531.2	98	70-130		
	METHOMYL	61.3	61.4	ug/L	531.2	100	70-130		
	OXYMAL	52.7	59.9	ug/L	531.2	88	70-130		
	PROPOXUR (BAYGON)	100	96.7	ug/L	531.2	103	70-130		
COD_081229	CHEMICAL OXYGEN DEMAND	89	92	mg/L	SM5220 D	97	80-120	QCS	
D061230A	BROMATE	0.0187	0.0182	mg/L	300.1	103	75-125	QCS	
ec_081223	ELECTRICAL CONDUCTIVITY	158	150.5	uS/cm	SM2510 B	105	80-120	QCS	
ec_081223	ELECTRICAL CONDUCTIVITY	157	150.5	uS/cm	SM2510 B	104	80-120	QCS	
ec_081223	ELECTRICAL CONDUCTIVITY	159	150.5	uS/cm	SM2510 B	106	80-120	QCS	
ec_081223	ELECTRICAL CONDUCTIVITY	157	150.5	uS/cm	SM2510 B	104	80-120	QCS	
108121 7 A	CHLORIDE	28.6	30 .0	mg/L	300.0	95	80-120	QCS	
	NITRATE-N	2.53	2.50	mg/L	300.0	101	80-120		
OPHOS-081217	ORTHO-PHOSPHATE	0.47	0.49	mg/L	SM4500-P F	96	70-130	QCS	

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SAMPLE INDEPENDENT QUALITY CONTROL REPORT

Quality Control Sample

Reference Number: 08-17751

Report Date: 01/28/09

			True			%	QC	
Batch	Analyte	Result	Value	Units	Method	Recovery Limits	Qualifier Type*	Comment
TURB_081217	TURBIDITY	0.98	1.00	NTU	180.1	98 70-13	0 QCS	

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Reference Number: 08-17751

QUALITY CONTROL REPORT

Dunitate and Matrix Shike/Matrix Shike Dunitate Depart

	Didna	ate ar	ıd Matrix Spik	Duplicate and Matrix Spike/Matrix Spike Duplicate Report	ָ בס			Report Date: 1/28/2009
Duplicate								
			Duplicate				30	
	Sample Analyte	Result	Result	Units	%RPD	Limits	QueiMer	Comments
200.7-081223A								
	37079 HARDNESS	119	117	mg CaCO3/L	1.7	54-0	ф	
	37234 HARDNESS	74.0	73.5	mg CaCO3/L		0.45	PUP	
515_081222			'te					
	37091 DCPA (ACID METABOLITES)	1.2	O.BMI.	ng/L	40.0	0-20	dua	
	37091 2,4 - DCAA (SURR)	95	72	*	27.5	<u>}</u>	đực	
525_081222								
- I	36873 1,3-DIMETHYL-2-NITROBENZENE (Sun	100	102	%	2.0	0-45	qua	
	38873 PYRENE-D10 (Sum)	83	88	*	0.0	0-45	DUP	
	36873 PERYLENE-D12 (Sum)	5	110	*	6.6	0.45	ana	
	38873 TRIPHENYLPHOSPHATE (Sur)	107	105	*	1.9	0-45	920	
	37230 BROMACIL	0.09	60'0	ng/L	0.0	0.45	DUP	
COD_081229								
D081230A								
EC_081223								
ı	37233 ELECTRICAL CONDUCTIVITY	171	167	us/cm	2.4	0.45	OUP	
	37385 ELECTRICAL CONDUCTIVITY	181	181	uS/cm	0.0	£	DUP	
	37476 ELECTRICAL CONDUCTIVITY	16.6	16.7	nS/cm	9.0	245	DUP	
1081217A				-				
	37252 CHLORIDE	8	8	mg/L	0,0	54.0	ana	
	37278 NITRATE-N	77.0	0.8	mg/L	3.8	0-45	DO	
OPHOS-081217								
	37235 ORTHO-PHOSPHATE	0.16	0.16	mg/L	0.0	0-20	PUP	
PH_081217						-		
	37235 HYDROGEN ION (pH)	6.58	6.54	pH Cnits	9.0	949	ana	
TDS_081219		ģ	3	£	,	Ų	Š	
	37056 (OIAL DISSOLVED SOLIDS	gg	115	Tright.	P.	£	Š	

%RPD = Relative Percent Difference

NA = indicates %RPD could not be calculated

Matrix Spike (MSVMatrix Spike Duplicate (MSD) analyses are used to determine the accuracy (MSD) and precision (MSD) of a analytical method in a given sample matrix. Therefore, the usefulness of this report is limited to samples of similar matrices analyzed in the same analytical batch.

. 3



Duplicate

Comments 출출 5 8 Cimils 3 9 %RPD 2 캶 Шg/L щď Z Duplicate Result 267 158 5.84 Result 5.78 168 37331 TOTAL DISSOLVED SOLIDS 37177 TOTAL DISSOLVED SOLIDS 37215 TURBIDITY Sample Analyte TURB_081217 Baich

Page 2 of 7 Reference Number: 08-17751 Report Date: 1/28/2009

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Matrix Spike (MS)/Matrix Spike Duplicate (MSD) analyses are used to determine the accuracy (MS) and practation (MSD) of a analytical method in a given sample matrix. Therefore, the usefulness of this report is limited to samples of similar matrices analyzed in the same analytical batch.



Batch

Report Date: 1/28/2009 Reference Number: 08-17751 Page 3 of 7

Comments ΞΞ Ę Σ F F F Ę Σ ΣĽ Ξ E ¥ Σ, Σ Œ, Ξ Ę Ξ Ę Ę Œ, Ę Ξ 7 Ξ Z Ę Ę Qualifier 뎣 90 Ŗ 9-0 8 8 F 9 8 8 윷 윩 윩 윩 9 윷 8 9 異 SRPO ٤ ŧ. ≶ \$ 85-135 85-135 70-130 65-135 55-135 85-135 85-135 35-135 85-135 35-135 35-135 35-135 80-120 85-135 65-135 65-135 55-135 70-130 65-135 65-135 Limits 65-135 65-135 85-135 85-135 65-135 65-135 65-135 65-135 65-135 Percent Recovery WSD ₹ ş ₹ ¥ ş ş ž ž ¥ ¥ ş ₹ ₹ ¥ 88 mg CaCO3/L 96 mg CaCO3/L 98 5 Ş 5 Ξ mg/L ПgЙ Ę 퍨 뎚 Splke 69.5 69.5 Duplicate Result Spilke **高** 左 Result Spike 0.71 0.97 0.95 3.03 5 첮 .65 9 0.81 8 4 Result 119 99 99 999 9 9 9 9993 身 TOTAL (DCPA & Metabolites) TOTAL (DCPA & Metabolites) PENTACHLOROPHENOL PENTACHLOROPHENOL 2,4 - DCAA (SURR) 2,4,5 - TP (SILVEX) 2,4,5 - TP (SILVEX) 2,4 - DCAA (SURR) DICHLORPROP DICHLORPROP CHLORAMBEN CHLORAMBEN ACIFLUORFEN ACIFLUORFEN 37079 HARDNESS 37234 HARDNESS BENTAZON BENTAZON PICLORAM PICLORAM DALAPON DALAPON DINOSEB DINOSEB DICAMBA DICAMBA 37231 ENDRIN 2,4,5 T 2,4 DB 2,4 DB 2,4,5 T 2.4.D Sample Analyte 37077 37077 37078 37079 37079 37079 37079 37079 37078 37079 **9707**B 3707B 37079 37079 37079 37079 37077 7,075 7707E 37077 37077 37075 37077 37077 37077 77078 37077 3707E Matrix Spike 200.7-081223A 525_081222 515_081222

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Report Date: 1/28/2009

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Matrix Spike

Batch

	ı																																		
	ı																																		
20 6	Qualifier	LFM	E I			MEI.	LFM	LFM	LFIM	LFM	LFM	LFIM	M.	LFM	LFIN	LFM	LFIX	LFM	LFM	FE	LFM	LFM	Ma	LFM	LFM.	LFM	LFM	LFM	LFM	LFM	LFM	FM	LFM	E.	LFM
i i	Lenits	2	6	2 2	8	0-80	0-80	0-60	09-0	0-20	9-60	09-0	09-0	0-20	09-0	0-60	9	0-60	0-60	09-0	0-20	09-0	9	9 9	9-60	90	0-60	09-0	9	0-60	9	9-90	9-0	09-0	9 0
9	%RPD	ΑN	¥ :	¥ ž	€	¥	¥	Ą	¥	¥	¥	Ϋ́	Ą	Ą	NA	NA	¥	¥	Ϋ́	¥.	ž	Ä	¥	¥	¥	Ϋ́	¥	ΑN	¥	ž	Ą	¥2	¥	Ϋ́	ž
# # #	Limits	70-130	70-130	70-130	70-130	70-130	70-130	70-130	70-130	70-130	70-130	70-130	70-130	70-130	70-130	70-130	70-130	70-130	70-130	70-130	70-130	70-130	70-130	70-130	70-130	70-130	70-130	70-130	70-130	70-130	70-130	70-130	70-130	70-130	70-130
Percent Recovery	MSD	¥	¥:	ě š	. ₹	¥	¥.	¥.	¥	NA A	¥	¥.	X X	¥	NA A	A A	¥.	NA N	¥	Ą	¥	Y	¥.	¥	¥	¥	۷ ۲	Ą V	¥	Ä	ΝA	Ϋ́	ΑN	Ą.	ΝA
Percent	MS	108	£	107	87	106	82	82	103	109	104	66	118	92	104	109	107	Ξ	8	128	104	106	102	114	118	114	114	5	95	108	9	120	103	112	104
a strait	Units	γđη	ug/L	1 5	, d	nov.	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ugvL	ug/L	ηđγ	1 .6 0/L	n b /L	- ng/L	ug/L	MD/L	ug/L	ng/L	ug/L	Ng.	rg/L	ug/L	ug/L	ug/L	ng/L	ug/L	ng∕l.	ugy	UgV.	ng/L	7,00
	Conc	-	- (ο ι ε		_	<u>-</u>	-	-	-	-	-	-	4		-	-	_	,-	-	10	-	-	60	-	-	60	-	-	_	_	8	8	8	_
Duplicate Spike	Result																															٠			
Spike	Result	1.08	1.15	2.13	0.87	1.08	0.82	0.85	1.03	1.09	1.04	0.99	1.18	3.67	1.04	1.09	1.07	1,11	0.81	1.28	5.2	1.06	1.02	3.42	1.18	1.14	3.42	-	0.92	1,08	-	2,4	2.08	2.24	8
Ę.	Result	₽	2	2 2	2 2	9	皇	2	2	Q	2	욮	2	2	2	2	2	2	2	2	9	₽	9	오	2	ð	오	오	2	2	ð	ջ	Z	2	2
Amenta de A				ALACHLOR	BENZO(A)PYRENE	CHLORDANE, TECHNICAL	DI(ETHYLHEXYL)-ADIPATE	DI(ETHYLHEXYLPHTHALATE	HEPTACHLOR	HEPTACHLOR EPOXIDE	HEXACHLOROBENZENE	HEXACHLOROCYCLO-PENTADIENE	SIMAZINE	PENTACHLOROPHENOL	ALDRIN	BUTACHLOR	DIELDRIN	METOLACHLOR	METRIBUZIN	PROPACHLOR	BISPHENOL-A	BROMACIL	TERBACIL	DIAZINON	SIMAZINE	EPTC	DIAZINON	4,4-000	4,4-DDE	LINDANE (BHC - GAMMA)	4,4-DDT	CYANAZINE	MALATHION	PARATHION	
Spike	Sample	37231	37231	37231	37231	37231	37231	37231	37231	37231	37231	37231	37231	37231	37231	37231	37231	37231	37231	37231	37231	37231	37231	37231	37231	37231	37231	37231	37231	37231	37231	37231	37231	37231	42024

%RPD = Relative Percent Difference

NA = Indicates %RPD could not be calculated

Martix Spike (MS)Martix Spike Duplicate (MSD) analyses are used to determine the accuracy (MS) and precision (MSD) of a analytical method in a given sample matrix. Therefore, the usefulness of this report is limited to samples of similar matrices analyzed in the same analytical batch.



Batch

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Report Date: 1/28/2009 Reference Number: 08-17751

pike				Duplicate				-							
			Spike	Spike	Spike		Percent	Percent Recovery				8			
Sample Analyte		Result	Result	Result	Солс	Units	MS.	MSD	Limits	%RPD	Limits	Qualifier	o	Соттепта	
37231	4,4-DDE	S	0.92		-	ug/L	92	NA A	70-130	ΑN	星		LFM		
37231	4,4-DDT	₽	Ψ.		_	ug/L	100	¥.	70-130	¥	0-60	-	Σ		
37231	MALATHION	욛	2.06		8	ng/L	103	¥	70-130	¥	09-0	_	Σ		
37231	PARATHION-ETHYL	Ş	2.24		CI	Jigu	112	¥.	70-130	¥	0-60	_	¥.		
37231	FLUORENE	2	1.09		-	ug/L	109	ΑĀ	70-130	¥	09-0	_	FM		
37231	ACENAPHTHYLENE	잎	1.07		.	ug/L.	107	¥.	70-130	≨	09-0	-	LFM		
37231	ANTHRACENE	모	1.05		-	ug/L	105	¥	70-130	¥	09-0	_	LFIN		
37231	BENZ(A)ANTHRACENE	9	1.04		-	ug/L	104	¥	70-130	¥	0-60	_	LFIN		
37231	BENZO(B)FLUORANTHENE	2	0.89		-	ug/L	68	¥	70-130	¥	09-0	_	Ę		
37231	BENZO(G,H,I)PERYLENE	2	0.54		_		2	Ā	70-130	ž	9-60	뿔	LFM		
37231	BENZO(K)FLUORANTHENE	오	0.89		_	ug/L	89	NA	70-130	Ā	0-60	_	LFM		
37231	CHRYSENE		1.05		_	ug/L	105	NA A	70-130	¥	0-80	_	LFM		
37231	DIBENZO(A,H)ANTHRACENE	 2	0.83		_	ug/L	63	N A	70-130	¥	090	ME	LFM		
37231	INDENO(1,2,3-CD)PYRENE	2	0.54		_	ug/L	3	¥	70-130	¥	0.60	₩	LFM		
37231	PHENANTHRENE	S	1.09		_	ug/L	109	¥.	70-130	¥	09-0		LF.M		
37231	PYRENE	2	1.06		_	ug/L	106	NA A	70-130	¥	09-0	_	LFIN		
37231	BENZYL BUTYL PHTHALATE	Q	1.05		-	ug/L	105	Ą2	70-130	¥	0-60	_	LFM		
37231	DI-N-BUTYL PHTHALATE	2	1.15		γ	ug/L	115	NA A	70-130	¥	09-0	_	LFM		
37231	DIETHYL PHTHALATE	ç	1.45		_	₩.	145	NA NA	70-130	¥	09-0	82	LFM		
37231	DIMETHYL PHTHALATE	2	1.08		_	ug/L	108	Ϋ́ Y	70-130	¥	0-60	_	LFM		
37231	1,3-DIMETHYL-2-NITROBENZËNË (Sum	109	104			*		٧X	70-130	¥	0-60	_	LF.N		
37231	PYRENE-D10 (Sur)	93	85			*		¥	70-130	¥	09-0	_	LFM		
37231	PERYLENE-D12 (Sun)	102	86			%		AN A	70-130	¥	09-0	_	LFM		
37231	TRIPHENYLPHOSPHATE (Sur.)	107	108			%		NA	70-130	Ą	0-60	_	E.F.		
37231	HEXAZINONE	Ş	1.2		-	ug/L	2	Υ Y	70-130	¥ Z	- 20		E E		
37231	HEXAZINONE (Vetpar)	D Z	1		-	ug/L	120	Ā	70-130	ž	0-60	_	LFM		
36873	OXYMAL	Q	10.5	7	5	ug/L	105	110	70-130	4.7	6 50	_	LFM		
38873	CARBOFURAN	2	10.4	10.5	5	ug/L	104	105	70-130	1.0	0-20	_	EN		
36873	ALDICARB SULFOXIDE	õ	10.6	11.6	무	ug/L	106	116	70-130	9.0	0-20		LFM		
36873	ALDICARB SULFONE	9	10.8	11.7	₽	√gn	108	117	70-130	8.0	0-90		LFM		
36873	METHOMYI.	2	10.8	11.4	5	ug/L	106	114	70-130	7.3	0-20		LFK		
36873	3-HYDROXYCARBOFURAN	S	10.9	;-	무	₩.	109	110	70-130	6.0	0-50		LFM		
36873	ALDICARB	Q	1.1	10.9	무	T/Gn	11	109	70-130	1.6	0-50	_	Ē		
36873	CARBARYL	₽.	F	11.2	6	ng/L	110	112	70-130	1.8	0-20		LF.M		

%RPD = Relative Percent Difference

525X_081222

531_081217

NA = Indicates %RPD could not be calculated
Matrix Spike (MSD) analyses are used to determine the accuracy (MS) and precision (MSD) of a analytical matrices analytical matrix. Therefore, the usefulness of this report is limited to samples of similar matrices analyzed in the same analytical batch.



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Report Date: 1/28/2009 Reference Number: 08-17751

Matrix Spike	sike				Duplicate									
				Spike	Spike	Spike		Percent Recovery	ecovery				ည္က	
Batch	Sample	Sample Analyte	Result	Result	Result	Conc	Units	MS	MSD		%RPD	Limits	Qualifier	Comments
	36873	PROPOXUR (BAYGON)	2	11.1	11.1	2	ngv	111	111	70-130	0.0	0-50	LFM	The state of the s
	36873	METHIOCARB	2	1.1	=	우	ug/L	11	1	70-130	0.0	0-20	LFR	
	36873	BOMC (SURR)	5	103	8		*		¥	70-130	¥2	0-50	LFE	
	37232	OXYMAL	ð	11.5	::	10	ug/L.	115	11	70-130	3.5	0-50	LFM	
	37232	CARBOFURAN	ç	10.4	11.5	9	ng/L	104	115	70-130	10.0	0-20	LFM	
	37232	ALDICARB SULFOXIDE	g	12.1	11.9	6	ug/L	121	119	70-130	1.7	0-50	Ę	
	37232	ALDICARB SULFONE	오	12.2	11.9	5	ug/L	122	119	70-130	2.5	0-50	FR	
	37232	METHOMYL	2	12.2	11.7	무	ng/L	122	117	70-130	4.2	0-50	LF.	
	37232	3-HYDROXYCARBOFURAN	2	11.8	11.7	우	ug/L	118	117	70-130	6.0	0-50	LFM	
	37232	ALDICARB	2	12	11.7	5	ug/L	120	117	70-130	2.5	0-20	LFM	
	37232	CARBARYL	모	12.1	12.4	\$	ug/L	121	124	70-130	2.4	0-20	LFM	
	37232	PROPOXUR (BAYGON)	2	, 11.7	12.2	5	ug/L	117	122	70-130	4.2	0-20	LFR	
	37232	METHIOCARB	2	4. 3.	=	10	ug/L	115	110	70-130	4.4	0-20	LFM	
	37232	BDMC (SURR)	110	105	109		%		NA NA	70-130	¥	0-20	LFM	_
531 090113												٠		
•	37742	OXYMAL	2	11.5	11.8	6	ngv.	115	118	70-130	2.6	0-20	LFM	
	37742	CARBOFURAN	2	11.1	-	6	ug/L	111	110	70-130	6'0	05-0	LFM	_
	37742	ALDICARB SULFOXIDE	2	11.7	12.1	우	ugu	117	121	70-130	9.4 4.	0-20	LFM	_
	37742	ALDICARB SULFONE	2	12.1	11.7	무	ng/L	121	117	70-130	3.4	99-0	LFM	
	37742	METHOMYL	Ş	12.1	11.9	무	ug/L	121	119	70-130	1.7	05-0	<u> </u>	
	37742	3-HYDROXYCARBOFURAN	2	11.8	11.1	₽	ug/L	118	11	70-130	6.1	0-20	LFR	_
	37742	ALDICARB	2	11.7	11.7	무	og/L	117	117	70-130	0.0	0-20	E.	_
	37742	CARBARYL	S	11,7	11.7	욘	ng/L	117	117	70-130	0.0	99	EA	-
	37742	PROPOXUR (BAYGON)	Ð	11.8	11.6	5	ug/L	118	116	70-130	1.7	0-20	LFR	_
	37742	METHIOCARB	2	10.9	#. #:	5	ug/L	109	115	70-130	5.4	0-20	E	-
	37742	BDMC (SURR)	103	102	5		*		¥	70-130	¥	0-50	<u> </u>	-
COD_081229														
ı	37235	CHEMICAL OXYGEN DEMAND	2	52	55	20	ш д/L	12	110	80-120	<u>.</u>	- -	E.	_
D081230A			!	;		!		1	:	ļ	;	8	į	
	37233	BROMATE	2	0.008		0.010	mg/L	96	⊈	(2-125	€ Z	3	ב.	_
	37255	BROMATE	9	0.0108		0.010	шg/L	106	₹	75-125	¥	8	Ē	₹
1081217A										;	;		į	
	37252	NITRATE-N	읒	1.04		8.	mg/L	104	¥	80-120	Ą Z	8	E L	•
	37252	CHLORIDE	8	31		1,00	т9/.	9	ΥN	80-120	¥	<u>8</u>		
	37278	NITRATE-N	0.77	2.84		1.00	mg/L	202	٧	80-120	ž	0 9- 0	M LFM	A Chlorinated

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Reference Number: 08-17751 Page 7 of 7

Report Date: 1/28/2009

pike			Duplicate										
		Spike	Spike	Spike		Percent Recover	RECOVERY				oc		
Sample Analyte	Result	Result	Result	Conc	Units	MS	MSD	Limits	%RPD Limits	Limits	Quelifier	Comments	
117 37235 ORTHO-PHOSPHATE	0.16	1.20	1.20	1.00	mg/L	104	104	70-130	9.6	0-20	LFM		1

OPHOS-081217

Balch

%RPD = Relative Percent Difference

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Matrix Spike (MS)/Matrix Spike Duplicate (MSD) and precision (MSD) of a analytical method in a given sample matrix. Therefore, the usefulness of this report is limited to samples of similar matrices analytical batch.



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Qualifier Definitions

Reference Number: 08-17751

Report Date: 01/28/09

Qualifier	Definition
AC	Ambient contamination during fortification of samples.
B5	The compound was detected in the sample below the State Reporting Limit, result is biased high.
HQ	High QCS recovery due to increased detector response of the sample extract. The continuing calibration checks are within acceptance limits.
J	Indicates an estimated concentration. This occurs when an analyte concentration is below the calibration curve but is above the method detection limit.
M	Matrix induced bias assumed.
ME	Matrix spike shows a possible matrix induced bias. The LFB was within acceptance limits, results for this compound are suspect.
ML	Indicates mechanical loss during extraction.
N1	Acceptance limits have not been established, the limits listed are for guidance only.
S	Spiking amount was lower than the 5:1 spike to background (sample amount) basis for performance criteria. The reported criteria does not apply due to increased errors in measurement of both sample and spike concentration.

Note: Some qualifier definitions found on this page may pertain to results or QC data which are not printed with this report.

FORM: QualifierDets



Cornorate Office 800.755.9295 • 360.757.1400 • 360.757.1402ta-Bellingham WA 805 Orchard Dr Suite 4 - 982.

1360.671.0688 • 360.671.15771.



March 5, 2009

Page 1 of 1

Mr. Troy Baker Walla Walla Basin Watershed Council 810 South Main Street Milton-Freewater, OR 97862

RE: 09-02151 - Locher Road Recharge Sites

Dear Mr. Troy Baker,

Your project: Locher Road Recharge Sites, was received on Friday February 13, 2009.

All samples were analyzed within the accepted holding times, were appropriately preserved and were analyzed according to approved analytical protocols. The quality control data was within laboratory acceptance limits, unless specified in the QA reports.

If you have questions phone me at 800 755-9295.

Respectfully Submitted,

Lawrence | Henderson, PhD Director of Laboratories

Enclosures Data Report

QC Reports Chain of Custody



Burington WA. 1 1620 S Walnut St - 9820.

Converse: Other 800.755.9295 • 360.757.1400 • 360.757.1400

Bellingnam WA | 805 Orchard Dr Suite 4 - 9222 Micromata | 360.671.0688 • 360.671.1577.

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Data Report

Client Name: Walla Walla Basin Watershed Council

810 South Main Street

Milton-Freewater, OR 97862

Reference Number: 09-02151

Project: Locher Road Recharge Sites

Report Date: 3/5/09

Date Received: 2/13/09

Peer Review:

Sample Description: L1 - Locher 1 Sai
Lab Number: 4470 Co

Sample Date: 2/12/09 Collected By: Unknown

Parameter	Result	PQL	MDL	Unița	DF	Method	Analyzed	Anaiy	st Batch	Comment
BROMATE	ND	0.005	0.0016	mg/L	1	300.1	3/4/00	MVP	D090304A	
HARDNESS	161	3.30	0.055	mg CaCO3/L	1	200.7	2/16/09	ВI	200.7-090216A	
NITRATE-N	6.08	0.100	0.015	mg/L	1	200,0	2/13/09	BJ	(000213A	
CHEMICAL OXYGEN DEMAND	ND	8.0	2.47	mg/L	1	SM5220 D	2/16/09	MAK	COD_000216	
TOTAL DISSOLVED SOLIDS	259	10	6	mg/L	t	SM2540 C	2/18/09	CCN	TD8_090215	
CHLORIDE	6.7	0.1	0.012	mg/L	1	300.0	2/13/09	롸	1000213A	
ORTHO-PHOSPHATE	0.23	0.01	0.002	mg/L	1	SM4500-P F	2/13/09	so	OPHOS-000213	
HYDROGEN ION (pH)	6.78			pH Units	1	SM4500-H+ B	2/13/09	CON	PH_090213	
TURBIDITY	1.78	0.05	0.03	NTU	•	180.1	2/13/09	CCN	TURB_090213	
ELECTRICAL CONDUCTIMITY	418	10		uS/cm	1	SM2510 B	2/16/09	CCN	EC_090216	
	Parameter BROMATE HARDNESS NITRATE-N CHEMICAL OXYGEN DEMAND TOTAL DISSOLVED SOLIDS CHLORIDE ORTHO-PHOSPHATE HYDROGEN ION (pH) TURBIDITY	Parameter Result BROMATE ND HARDNESS 161 NITRATE-N 6.08 CHEMICAL OXYGEN DEMAND ND TOTAL DISSOLVED SOLIDS 259 CHLORIDE 6.7 ORTHO-PHOSPHATE 0.23 HYDROGEN ION (pH) 6.78 TURBIDITY 1.78	Parameter Result PQL BROMATE ND 0.005 HARDNESS 161 3.30 NITRATE-N 6.08 0.100 CHEMICAL OXYGEN DEMAND ND 8.0 TOTAL DISSOLVED SOLIDS 259 10 CHLORIDE 6.7 0.1 ORTHO-PHOSPHATE 0.23 0.01 HYDROGEN ION (pH) 6.78 TURBIDITY 1.78 0.05	Parameter Result PQL MOL BROMATE ND 0.005 0.0016 HARDNESS 161 3.30 0.055 NITRATE-N 6.08 0.100 0.015 CHEMICAL OXYGEN DEMAND ND 8.0 2.47 TOTAL DISSOLVED SOLIDS 259 10 6 CHLORIDE 6.7 0.1 0.012 ORTHO-PHOSPHATE 0.23 0.01 0.002 HYDROGEN ION (pH) 6.78 TURBIDITY 1.78 0.05 0.03	Parameter Result PQL MDL Units BROMATE ND 0.005 0.0016 mg/L HARDNESS 161 3.30 0.055 mg NITRATE-N 6.08 0.100 0.015 mg/L CHEMICAL OXYGEN DEMAND ND 8.0 2.47 mg/L TOTAL DISSOLVED SOLIDS 259 10 6 mg/L CHLORIDE 6.7 0.1 0.012 mg/L ORTHO-PHOSPHATE 0.23 0.01 0.002 mg/L HYDROGEN ION (pH) 6.78 pH Units TURBIDITY 1.78 0.05 0.03 NTU	Parameter Result PQL MOL Units DF BROMATE ND 0.005 0.0016 mg/L 1 HARDNESS 161 3.30 0.055 mg 1 NITRATE-N 6.08 0.100 0.015 mg/L 1 CHEMICAL OXYGEN DEMAND ND 8.0 2.47 mg/L 1 TOTAL DISSOLVED SOLIDS 259 10 6 mg/L 1 CHLORIDE 6.7 0.1 0.012 mg/L 1 ORTHO-PHOSPHATE 0.23 0.01 0.002 mg/L 1 HYDROGEN ION (pH) 6.78 pH Units 1 TURBIDITY 1.78 0.05 0.03 NTU 1	Parameter Result PQL MDL Units DF Method BROMATE ND 0.005 0.0016 mg/L 1 300.1 HARDNESS 161 3.30 0.055 mg 1 200.7 NITRATE-N 6.08 0.100 0.015 mg/L 1 300.0 CHEMICAL OXYGEN DEMAND ND 8.0 2.47 mg/L 1 SM5220 D TOTAL DISSOLVED SOLIDS 259 10 6 mg/L 1 SM2540 C CHLORIDE 6.7 0.1 0.012 mg/L 1 300.0 ORTHO-PHOSPHATE 0.23 0.01 0.002 mg/L 1 SM4500-P F HYDROGEN ION (pH) 6.78 pH Units 1 SM4500-H+B TURBIDITY 1.78 0.05 0.03 NTU 1 180.1	Parameter Result PQL MDL Units DF Method Analyzed	Parameter Result PQL MDL Units DF Method Analyzed Analyzed Analyzed BROMATE ND 0.005 0.0016 mg/L 1 300.1 34/00 MVP	Parameter Result PQL MDL Units DF Method Analyzed Analyst Batch

•	ription: L2 - Locher 2 umber: 4471						•	ole Date: : cted By: !			
CAS ID#	Parameter	Result	PQL	MDL	Units	DF	Method	Analyzed	Analys	t Batch	Comment
15541-45-4	BROMATE	ND	0.005	0.0016	mg/L	1	300.1	3/4/09	MVP	D090304A	·
E-11778	HARDNESS	109	3.30	0.055	mg CaCO3/L	1	200.7	2/16/09	BJ	200.7-090216A	
14797-55-8	NITRATE-N	4.25	0.100	0.015	mg/L	1	300.0	2/14/09	BJ	1090213A	
E-10117	CHEMICAL OXYGEN DEMAND	ND	0.8	2.47	mg/L	1	SM5220 D	2/18/09	MAK	COD_090216	
E-10173	TOTAL DISSOLVED SOLIDS	193	10	6	mg/L .	1	SM2540 C	2/18/09	CCN	TDS_090218	
16887-00-6	CHLORIDE	5.8	0.1	0.012	mg/L	1	300.0	2/13/09	BJ.	1090213A	
14265-44-2	ORTHO-PHOSPHATE	0.23	0.01	0.002	mg/L	1	SM4500-P F	2/13/09	so	OPHOS-090213	
E-10139	HYDROGEN ION (pH)	6.77			pH Units	1	SM4500-H+ B	2/13/09	CCN	PH_000213	
E-10617	TURBIDITY	0.76	0.05	0.03	NTU	1	180.1	2/13/09	CCH	TURB_090213	
E-10184	ELECTRICAL CONDUCTIVITY	301	10		uS/cm	1	SM2510 B	2/16/09	CCN	EC_090216	

Notes

D.F. - Dilution Factor

ND = Not detected above the listed practical quantitation limit (PQL) or not above the Method Detection Limit (MDL), if requested

PQL = Practical Quantitation Limit is the lowest level that can be acheived within specified limits of precision and accuracy during routine laboratory operating conditions



Page 2 of 2 Reference Number: 09-02151 Report Date:3/5/09

Data Report

•	ription: L3 - Locher 3 umber: 4472						•	ile Date: 2 cled By: \		វា	
AS ID#	Parameter	Result	PQL	MDL	Units	DF	Method	Analyzed	Analysi	Batch	Comment
5541-45-4	BROMATE	ND	0.005	0.0016	mg/L	1	300.1	3/4/09	MVP I	0000304A	_
-11778	HARDNESS	84.4	3.30	0.055	mg CeCO3/L	1	200.7	2/16/09	B) 2	100.7-090216A	
4797-55-8	NITRATE-N	5.36	0.100	0.015	mg/L	1	300.0	2/14/00	BJ I	090213A	
-10117	CHEMICAL OXYGEN DEMAND	ND	8.0	2.47	mg/L	1	SM5220 D	2/16/09	MAK (COD_090216	
-10173	TOTAL DISSOLVED SOLIDS	169	10	6	mg/L	,	SM2540 C	2/15/09	CCN '	TD\$_090218	
6887-00-6	CHLORIDE	4.2	0.1	0.012	mg/L	1	300.0	2/13/00	BJ I	090213A	
4265-44-2	ORTHO-PHOSPHATE	0.21	0.01	0.002	mg/L	1.	SM4500-P F	2/13/09	60 (OPHOS-090213	
-10139	HYDROGEN (ON (pH)	6.81			pH Units	1	SM4500-H+ B	2/13/09	CCN I	H_090213	
-10617	TURBIDITY	1.62	0.05	0.03	NTU	3	180.1	2/13/09	CCN 1	TUFIB_090213	
10184	ELECTRICAL CONDUCTIVITY	231	10		uS/cm		SM2510 B	2/16/09	CCN I	C_090216	

•	ription: LSW1 - Locher SW1 lumber: 4473							ole Date:			
CAS ID#	Parameter	Result	PQL	MDL	Units	DF	Method	Analyzed	Analys	st Batch	Comment
15541-45-4	BROMATE	ND	0.005	0.0016	mg/L	1	300.1	3/4/00	MVP	D090384A	
E-11778	HARDNESS	141	3.30	0.055	mg CaCO3/L	1	200.7	2/16/09	BJ	200.7-090218A	
14797-55-8	NITRATE-N	2.57	0.100	0.015	mg/L	1	300.0	2/14/09	BJ	1090213A	
E-10117	CHEMICAL OXYGEN DEMAND	ND	8.0	2.47	mg/L	1	SM5220 D	2/16/09	HAK	COD_090215	
E-10173	TOTAL DISSOLVED SOLIDS	224	10	6	mg/L	1	SM2540 C	2/18/00	CCN	TDS_090218	
16887-00-6	CHLORIDE	7.2	0.1	0.012	mg/L	1	300.0	2/13/09	BN	I090213A	
14265-44-2	ORTHO-PHOSPHATE	0.18	0.01	0.002	mg/L	,	SM4500-P F	2/13/00	so	OPHOS-090213	•
E-10139	HYDROGEN ION (pH)	7.63			pH Units	1	SM4500-H+ B	2/13/09	CCN	PH_000213	
E-10817	TURBIDITY	0.83	0.05	0.03	NTU	T	180.1	2/13/09	CCN	TURB_090213	-
E-10184	ELECTRICAL CONDUCTIVITY	378	10		uS/cm	1	SM2510 B	2/16/09	CCN	EC_090216	

Notes:

ND = Not detected above the listed practical quantitation limit (POL) or not above the Method Detection Limit (MDL), if requested.

PQL = Practical Quantitation Limit is the lowest level that can be acheived within specified limits of practical quantitation Limit is the lowest level that can be acheived within specified limits of practical quantitation Limit.

D.F. - Dilution Factor



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HERBICIDES IN DRINKING WATER

Client Name: Walla Walla Basin Watershed Council

810 South Main Street

Milton-Freewater, OR 97862

Project::

Field ID: L1

Sample Description: Locher 1

Sampled By: Unknown

Sample Date: 2/12/09 Source Type:

Sampler Phone:

Lab Number: 04470

Report Date: 2/26/09

. Reference Number: 09-02151

Date Analyzed: 02/20/09

Date Extracted: 515.4_090219

Analyst:

Peer Review. Analytical Method: 515.4

Herbicides in Drinking Water

Project: Locher Road Recharge Sites

CAS	COMPOUND	RESULTS	UNITS	PQL	MDL	MCL	COMMENT
EPA Re	egulated				-		
94-75-7	2,4 - D	ND	ug/L	0.5	0.2	70	
93-72-1	2,4,5 - TP (SILVEX)	ND	ug/L	0.25	0.04	50	
87-86-5	PENTACHLOROPHENOL.	ND	ug/L	0.25	0.05	1	
75-99-0	DALAPON	ND	ug/L	3.25	0.9	200	
88-85-7	DINOSEB	ИD	ug/L	0.5	0.2	7	
1918-02-1	PICLORAM	ND	ug/L	0.25	0.07	500	
Other							
E-14028	DCPA (ACID METABOLITES)	ND	ug/L	0.25	0.1		•
1918-00- 9	DICAMBA	NÐ	ug/L	0.25	0.05		
94-82-6	2,4 DB	ND	ug/L	2			
93-76-5	2,4,5 T	ND	ug/L	0.25	0.04		
25057-89-0	BENTAZON	ND	ug/L	0.5	0.2		
120-36-5	DICHLORPROP	ND	ug/L	0.75	0.5		
50594-66-6	ACIFLUORFEN	ND	ug/L	0.25	0.1		
51-3 6 -5	3,5 - DICHLOROBENZOIC ACID	ND	ug/L	0.25	0.2		

NOTES:
If a compound is detected > or = to the State Reporting Level, SRL, specified increased mositoring frequencies may occur par DOH.
MCL (Maximum Conteminant Level) maximum pormissible level of a contaminant in water established by EPA; Federal Action Levels are 0.015 mg/L for Lead and 1.3 mg/L for Copper. Sedium has a recommended limit of 20 mg/L. A blank MCL value indicates a level is not currently established.
Trigger Level: DOH Drinking Water Respirate level. Systems with compounds detected in excess of this level are required to take additional samples. Contact your regional DOH office.

ND (Not Detected): indicates that the parameter was not detected above the State Reporting Limit (SRL).



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HERBICIDES IN DRINKING WATER

Client Name: Walla Walla Basin Watershed Council

810 South Main Street

Milton-Freewater, OR 97862

Project::

Field ID: L2

Sample Description: Locher 2

Sampled By: Unknown

Sample Date: 2/12/09

Source Type:

Sampler Phone:

Reference Number: 09-02151

Project: Locher Road Recharge Sites

Lab Number: 04471

Report Date: 2/26/09

Date Analyzed: 02/20/09

Date Extracted: 515.4_090219

Analyst: CØ

Peer Review:

Analytical Method: 5\5.4

Herbicides in Drinking Water

CAS	COMPOUND	RESULTS	UNITS	PQL	MDL	MCL	COMMENT
EPA R	egulated						
94-75-7	2,4 - D	ND	ugň.	0.5	0.2	70	
93-72-1	2,4,5 - TP (SILVEX)	ND	ug#L	0.25	0.04	50	
37-86- 5	PENTACHLOROPHENOL	ND	n ā ⁄ι	0.25	0.05	1	
75-99-0	DALAPON	ND	u g/L	3.25	0.9	200	
88-85-7	DINOSEB	ND	u g/L	0.5	0.2	7	
1918-02-1	PICLORAM	ND	⊔ g/L	0.25	0.07	500	
Other							
E-1402B	DCPA (ACID METABOLITES)	ND	ug/L	0.25	0.1		
918-00-9	DICAMBA	ND	u g/ L	0.25	0.05		
4-82-6	2.4 DB	ND	ug/L				qualitative analysis
3-76-5	2,4,5 T	ND	ug/L	0.25	0.04		
25057-89-0	BENTAZON	ND	ug/L	0.5	0.2		
20-36-5	DICHLORPROP	ND	ug/L	0.75	0.5		
50594-66-6	ACIFLUORFEN	ND	ug/L	0.25	0.1		
1-36-5	3,5 - DICHLOROBENZOIC ACID	ND	ug/L	0.25	0.2		

NOTES:
If a compound is detected > or w to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.

MCL (Maximum Contaminant Level) maximum permissible level of a contaminant in water established by EPA; Federal Action Levels are 0.015 mg/L for Lead and 1.3 mg/L for Copper. Sodium has a recommended Smit of 20 mg/L. A blank MCL value indicates a level is not currently established.

Trigger Level: DOH Orinking Water Response level. Systems with compounds detected in excess of this level are required to take additional samples. Contact your regional DOH office.

NO (Not Detected): indicates that the parameter was not detected above the State Reporting Limit (SRL).



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HERBICIDES IN DRINKING WATER

Client Name: Walla Walla Basin Watershed Council

810 South Main Street

Milton-Freewater, OR 97862

Prolect::

Field ID: L3

Sample Description: Locher 3

Sampled By: Unknown

Sample Date: 2/12/09

Source Type:

Sampler Phone:

Reference Number: 09-02151

Project: Locher Road Recharge Sites

Lab Number: 04472

Report Date: 2/26/09

Date Analyzed: 02/20/09

Date Extracted: 515.4_090219

Analyst: CQ

Peer Review: Analytical Method: 515.4

Herbicides in Drinking Water

CAS	COMPOUND	RESULTS	UNITS	PQL	MDL	MCL	COMMENT
EPA R	egulated						······································
94-75-7	2,4 - D	ND	ug/L	0.5	0.2	70	
93-72-1	2,4,5 - TP (SILVEX)	ND	ug/L	0.25	0.04	50	
37-86-5	PENTACHLOROPHENOL	ND	ug/L	0.25	0.05	1	·
75-99-0	DALAPON	ND	ug/L	3.25	0.9	200	
38-8 5-7	DINOSEB	ND	ug/L	0.5	0.2	7	
1918-02-1	PICLORAM	ND	ug/L	0.25	0.07	500	
Other							
-14028	DCPA (ACID METABOLITES)	ND	ug/L	0.25	0.1		
918-00-9	DICAMBA	ND	ug/L	0.25	0.05		
4-82-6	2,4 DB	ND	ug/L				qualitative analysis
3-76-5	2,4,5 T	ND	ug/L	0.25	0.04		•
5057-89-0	BENTAZON	ND .	ug/L	0.5	0.2		
20-36-5	DICHLORPROP	ND	ug/L	0.75	0.5		•
0594-66-6	ACIFLUORFEN	ND	ug/L	0.25	0.1		
1-36-5	3,5 - DICHLOROBENZOIC ACID	ND	ug/L	0.25	0.2		

NOTES:
If a compound is detected > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOM.

MCL (Maximum Contaminant Level) maximum permissible level of a contaminant in water established by EPA; Federal Action Levels are 0.015 mg/L for Lead and 1.3 mg/L for Coppur. It blank MCL value indicates a level is not currently established.

Trigger Level: DOH Drinking Water Response level. Systems with compounds detected in excess of this level are required to take additional samples. Contact your regional DOH office. d is detected > or = to line State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOM.

Unification in the state Reporting Level, SRL, specified increased monitoring frequencies may occur per DOM.

Levels are 0.015 mg/L for Lead and 1.3 mg/L for Copper. Socium has a recommended limit of 20 mg/L. A

ND (Not Detected): indicates that the parameter was not detected above the State Reporting Limit (SRL).



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SYNTHETIC ORGANIC COMPOUNDS (SOC) REPORT

Client Name: Walla Walla Basin Watershed Council

810 South Main Street

Milton-Freewater, OR 97862

Project::

Field ID: L1

Sample Description: Locher 1

Sampled By: Unknown

Sample Date: 2/12/09

Source Type:

Sampler Phone:

Reference Number: 09-02151

Project: Locher Road Recharge Sites

Lab Number: 04470

Report Date: 2/24/09

Date Analyzed: 02/19/09

Date Extracted: 525_090217

Analyst: CQ

Peer Review.

Analytical Method: 525.2

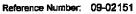
Synthetic Organics

CAS -	COMPOUND	RESULTS	UNITS	, PQL	MDL	MCL	COMMENT
EPA Re	gulated						
72-20-8	ENDRIN	ND	ug/L	0.1	0.030	2	
58-89-9	LINDANE (BHC - GAMMA)	ND	ug/L	0.1	0.028	0.2	
72-43-5	METHOXYCHLOR	ND	ug/L	0.1	0.015	40	
15972-60-8	ALACHLOR	ND	ug/L	0.1	0.044	2	
1912-24-9	ATRAZINE	0.04 J	ոգմե	0.1	0.030	3	
50-32-8	BENZO(A)PYRENE	ND	ug/L	0.1	0.012	0.2	
57-74-9	CHLORDANE, TECHNICAL	ND	ug/L	0.1	0.3	2	
103-23-1	DI(ETHYLHEXYL)-ADIPATE	ND	ug/L	0.1	0.022	400	
117-81-7	DI(ETHYLHEXYL)-PHTHALATE	ND	ug/L	0.1	0.063	6	
76-44-8	HEPTACHLOR	ND	ug/i.	0.1	0.022	0.4	
1024-57-3	HEPTACHLOR EPOXIDE	ND	ug/L	0.1	0.02	0.2	
118-74-1	HEXACHLOROBENZENE	ND	ug/L	0.1	0.025	1	
77-47-4	HEXACHLOROCYCLO-PENTADIENE	ND	ug/L	0.1	0.024	50	
122-34-9	SIMAZINE	ND	ug/L	0.1	0.030	4	
37-86-5	PENTACHLOROPHENOL	ND	ug/L	0.4	0.08	1 .	screening only / compliance by 515,1
EPA Un	regulated						
309-00-2	ALDRIN	ND	ug/L	0.1	0.022		
23184-66-9	BUTACHLOR	ND	ug/L	0.1	0.024		
60-57-1	DIELDRIN	ND	ug/L	0.1	0.031		
51218-45-2	METOLACHLOR	ND	ug/L	0.1	0.024		
21087-64-9	METRIBUZIN	ND	ug/L	0.1	0.030		
1918-16-7	PROPACHLOR	ND	ug/L	0.1	0.031		
State U	nregulated - Other						
314-40-9	BROMACIL	0.21 JJ	ug/L	0.1	0.031		

DRINK WILL Visual misseries a reversite from cannon my executation.

Trigger Lavel: DOH Drinking Water Response level. Systems with compounds detected in excess of this NO (Not Detected): indicates that the parameter was not detected above the State Reporting Limit (SRL).

If a compound is detected > or = to the State Reportin MCL (Maximum Contaminant Level) maximum permis blank MCL value indicates a level is not currently estal Lavel, SRL, specified increased monitoring frequencies may occur per DOH. Side level of a conteminant in water established by EPA; Federal Action Levels are 0.015 mg/L for Lead and 1.3 mg/L for Copper. Sodium has a recommended limit of 20 mg/L. A MCL value indicates a level is not currently established.





Lab Number: 04470

Report Date: 2/24/09 15:37



SYNTHETIC ORGANIC COMPOUNDS (SOC) REPORT

CAS	COMPOUND	RESULTS	UNITS	PQL	MDL	MCL	COMMENT
5902-51-2	TERBACIL	ND	ug/L	0.1	0.043		
333-41-5	DIAZINON	ND	ug/L	0.1	0.035		Unstable in Acidified Sample Matrix
759-94-4	EPTC	ND	ug/L	0.1	0.028 .		
72-54-8	4,4-DDD	ND	ug/L	0.1	0.024		
72-55-9	4,4-DDE	ND	ug/L	0.1	0.024		
50-29-3	4,4-DDT	ND	ug/L	0.1	0.022		
21725-46-2	CYANAZINE	ND	ug/L	0.1	0.13		Qualitative Analysis Only
21-75-5	MALATHION	ND	ug/L	0.1	0.015		
6-38-2	PARATHION	ND	ug/L	0.1	0.022		
1582-09-8	TRIFLURALIN	ND	ug/L	0.1	0.024		
- PAHs			•				
1-20-3	NAPTHALENE	ND	ug/L	0.1	0.1^		
6-73-7	FLUORENE	ND	ug/L	0.1	0.026		•
08-96-8	ACENAPHTHYLENE	ND	ug/L	0.1	0.025		
3-32-9	ACENAPHTHENE	ND	ug/L	0.1	0.1^		
20-12-7	ANTHRACENE	ND	ug/L	0.1	0.012		
6-55-3	BENZ(A)ANTHRACENE	ND	ug/L	0.1	0.012		
05-99-2	BENZO(B)FLUORANTHENE	ND	ug/L	0.1	0.025		
91-24-2	BENZO(G,H,I)PERYLENE	ND	ug/L	0.1	0.025		
07-08-9	BENZO(K)FLUORANTHENE	ND	ug/L	0.1	0.022	-	
18-01-9	CHRYSENE	ND	ug/L	0.1	0.022		
3-70-3	DIBENZO(A,H)ANTHRACENE	ND	ψg/L	0.1	0.024		
06-44-0	FLUORANTHENE	ND	ug/L	0.1	0.1^		
93-39-5	INDENO(1,2,3-CD)PYRENE	ND	ug/L	0.1	0.040		
5-01-8	PHENANTHRENE	ND	ug/L	0.1	0.015		
29-00-0	PYRENE	ND	ug/L	0.1	0.022		
- Phthai	lates						
5-68-7	BENZYL BUTYL PHTHALATE	ND	ug/L	0.1	0.022		
1-74-2	DI-N-BUTYL PHTHALATE	ND	ug/L	0.1	0.085		
4-6 6- 2	DIETHYL PHTHALATE	ND	· ug/L	0.1	0.044		
31-11-3	DIMETHYL PHTHALATE	ND	ug/L	0.1	0.015		

NOTES:

If a compound is detected > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.

MCL (Maximum Contaminant Level) maximum permissible level of a contaminant in water established by EPA; Federal Action Levels are 0.015 mg/L for Lead and 1.3 mg/L for Copper, Sodium has a recommended limit of 20 mg/L. A blank MCL value indicates a level is not currently established.

Trigger Level: DOH Dirinking Water Response level. Systems with compounds detected in excess of this level are required to take additional samples. Contact your regional DOH office.

ND (Not Detected): ladicates that the parameter was not detected above the State Reporting Umit (SRL).



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Page 1 of 2

SYNTHETIC ORGANIC COMPOUNDS (SOC) REPORT

Client Name: Walla Walla Basin Watershed Council

810 South Main Street

Milton-Freewater, OR 97862

Project::

Field ID: L2

Sample Description: Locher 2

Sampled By: Unknown

Sample Date: 2/12/09

Source Type:

Sampler Phone:

Reference Number: 09-02151

Project: Locher Road Recharge Sites

Lab Number: 04471

Report Date: 2/24/09 Date Analyzed: 02/19/09

Date Extracted: 525_090217

Analyst: r-QO Peer Review:

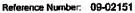
Analytical Method: 525.2

Synthetic Organics

CAS	COMPOUND	RESULTS	UNITS	PQL	MDL	MÇL	COMMENT.
EPA Re	egulated						
72-20-8	ENDRIN	ND	ug/L	0.1	0.030	2	
58 - 89-9	LINDANE (BHC - GAMMA)	ND	ug/L	0.1	0.028	0.2	
72-43-5	METHOXYCHLOR	ND	ug/L	0.1	0.015	40	•
15972-60-8	ALACHLOR	ND	ug/L	0.1	0.044	2	
1912-24-9	ATRAZINE	ND	ug/L	0.1	0.030	3	
50-32-8	BENZO(A)PYRENE	ND	ug/L	0.1	0.012	0.2	
57-74-9	CHLORDANE, TECHNICAL	ND	ug/L	0.1	0.3	2	
103-23-1	DI(ETHYLHEXYL)-ADIPATE	ND	ug/L	0.1	0.022	400	
117-81-7	DI(ETHYLHEXYL)-PHTHALATE	ND	ug/L	0.1	0.063	6	
76-44-8	HEPTACHLOR	ND	ug/L	0.1	0.022	0.4	
1024-57-3	HEPTACHLOR EPOXIDE	ND	ug/L	0.1	0.02	0.2	
118-74-1	HEXACHLOROBENZENE	ND	ug/L	0.1	0.025	1	
77-47-4	HEXACHLOROCYCLO-PENTADIENE	ND	ug/L	0.1	0.024	50	
122-34-9	SIMAZINE	ND	ug/L	0.1	0.030	4	
87-86-5	PENTACHLOROPHENOL	ND	ug/L	0.4	80.0	1 .	screening only / compliance by 515.1
EPA Un	ıregulated						
309-00-2	ALDRIN	ND	u g/L	0.1	0.022		
23184-66-9	BUTACHLOR	ND	u g/L .	0.1	0.024		
60-57-1	DIELDRIN .	ND	ug/L	0.1	0.031		
51218-45-2	METOLACHLOR	ND	ug/L	0.1	0.024		
21087-64-9	METRIBUZIN	ND	ug/L	0.1	0.030		
1918-16-7	PROPACHLOR	МÐ	ug/L	0.1	0.031		
State U	nregulated - Other					٠	•
314-40-9	BROMACIL	ND	ug/L	0.1	0.031		

g Level, SRL, specified increased monitoring frequencies may occur per DOH.

sebie level of a contaminant in water established by EPA; Federal Action Levels are 0.015 mg/L for Lead and 1.3 mg/L for Copper. Sodium has a recommended limit of 20 mg/L. A





Lab Number: 04471

Report Date: 2/24/09 15:37



SYNTHETIC ORGANIC COMPOUNDS (SOC) REPORT

CAS	COMPOUND	RESULTS	UNITS	PQL	MIDL,	MCL	COMMENT
5902-51-2	TERBACIL	ND	ug/L	0.1	0.043		
333-41-5	DIAZINON	ND	ug/L	0.1	0.035		Unstable in Acidified Sample Matrix
759-94-4	EPTC	ND	ug/L	0.1	0.028		
72-54-8	4,4-DDD	ND	ug/L	0.1	0.024		,
72-55-9	4,4-DDE	ND	ug/L	0.1	0.024		
50-2 9- 3	4,4-DDT	ND	ug/L	0.1	0.022		
21725-46-2	CYANAZINE	ND	ug/L	0.1	0.13		Qualitative Analysis Only
121-75-5	MALATHION	ND .	ug/L	0.1	0.015		
56-38-2	PARATHION	ND	ug/L	0.1	0.022		
1582-09-8	TRIFLURALIN	ND	ug/L	0.1	0.024		
- PAHs							
91-20-3	NAPTHALENE	ND	ug/L	0.1	0.1^		
36-73-7	FLUORENE	ND	ug/L	0.1	0.026		
208-96-8	ACENAPHTHYLENE	ND	ug/L 1	0.1	0.025		<u>-</u>
33-32-9	ACENAPHTHENE	ND	ug/L	0.1	0.1^		
120-12-7	ANTHRACENE	ND	ug/L	0.1	0.012		
56-55-3	BENZ(A)ANTHRACENE	₩D	ug/L	0.1	0.012		
205-99-2	BENZO(B)FLUORANTHENE	ND	ug/L	0.1	0.025		
191-24-2	BENZO(G,H,I)PERYLENE	ND	ug/L	0.1	0.025		
207-08-9	BENZO(K)FLUORANTHENE	ND	ug/L	0.1	0.022		
218-01-9	CHRYSENE	ND	ug/L	0.1	0.022		
53-70-3	DIBENZO(A,H)ANTHRACENE	ND	ug/L	0.1	0.024		•
206-44-0	FLUORANTHENE	ND	ug/L	0.1	0.1^		
93-39-5	INDENO(1,2,3-CD)PYRENE	ND	ug/L	0.1	0.040		
5-01-8	PHENANTHRENE	ND	ug/L	0.1	0.015		
29-00-0	PYRENE	ND	ug/L	0.1	0.022		
- Phthal	lates						
5-68-7	BENZYL BUTYL PHTHALATE	ND	ug/L	0.1	0.022		
4-74-2	DI-N-BUTYL PHTHALATE	ND	ug/L	0.1	0.085		
4-66-2	DIETHYL PHTHALATE	ND	ug/L	0.1	0.044		
31-11-3	DIMETHYL PHTHALATE	ND	ug/L	0.1	0.015		

NOTES:
If a compound is detected > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.

MCI. (Maximum Contaminant Level) maximum permissible level of a contaminant in water established by EPA; Federal Action Levels are 0.015 mg/L for Lead and 1.3 mg/L for Copper. Sodium has a recommended limit of 20 mg/L. A blank MCI. Value indicates a level is not currently established.

Trigger Level: DOH Drinking Water Respense level. Systems with compounds detected in excess of this level are required to take additional samples. Contact your regional DOH office.



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SYNTHETIC ORGANIC COMPOUNDS (SOC) REPORT

Client Name: Walla Walla Basin Watershed Council

810 South Main Street

Milton-Freewater, OR 97862

Project::

Field ID: L3

Sample Description: Locher 3

Sampled By: Unknown

Sample Date: 2/12/09

Source Type:

Sampler Phone:

Reference Number: 09-02151

Project: Locher Road Recharge Sites

Lab Number: 04472

Report Date: 2/24/09

Date Analyzed: 02/19/09

Date Extracted: 525 090217

Analyst:

Peer Review:

Analytical Method: 525.2

Synthetic Organics

-CAS	COMPOUND	RESULTS	UNITS	PQL	MDL	MCL	COMMENT
EPA Re	egulated						
72-20-8	ENDRIN	ND	ug/L	0.1	0.030	2	
58-89-9	LINDANE (BHC - GAMMA) .	ND	ug/L	0.1	0.028	0.2	
72-43-5	METHOXYCHLOR	ND	ug/L	0.1	0.015	40	
15972-60-8	ALACHLOR	ND	ug/L	0.1	0.044	2	
1912-24-9	ATRAZINE	ND	ug/L	0.1	0.030	3	
50-32-8	BENZO(A)PYRENE	ND	ug/L	0.1	0.012	0.2	
57-74-9	CHLORDANE, TECHNICAL	ND	ug/L	0.1	0.3	2	
103-23-1	DI(ETHYLHEXYL)-ADIPATE	ND	ug/L	0.1	0.022	400	
117-81-7	DI(ETHYLHEXYL)-PHTHALATE	ND	ug/L	0.1	0.063	6	
76-44-8	HEPTACHLOR	ND	ug/L	0.1	0.022	0.4	
1024-57-3	HEPTACHLOR EPOXIDE	ND	ug/L	0.1	0.02	0.2	
118-74-1	HEXACHLOROBENZENE	ND	ug/L	0.1	0.025	1	
77-47-4	HEXACHLOROCYCLO-PENTADIENE	ND	ug/L	0.1	0.024	50	
122-34-9	SIMAZINE	ND	ug/L	0.1	0.030	4	
87-86-5	PENTACHLOROPHENOL	ND	ug/L	0.4	0.08	1	screening only / compliance by 515.1
EPA Ur	nregulated						
309-00-2	ALDRIN	ND	ug/L	0,1	0.022		
23184-66-9	BUTACHLOR	ND	ug/L	0.1	0.024		
60-57-1	DIELDRIN	ND	ug/L	0.1	0.031		
51218-45-2	METOLACHLOR	ND	ug/L	0.1	0.024		
21087-64-9	METRIBUZIN	ND	ug/L	0.1	0.030		
1918-16-7	PROPACHLOR	ND	ug/L	0.1	0.031		
State U	nregulated - Other						
314-40-9	BROMACIL	ND	ug/L	0.1	0.031		

NOTES:
If a compound is detected > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.

MCL (Maximum Contaminant Level) maximum permissible level of a contaminant in water established by EPA; Federal Action Levels are 0.015 mg/L for Level and 1.3 mg/L for Copper. Sodium has a recommended limit of 20 mg/L. A blank MCL vature indicates a level is not currently established.

Trigger Level: DOH Diriking Water Response level. Systems with compounds detected in excess of this level are required to take additional samples. Contact your regional DOH office.

ND (Not Detected): indicates that the parameter was not detected above the State Reporting Limit (SRL).





Lab Number: 04472

Report Date: 2/24/09 15:37

SYNTHETIC ORGANIC COMPOUNDS (SOC) REPORT

CAS	COMPOUND	RESULTS	UNITS	PQL	MDL	MCL	COMMENT
5902-51-2	TERBACIL	ND	ug/L	0.1	0.043		
333-41-5	DIAZINON	ND	ug/L	0.1	0.035		Unstable in Ackliffed Sample Matrix
759-94-4	EPTC	ND	ug/L	0.1	0.028		
72-54-8	4,4-DDD	ND	ug/L	0.1	0.024		
72-55-9	4,4-DDE	ND	ug/L	0.1	0.024		
50-2 9- 3	4,4-DDT	ND	ug/L	0.1	0.022		
21725-46-2	CYANAZINE	ND	ug/L	0.1	0.13		Qualitative Analysis Only
121-75-5	MALATHION	ND	ug/L	0.1	0.015		
56-38-2	PARATHION	ND	ug/L	0.1	0.022		
1582-09-8	TRIFLURALIN	ND	ug/L	0.1	0.024		
- PAHs	ì						
91-20-3	NAPTHALENE	ND	ug/L	0.1	0.1^		
36-73-7	FLUORENE	ND	ug/L	0.1	0.026		
208-96-8	ACENAPHTHYLENE	ND	'ug/L	0.1	0.025		•
33-32-9	ACENAPHTHENE	ND	ug/L	0.1	0.1^		
120-12-7	ANTHRACENE	ND	ug/L	0.1	0.012		
6-55-3	BENZ(A)ANTHRACENE	ND	ug/L	0.1	0.012		
205-99-2	BENZO(B)FLUORANTHENE	ND	ug/L	0.1	0.025		•
91-24-2	BENZO(G,H,I)PERYLENE	ND	ug/L	0.1	0.025		
07-08-9	BENZO(K)FLUORANTHENE	ND	ug/L	0.1	0.022		
:18-01- 9	CHRYSENE	ND	ug/L	0.1	0.022		
3-70-3	DIBENZO(A,H)ANTHRACENE	ND	ug/L	0.1	0.024		
06-44-0	FLUORANTHENE	ND	ug/L	0.1	0.1^		
93-39-5	INDENO(1,2,3-CD)PYRENE	ND	ug/L	0.1	0.040		
5-01-8	PHENANTHRENE	ND	ug/L	0.1	0.015		
29-00-0	PYRENE	ND .	ug/L	0.1	0.022		
- Phtha	lates						
5-68-7	BENZYL BUTYL PHTHALATE	ND	ug/L	0.1	0.022		
4-74-2	DI-N-BUTYL PHTHALATE	ND	ug/L	0.1	0.085		
4-66-2	DIETHYL PHTHALATE	ND	ug/L	0.1	0.044		
31-11-3	DIMETHYL PHTHALATE	ND	ug/L	0.1	0.015		

NOTES:
If a compound is defected > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.

MCI. (Maximum Contaminant Level) maximum permissible level of a contaminant is water established by EPA; Federal Action Levels are 0.015 mg/L for Lead and 1.3 mg/L for Copper. Sodium has a recommended limit of 20 mg/L. A blank MCI. value indicates a level is not currently established.

Trigger Level. DOH Drinking Water Response level. Systems with compounds detected in excess of this level are required to take additional samples. Contact your regional DOH office.

ND (Not Detected): Indicates that the parameter was not detected above the State Reporting Limit (SRL).



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SYNTHETIC ORGANIC COMPOUNDS (SOC) REPORT

Client Name: Walla Walla Basin Watershed Council

810 South Main Street

Milton-Freewater, OR 97862

Project::

Field ID: L1

Sample Description: Locher 1

Sampled By: Unknown

Sample Date: 2/12/09 Source Type:

Sampler Phone:

Reference Number: 09-02151

Project: Locher Road Recharge Sites

Lab Number: 04470

Report Date: 2/24/09 Date Analyzed: 02/17/09

Date Extracted: 508_090217

Analyst: GEB

Peer Review.

Analytical Method: 508.

Synthetic Organics

CAS	COMPOUND	RESULTS	UNITS	PQL	MDL	MCL	COMMENT
PCBs/1	Гохарћеле						
1336-36-3	PCBS (Total Arcclors)	ND	ug/iL	0.2		0.5	
11104-28-2	AROCLOR 1221	ND	ug/L	0.1	0.1^		
11141-16-5	AROCLOR 1232	ŃD	ug/L	0.1	0.1^		
53469-21-9	AROCLOR 1242	ND	ug/L	0.1	0.1^		
12672-29-6	AROCLOR 1248	ND	u g/ Ľ	0.1	0.1^		
11097-69-1	AROCLOR 1254	ND	ug/L	0.1	0.1^		
11096-82-5	AROCLOR 1260	ND	ug/L	0.1	0.08		
12674-11-2	AROCLOR 1016	ND	ug/L	0.1	0.1		
8001-35-2	TOXAPHENE	ND	ug/L	1	0.5	3	•

NOTES:

If a compound is detected > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.

MCI. (Maximum Contaminant Level) maximum permissible level of a centaminant in water established by EPA; Federal Action Levels are 0.015 mg/L for Lead and 1.3 mg/L for Copper. Sedium has a recommended limit of 20 mg/L. A blank MCL value indicates a level is not correspond established.

Tripger Level: DOH Dirikhing Water Response level. Systems with compounds detected in excess of this level are required to take additional samples. Contact your regional DOH office.

ND (Not Detected): Indicates that the parameter was not detected above the State Reporting Unit (SRL).



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Page 1 of 1

SYNTHETIC ORGANIC COMPOUNDS (SOC) REPORT

Client Name: Walla Walla Basin Watershed Council

810 South Main Street Milton-Freewater, OR 97862

Project::

Field ID: L2

Sample Description: Locher 2

Sampled By: Unknown Sample Date: 2/12/09

Source Type:

Sampler Phone:

Reference Number: 09-02151

Project: Locher Road Recharge Sites

Lab Number: 04471 Report Date: 2/24/09 Date Analyzed: 02/17/09

Date Extracted: 508 090217

Analyst: GEB ->>

Peer Review. Analytical Method: 508.T

Synthetic Organics

CAS	COMPOUND	RESULTS	UNITS	PQL	MDL	MCL	COMMENT
PCBs/	Toxaphene					.,	
1336-36-3	PCBS (Total Aroclors)	ND	ug/L	0.2		0.5	
11104-28-2	AROCLOR 1221	ND	ug/L	0.1	0.1^		
11141-16-5	AROCLOR 1232	ND	ug/L	0.1	0.1^		
53469-21-9	AROCLOR 1242	ND	ug/L	0.1	0.1^		
12672-29-6	AROCLOR 1248	ND	ug/L	0.1	0.1^		
11097 -69 -1	AROCLOR 1254	ND	ug/L	0.1	0.1^		
11096-82-5	AROCLOR 1260	ND.	ug/L	0.1	0.08		
12674-11-2	AROCLOR 1016	ND	ug/L	0.1	0.1		
8001-35-2	TOXAPHENE	ND	ug/L	1	0.5	3	

ncreased monitoring trequencies may occur per DOH. sent in water established by EPA; Federal Action Levels are 0.015 mg/L for t.ead and 1.3 mg/l. for Copper, Sodium has a recommended limit of 20 mg/L. A



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SYNTHETIC ORGANIC COMPOUNDS (SOC) REPORT

Client Name: Walla Walla Basin Watershed Council

810 South Main Street

Milton-Freewater, OR 97862

Project::

Field ID: L3

Sample Description: Locher 3

Sampled By: Unknown

Sample Date: 2/12/09

Source Type:

Sampler Phone:

Reference Number: 09-02151

Project: Locher Road Recharge Sites

Lab Number: 04472

Report Date: 2/24/09

Date Analyzed: 02/17/09

Date Extracted: 508 090217

Analyst: GEB

Peer Review:

Analytical Method: 508.1

Synthetic Organics

CAS	COMPOUND	RESULTS	UNITS	PQL	MDL	MCL	COMMENT
PCBs/I	Toxaphene				. , , , , , , , , , , , , , , , , , , ,		
1336-36-3	PCBS (Total Aroclors)	ND	ug/L	0.2		0.5	
11104-28-2	AROCLOR 1221	ND	ug/L	0.1	0.14		
11141-16-5	AROCLOR 1232	ND	ug/L	0.1	0.1^		
53469-21-9	AROCLOR 1242	ND	ug/L	0.1	0.1^		
12672-29-6	AROCLOR 1248	ND	ug/L	0.1	0.1^		
11097-69-1	AROCLOR 1254	ND	ug/L	0.1	0.1^		
11096-82-5	AROCLOR 1260	ND	ug/L	0.1	0.08		
12674-11-2	AROCLOR 1016	ND	ug/L	0.1	0.1		
8001-35-2	TOXAPHENE	ND	ug/L	1	0.5	3	

NOTES:
If a compound is delected > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.

MCI. (Maximum Contaminant Lavel) maximum permissible level of a contaminant in water established by EPA; Federal Action Levels are 0.015 mg/L for Lead and 1.3 mg/L for Copper. Sodium has a recommended first of 20 mg/L. A blank MCI. value indicates a level is not currently established.

Trigger Level: DOH Drinking Water Response level. Systems with compounds detected in excess of this level are required to take additional samples. Contact your regional DOH office.

ND (Not Detected): Indicates that the perameter was not detected above the State Reporting Limit (SRL).



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CARBAMATES IN DRINKING WATER

Client Name: Walla Walla Basin Watershed Council

810 South Main Street

Milton-Freewater, OR 97862

Project::

Field ID: L1

Sample Description: Locher 1

Sampled By: Unknown

Sample Date: 2/12/09

Source Type: Sampler Phone: Reference Number: 09-02151

Project: Locher Road Recharge Sites

Lab Number: 04470 Report Date: 2/24/09 Date Analyzed: 02/19/09

Date Extracted: 531_090219

Analyst: GO Peer Review: Analytical Method: 531.2

Carbamates

	COMPOUND	RESULTS	UNITS	PQL	MDL	MCL	COMMENT
EPA Re	gulated						
23135-22-0	OXYMAL	ND	υ ը/ ′∟	1.0	0.3	200	
1563-66-2	CARBOFURAN	ND	ug/L	1.0	0.2	40	
EPA Un	regulated						
1646-87-3	ALDICARB SULFOXIDE	.ND	ug/L	1.0	0.3		
1646-88-4	ALDICARB SULFONE	ND	ug/L	1.0	0.3		
16752-77-5	METHOMYL	ND	ug/L	1.0	0.3		
16655-82-6	3-HYDROXYCARBOFURAN	ND	ug/L	1.0	0.3		
116-06-3	ALDICARB	ND	ug/L	1.0	0.3		
63-25-2	CARBARYL	ND	ug/L	1.0	0.2		
State U	nregulated - Other						
114-26-1	PROPOXUR (BAYGON)	ND	пðуГ	1.0	0.4		
2032-65-7	METHIOCARB	ND	u g/L	1.0	0.3		

NO (Not Delected): indicates that the parameter was not detected above the State Reporting Limit (SRL).

If a compound is detected > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.

MCI. (Maximum Contaminant Level) maximum permissible level of a contaminant in water established by EPA; Federal Action Levels are 0.015 mg/L for Lead and 1.3 mg/L for Copper. Sedium has a recommended limit of 20 mg/L. A blank MCI. (value indicates a level is not currently established.

Trigger Level: DOH Drinking Water Resiponse level. Systems with compounds detected it excess of this level are required to take additional samples. Contact your regional DOH office.



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CARBAMATES IN DRINKING WATER

Client Name: Walla Walla Basin Watershed Council

810 South Main Street

Milton-Freewater, OR 97862

Project::

Field ID: L2

Sample Description: Locher 2

Sampled By: Unknown Sample Date: 2/12/09

Source Type: Sampler Phone: Reference Number: 09-02151

Project: Locher Road Recharge Sites

Lab Number: 04471 Report Date: 2/24/09 Date Analyzed: 02/19/09 Date Extracted: 531 090219

Analyst: CO

Peer Review: Analytical Method: 531.2

Carbamates

CAS	COMPOUND	RESULTS	UNITS	PQL	- MDL	MCL	COMMENT
EPA Re	gulated		:				
23135-22-0	OXYMAL	ND	ug/L	1.0	0.3	200	
1563-66-2	CARBOFURAN	ND	ug/L	1.0	0.2	40	
EPA Ur	regulated						
1646-87-3	ALDICARB SULFOXIDE	ND	ug/L	1.0	0.3		
1646-88-4	ALDICARB SULFONE	ND	ug/L	1.0	0.3		
16752-77-5	METHOMYI.	ND	ug/L	1.0	0.3		
16655-82-6	3-HYDROXYCARBOFURAN	ND	ug/L	1.0	0.3		
116-06-3	ALDICARB	ND	ug/L	1.0	0.3		
63-25-2	CARBARYL	ND	ug/L	1,0	0.2		•
State U	nregulated - Other						
114-26-1	PROPOXUR (BAYGON)	ND	ug/L	1.0	0.4		
2032-65-7	METHIOCARB	ND	ug/L	1.0	0.3		

NOTES:

NOT [ES]

If a compound is detacted > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.

MCL (Maximum Conteminant Level) maximum permissible level of a conteminant in water established by EPA; Federal Action Levels are 0.015 mg/L for Lead and 1.3 mg/L for Copper. Sodium has a recommended limit of 20 mg/L. A blank MCL value indicates a level is not currently established.

Tripger Level: DOH Orthogy Water Response level. Systems with compounds detected in excess of this level are required to take additional samples. Contact your regional DOH office.

ND (Not Detected): ladicates that the parameter was not detected above the State Reporting Limit (SRL).



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Page 1 of 1

CARBAMATES IN DRINKING WATER

Client Name: Walla Walla Basin Watershed Council

810 South Main Street

Milton-Freewater, OR 97862

Project::

Field ID: L3

Sample Description: Locher 3

Sampled By: Unknown

Sample Date: 2/12/09

Source Type:

Sampler Phone:

Reference Number: 09-02151

Project: Locher Road Recharge Sites

Lab Number: 04472

Report Date: 2/24/09

Date Analyzed: 02/19/09 Date Extracted: 531_090219

Analyst: CO

Peer Review: Analytical Method: 7531.2

Carbamates

CAS	COMPOUND	RESULTS	UNITS	PQL -	MIDL	MCL	COMMENT
EPA Re	gulated						
23135-22-0	OXYMAL	ND	ug/L	1.0	0.3	200	
1563-66-2	CARBOFURAN	ND	ug/L	1.0	0.2	40	
EPA Ur	regulated						
1646-87-3	ALDICARB SULFOXIDE	ND	ug/L	1.0	0.3		
1646-88-4	ALDICARB SULFONE	ND	ug/L	1.0	0.3		
16752-77-5	METHOMYL	ND	ug/L	1.0	0.3		
16655-82-6	3-HYDROXYCARBOFURAN	ND	ug/L	1.0	0.3		
116-06-3	ALDICARB	ND	ug/L	1.0	0.3		
63-25-2	CARBARYL	ND	ug/L	1.0	0.2		
State U	nregulated - Other						
114-26-1	PROPOXUR (BAYGON)	ND	ug/L	1.0	0.4		•
2032-65-7	METHIOCARB	ND	u g/L	1.0	0.3		•

NOTES:

NOTES:
If a compound is detected > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.

MCL (Maximum Contaminant Level) maximum permissible level of a contaminant in water catabisted by EPA; Federal Action Levels are 0.015 mg/l, for Lead and 1.3 mg/l, for Copper. 5 blank MCL value in indicates a level is not currently established.

Trigger Level: DOH Driving Water Response level. Systems with compounds detected in excess of this level are required to take additional samples. Contact your regional DOH office.

NO (Not Detected): Indicates that the parameter was not detected above the State Reporting Limit (SRL). erel, SRL, specified increased monitoring frequencies may accur per DOH. Is level of a contaminant in water catabilished by EPA; Federal Action Levels are 9.015 mg/L for Lead and 1.3 mg/L for Capper. Sediam has a recommended limit of 20 mg/L. A







QUALITY CONTROL REPORT SURROGATE REPORT

Reference Number: 09-02151 Report Date: 03/05/09

Lab No	Analyte	Result Qualifier	Units	Method	Limit
508_090217 4470 525_090217	TETRACHLORO-M-XYLENE (SURR)	116	%	508.1	Acceptance Limits 70%-130%
4470	1,3-DIMETHYL-2-NITROBENZENE (Surr) PYRENE-D10 (Surr) PERYLENE-D12 (Surr) TRIPHENYLPHOSPHATE (Surr)	98 92 105 93	% % %	525.2	Acceptance Range is 70% to 130% Acceptance Range is 70% to 130% Acceptance Range is 70% to 130% Acceptance Range is 70% to 130%
515.4_090219 4470	2,4 - DCAA (SURR)	87	*	515.4	Acceptance Range is 70 - 130%
531_090219 4470	BDMC (SURR)	103	%	531.2	
508_090217 4471 525_090217	TETRACHLORO-M-XYLENE (SURR)	82	%	508.1	Acceptance Limits 70%-130%
4471	1,3-DIMETHYL-2-NITROBENZENE (Surr) PYRENE-D10 (Surr) PERYLENE-D12 (Surr) TRIPHENYLPHOSPHATE (Surr)	96 91 105 89	% % %	525.2	Acceptance Range is 70% to 130% Acceptance Range is 70% to 130% Acceptance Range is 70% to 130% Acceptance Range is 70% to 130%
515.4_090219 4471	2,4 - DCAA (SURR)	78	%	515.4	Acceptance Range is 70 - 130%
531_090219 4471	BDMC (SURR)	111	%	531.2	
508_090217 4472 525_090217	TETRACHLORO-M-XYLENE (SURR)	80	% .	508.1	Acceptance Limits 70%-130%
4472	1,3-DIMETHYL-2-NITROBENZENE (Sum) PYRENE-D10 (Sum) PERYLENE-D12 (Sum) TRIPHENYLPHOSPHATE (Sum)	96 91 104 99	% % % %	525.2	Acceptance Range is 70% to 130% Acceptance Range is 70% to 130% Acceptance Range is 70% to 130% Acceptance Range is 70% to 130%
515.4_090219 4472	2,4 - DCAA (SURR)	84	%	515.4	Acceptance Range is 70 - 130%
531_090219 4472	BDMC (SURR)	121	%	531.2	

^{*}Notation:



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SAMPLE INDEPENDENT **QUALITY CONTROL REPORT**

Laboratory Fortified Blank

Reference Number: 09-02151

			True			%		QC	
Batch	Analyte	Result	Value	Units	Method	Recove	ry Limits	Qualifier Type*	Comment
200.7-090216A	HARDNESS	70	69.5	mg/L	200.7	101	80-120	LFB	
508_090217	AROCLOR 1260	1.5	2	ug/L	508.1	75	60-140	LFB	
	TETRACHLORO-M-XYLENE (SURR)	74		%	508.1		70-130		
515.4_090219	2,4 - D	7.8	10	ug/L	515.4	78	70-130	LFB	
	2,4 - DCAA (SURR)	89		%	515.4		70-130		
	2,4 DB	33.4	40	ug/L	515.4	84	70-130		
	2,4,5 - TP (SILVEX)	4.3	5	ug/L	515.4	86	70-130		
	2,4,5 T	4.4	5	ug/L	515.4	88	70-130		
	ACIFLUORFEN	4.2	5	ug/L	515.4	84	70-130		
	BENTAZON	9	10	ug/L	515.4	90	70-130		
	DALAPON	52.3	65	u g/L	515.4	80	70-130		
	DCPA (ACID METABOLITES)	3.5	5	ug/L	515.4	70	70-130		
	DICAMBA	4.3	5	ug/L	515.4	86	70-130		
	DICHLORPROP	11.7	15	ug/L	515.4	78	70-130		
	DINOSEB	8.9	10	ug/L	515.4	89	70-130		
	PENTACHLOROPHENOL	4.6	5	ug/L	515.4	92	70-130		
	PICLORAM	4.5	5	ug/L	515.4	90	70-130		•
515.4_090219	2,4 - D	4	5	ug/L	515.4	80	70-130	LFB	
	2,4 - DCAA (SURR)	81		%	515.4		70-130		
	2,4 DB	18.2	20	ug/L	515.4	91	70-130		
	2,4,5 - TP (SILVEX)	2.4	2.5	ug/L	515.4	96	70-130		
	2,4,5 T	2.4	2.5	ug/L	515.4	96	70-130		
	ACIFLUORFEN	2.4	2.5	ug/L	515.4	96	70-130		
	BENTAZON	4.8	5	ug/L	515.4	96	70-130		
	DALAPON	28.7	32.5	ug/L	515.4	88	70-130		•
	DCPA (ACID METABOLITES)	2.6	2.5	ug/L	515.4	104	70-130		
	DICAMBA	2.5	2.5	ug/L	515.4	100	70-130		
	DICHLORPROP	6.35	7.2	ug/L	515.4	88	70-130		
	DINOSEB	4.9	5	ug/L	515.4	98	70-130		
	PENTACHLOROPHENOL	2.6	2.5	ug/L	515.4	104	70-130		
	PICLORAM	2.7	2.5	ug/L	515.4	108	70-130		

[%] Recovery = (Result of Analysis)/(True Value) * 100

NA = Indicates % Recovery could not be calcutated.

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MB or LRB. Melhod Blank or Laboratory Reagant Blank, an adjust of reagent matrix is analyzed exactly like a sample, and its purpose is to determine if there is background contamination.



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SAMPLE INDEPENDENT QUALITY CONTROL REPORT

Laboratory Fortified Blank

Reference Number: 09-02151 Report Date: 03/05/09

			True			%		QC	
Batch	Analyte	Result	Value	Units	Method	Recovery	Limits	QualiflerType*	Comment
525_090217	1,3-DIMETHYL-2-NITROBENZENE (Sum)	97		=1	202.0		70.400	1.50	
020_00021.	4,4-DDD	1.05	1	% ***	525.2 5 2 5.2	105	70-130 70-130		
	4,4-DDE	0.98	1	ug/L	525.2 525.2	98			
	4,4-DDT	1.02	1	ug/L		102	70-130		
	ACENAPHTHYLENE	10.1	† †	ug/L	525.2 525.2		70-130		
	ALACHLOR	2.12	2	ug/L		1,010	70-130		
	ALDRIN			ug/L	525.2	106	70-130		
	ANTHRACENE	0.97	1	ug/L	525.2	97	70-130		
	ATRAZINE	0.98	1	ug/L	525.2	98	70-130		
		2.2	2	ug/L	525.2	110	70-130		,
	BENZ(A)ANTHRACENE	1.05	1	ug/L	525.2	105	70-130		
	BENZO(A)PYRENE	1.15	1	ug/L	525.2	115	70-130		
	BENZO(B)FLUORANTHENE	1.27	1.	ug/L	525.2	127	70-130		
	BENZO(G,H,I)PERYLENE	0.96	1	ug/L	525.2	96	70-130		
	BENZO(K)FLUORANTHENE	1.04	1	ug/L	525.2	104	70-130		
	BENZYL BUTYL PHTHALATE	1.13	1	ug/L	525.2	113	70-130		
	BROMACIL.	1.07	1	ug/L	525.2	107	70-130		
	BUTACHLOR	1.2	1	ug/L	5 2 5.2	120	70-130		
	CHLORDANE, TECHNICAL	0.77	1	ug/L	525.2	77	70-130		
	CHRYSENE	1.04	1	ug/L	525.2	104	70-130		
	CYANAZINE	2.46	2	ug/L	525.2	123	70-130		
	DI(ETHYLHEXYL)-ADIPATE	1.2	1	ug/L	525.2	120	70-130		
	DI(ETHYLHEXYL)-PHTHALATE	1.38	1	ug/L	525.2	138	70-130	B5	
	DIAZINON	3.19	3	ug/L	525.2	106	70-130		
	DIBENZO(A,H)ANTHRACENE	1.09	1	ug/L	525.2	109	70-130		
	DIELDRIN	1.17	1	ug/L	525.2	117	70-130		
	DIETHYL PHTHALATE	1.08	1	ug/L	525.2	108	70-130		
	DIMETHYL PHTHALATE	1	1	ug/L	525.2	100	70-130		
	DI-N-BUTYL PHTHALATE	1.08	1	ug/L	525.2	108	70-130		
	ENDRIN	1.02	1	սց/Լ	525.2	102	70-130		
	EPTC	1.01	1	ug/L	525.2	101	70-130		
	FLUORENE	1.04	1	ug/L	525.2	104	70-130		
	HEPTACHLOR	1.02	1	ug/L	525.2	102	70-130		
	HEPTACHLOR EPOXIDE	1.08	1	ug/L	525.2	108	70-130		
	HEXACHLOROBENZENE	0.89	1	ug/L	525.2	89	70-130		
	HEXACHLOROCYCLO-PENTADIENE	0.96	1	ug/L	525.2	96	70-130		

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SAMPLE INDEPENDENT QUALITY CONTROL REPORT

Laboratory Fortified Blank

Reference Number: 09-02151

			True			%		QC	
Batch	Analyte	Result	Value	Units	Melhod	Recovery	Limits	Qualifier Type*	Comment
525_090217	INDENO(1,2,3-CD)PYRENE	1.19	1	ug/L	525.2	119	70-130	LFB	
	LINDANE (BHC - GAMMA)	0.97	1	ug/L	525.2	97	70-130		
	MALATHION	2.62	3	ug/L	525.2	87	70-130		
	METHOXYCHLOR	1.3	1	ug/L	525.2	130	70-130		
•	METOLACHLOR	1.13	1	ug/L	525.2	113	70-130		
	METRIBUZIN	0.86	1	ug/L	525.2	86	70-130		
	PARATHION	2.17	2	⊔g/L	525.2	109	70-130		
	PENTACHLOROPHENOL	3.87	4	ug/L	525.2	97	70-130		
	PERYLENE-D12 (Surr)	106		%	525.2		70-130		
	PHENANTHRENE	0.98	1	ug/L	525.2	98	70-130		
	PROPACHLOR	1.09	1	ug/L	525.2	109	70-130		
	PYRENE	0.97	1	ug/L	525.2	97	70-130		
	PYRENE-D10 (Surr)	94		%	525.2		70-130		
-	SIMAZINE	1.08	1	ug/L	525.2	108	70-130		
	TERBACIL	1.32	1	ug/L	525.2	132	70-130	HQ	
	TRIFLURALIN	1.02	1	ug/L	525.2	102	70-130		
	TRIPHENYLPHOSPHATE (Surr)	93		%	525.2		70-130		•
531_090219	3-HYDROXYCARBOFURAN	5.7	5	ug/L	531.2	114	70-130	LFR	
	ALDICARB	5.4	5	ug/L	531.2	108	70-130		
	ALDICARB SULFONE	5.6	5	ug/L	531.2	112	70-130		
	ALDICARB SULFOXIDE	5.5	5	ug/L	531.2	110	70-130		
	BDMC (SURR)	94		%	531.2		70-130		
	CARBARYL	5.2	5	ug/L	531.2	104	70-130		
	CARBOFURAN	5.3	5 ′	ug/L	531.2	106	70-130		
	METHIOCARB	5.3	5	ug/L	531.2	106	70-130		
	METHOMYL	5. 1	5	ug/L	531.2	102	70-130		
	OXYMAL	5	5	ug/L	531.2	100	70-130	-	
	PROPOXUR (BAYGON)	5.5	5	ug/L	531.2	110	70-130		
531_090219	3-HYDROXYCARBOFURAN	22.7	20	ug/L	531.2	114	70-130	LFB	
	ALDICARB	19.7	20	ug/L	531.2	99	70-130		
	ALDICARB SULFONE	21.9	20	ug/L	531.2	110	70-130		
	ALDICARB SULFOXIDE	22.4	20	ug/L	531.2	112	70-130		
	BDMC (SURR)	107		%	531.2		70-130	-	
	CARBARYL	19.7	20	սց /L	531.2	99	70-130		

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SAMPLE INDEPENDENT QUALITY CONTROL REPORT

Laboratory Fortified Blank

Reference Number: 09-02151

			eunT			%		QC	
Batch	Analyte	Result	Value	Units	Method	Recovery	Limits	Qualifier Type*	Comment
531_090219	CARBOFURAN	21.2	20	ug/L	531.2	106	70-130	LFB	
	METHIOCARB	20.4	20	ug/L	531.2	102	70-130		
	METHOMYL	22.7	20	ug/L	531.2	114	70-130		
	OXYMAL	22.5	20	ug/L	531.2	113	70-130		
	PROPOXUR (BAYGON)	21	20	ug/L	531.2	105	70-130		
COD_090216	CHEMICAL OXYGEN DEMAND	47	50	mg/L	SM5220 D	94	80-120	LFB	
OPHOS-090213	ORTHO-PHOSPHATE	1.00	1.00	mg/L	SM4500-P F	100	70-130	LFB	
TDS_090218	TOTAL DISSOLVED SOLIDS	494	500	mg/L	SM2540 C	99	80-120	LF8	
TDS_090218	TOTAL DISSOLVED SOLIDS	480	500	mg/L	SM2540 C	96	80-120	LFB	

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SAMPLE INDEPENDENT QUALITY CONTROL REPORT

Low Level Laboratory Fortified Blank

Reference Number: 09-02151

			True			%		QC	,
Batch	Analyte	Result	Value	Units	Method	Recovery	Limits	Qualifier Type"	Comment
515.4_090219	2,4 - D	0.32	0.5	ug/L	515.4	64	50-150	LFBD	
	2,4 - DCAA (SURR)	87		%	515.4		70-130		
	2,4 DB	1.1	2	ug/L	515.4	55	50-150		
	2,4,5 - TP (SILVEX)	0.25	0.25	ug/L	515.4	100	50-150		
	2,4,5 T	0.2	0.25	ug/L	515.4	80	50-150		
	ACIFLUORFEN	0.26	0.25	ug/L	515.4	104	50-150		
	BENTAZON	0.52	0.5	ug/L	515.4	104	50-150		
	DALAPON	3.9	3,25	ug/L	515.4	120	50-150		
	DCPA (ACID METABOLITES)	0.2	0.25	ug/L	515.4	80	50-150		
	DICAMBA	0.26	0.25	ug/L	515.4	104	50-150		
	DICHLORPROP	0.7	0.75	ug/L	515.4	93	50-150		
	DINOSEB	0.48	0.5	ug/L	515.4	96	50-150		
	PENTACHLOROPHENOL	0.23	0.25	ug/L	515.4	92	50-150		
	PICLORAM	0.22	0.25	ug/L	515.4	88	50-150		
531_090219	3-HYDROXYCARBOFURAN	1.1	1	ug/L	531.2	110	50-150	LFBD	
	ALDICARB	0.78	1	ug/L	531.2	78	50-150	•	
	ALDICARB SULFONE	1	1	ug/L	531.2	100	50-150		
	ALDICARB SULFOXIDE	0.71	1	ug/L	531.2	71	50-150		
	BDMC (SURR)	120	-	%	531.2	• •	50-150		
	CARBARYL	1.05	1	ug/L	531.2	105	50-150		
	CARBOFURAN	1.3	1	ug/L	531.2	130	50-150		
	METHIOCARB	1.3	1	ug/i.	531.2	130	50-150		
	METHOMYL	0.8	1	ug/L	531.2	80	50-150		
	OXYMAL	D.95	1	ug/L	531.2	95	50-150		
	PROPOXUR (BAYGON)	1.1	1	ug/L	531.2	110	50-150		

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SAMPLE INDEPENDENT **QUALITY CONTROL REPORT**

Laboratory Reagent Blank

Reference Number: 09-02151

			True			%		QC	
Batch	Analyle	Result	Value	Units	Method	Recovery	Limits	Qualifier Type*	Comment
200.7-090216A	HARDNESS	ND		mg/L	200.7		10.0000	C LRB	
COD_090216	CHEMICAL OXYGEN DEMAND	ND		mg/L	SM5220 D		4.00000	LRB	
D090304A	BROMATE	ND		mg/L	300.1		0.00500	LRB	
1090213A	CHLORIDE NITRATE-N	ND ND		mg/L mg/L	300.0 300.0		9.10000 9.10000		
OPHOS-090213	ORTHO-PHOSPHATE	ND		mg/L	SM4500-P F		0.10000	LRB	

^{*}Notation:

[%] Recovery = (Result of Analysis)/(True Value) * 100

NA = Indicates % Recovery could not be calculated.

QCS: Quality Control Sample, a solution containing known concentrations of method analytes which is used to fortify an aliquot of reagent matrix. The QCS is obtained from an external source and is used to check lab performance.

LFB: Laboratory Fortified Blank, an aliquot of reagent matrix to which known quantities of method analytes are added in the lab. The LFB is analyzed exactly like a sample, and its purpose is to determine whether method performance is within accepted control limits.

MB or LRB: Method Blank or Laboratory Reagent Blank, an aliquot of reagent matrix is analyzed exactly like a sample, and its purpose is to determine if there is background contamination.



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SAMPLE INDEPENDENT QUALITY CONTROL REPORT

Method Blank

Reference Number: 09-02151

			True		·	%	QC	
Batch	Analyte	Result	Value	Units	Method	Recovery Limits	Qualifier Type*	Comment
200.7-090216A	HARDNESS	ND		mg/L	200.7	0.8200	The state of the s	
						•		
508_090217	AROCLOR 1016	ND		ug/L	508.1	0.0200) MB	
	AROCLOR 1221	ND		ug/L	508.1	0.1200		
	AROCLOR 1232	ND		ug/L	50B.1	0.0200		
	AROCLOR 1242	ND		ug/L	508.1	0.0200		
-	AROCLOR 1248	ND		ug/L	508.1	0.0200		
	AROCLOR 1254	ND		ug/L	506.1	0.02000		
	AROCLOR 1260	ND		ug/L	508.1	0.02000		
	TETRACHLORO-M-XYLENE (SURR)	82		~ 	508.1	0.00000	•	
	•							
515.4_090219	2,4 - D	NĐ		υg/L	515.4	0.10000) MB	
	2,4 - DCAA (SURR)	98		%	515.4	3.150gL	1410	
	2,4 DB	NĐ		ug/L	515.4	0.60000)	
	2,4,5 - TP (SILVEX)	ND	-	ug/L	515.4	0.08000		
	2,4,5 T	NĐ		ug/L	515.4	0.08000		
	ACIFLUORFEN	ND		ug/L	515.4	0.08000		
	BENTAZON	ND		ug/L	515.4	0.10000		
	DALAPON	ND		ug/L	515.4	1.00000		
	DCPA (ACID METABOLITES)	ND		ug/L	515.4	0.08000		
	DICAMBA	ND		ug/L	515.4	0.08000		
	DICHLORPROP	ND		ug/L	515.4	0.20000		
	DINOSEB	ND		ug/L	515.4	0.10000		
	PENTACHLOROPHENOL	ND		ug/L	515.4	0.08000		
•	PICLORAM	ND		ug/L	515.4	0.08000		
25_090217	1,3-DIMETHYL-2-NITROBENZENE (Surr)	99		%	525.2		MB	
	4,4-DDD	ND		ug/L	525.2	0.05000		
	4,4-DDE	ND		ug/L	525.2	0.05000		
	4,4-DOT	ND		ug/L	525.2	0.05000		
	ACENAPHTHENE	ND		ug/L	525.2	0.05000		
	ALACHLOR	ND		ug/L	525.2	0.02000		
	ALDRIN	ND		ug/L	525.2	0.05000		

[&]quot;Notation:

[%] Recovery = (Result of Analysis)/(True Value) * 100

NA = indicates % Recovery could not be calculated.

QCS: Quality Control Sample, a solution containing known concentrations of method analytes which is used to fortify an aliquot of reagent matrix. The QCS is obtained from an external source and is used to check lab performance.

LFB: Laboratory Fortified Blank, an aliquot of reagent matrix to which known quantities of method analytes are added in the lab. The LFB is analyzed exactly like a sample, and its purpose is to determine whether method performance is within accepted control limits.

MB or LRB: Method Blank or Laboratory Reagent Blank, an aliquot of reagent matrix is analyzed exactly like a sample, and its purpose is to determine if there is background contamination.



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SAMPLE INDEPENDENT QUALITY CONTROL REPORT

Method Blank

Reference Number: 09-02151

Batch				True			%		QC	
ATRAZINE ND UJL 525.2 0.02000 BENZO(A)-PYRENE ND UJL 525.2 0.02000 BENZO(B)-PYRENE ND UJL 525.2 0.02000 BENZO(B)-PYRENE ND UJL 525.2 0.02000 BENZO(B)-FLUORANTHENE ND UJL 525.2 0.05000 CHLORDANE, TECHNICAL ND UJL 525.2 0.05000 CHETHYLHEXYL)-ADIPATE ND UJL 525.2 0.05000 DKETHYLHEXYL)-PHTHALATE 0.39 UJL 525.2 0.05000 DKETHYLHEXYL-PHTHALATE 0.39 UJL 525.2 0.05000 DIBENZO(A,H)ANTHRACENE ND UJL 525.2 0.05000 DIBENZO(A,H)ANTHRACENE ND UJL 525.2 0.05000 DIBENZO(A,H)ANTHRACENE ND UJL 525.2 0.05000 DIEDRIN ND UJL 525.2 0.05000 DIEDRIYL PHTHALATE ND UJL 525.2 0.05000 DI-HUTYL PHTHALATE ND UJL 525.2 0.05000 DI-HUTYL PHTHALATE ND UJL 525.2 0.05000 DI-HUTYL PHTHALATE ND UJL 525.2 0.05000 EPTC ND UJL 525.2 0.05000 PHENDRIN ND UJL 525.2 0.05000 PHENDRIC -CAMMA) ND UJL 525.2 0.05000	Batch	Analyte	Result	Value	Units	Method	Recovery L	imits.	Qualifier Type*	Comment
BENZ(A)ANTHRACENE ND	525_090217	ANTHRACENE	ND		ug/L	525.2	C	0.05000	MB	
BENZO(A)PYRENE ND		ATRAZINE	ND		ug/L	525.2	C	0.02000		
BENZO(B)FLUCRANTHENE ND		BENZ(A)ANTHRACENE	ND		ug/L	525.2	C	0.02000		
BENZO(G.H.I)PERYLENE ND		BENZO(A)PYRENE	ND		ug/L	525.2	C	0.02000		
BENZO(K)FLUORANTHENE ND ug/L 525.2 0.05000 BENZYL BUTYL PHTHALATE ND ug/L 525.2 0.05000 BROMACIL ND ug/L 525.2 0.05000 BUTACHLOR ND ug/L 525.2 0.05000 CHLORDANE, TECHNICAL ND ug/L 525.2 0.05000 CHLORDANE, TECHNICAL ND ug/L 525.2 0.05000 CHRYSENE ND ug/L 525.2 0.05000 CKYANAZINE ND ug/L 525.2 0.05000 DKETHYLHEXYL)-PHTHALATE 0.39 ug/L 525.2 0.05000 DIETHYLHEXYL)-PHTHALATE 0.39 ug/L 525.2 0.05000 DIBENZO(A,H)ANTHRACENE ND ug/L 525.2 0.05000 DIETHYL PHTHALATE ND ug/L 525.2 0.05000 DIETHYL PHTHALATE ND ug/L 525.2 0.05000 DIMETHYL PHTHALATE ND ug/L 525.2 0.05000 DIMETHYL PHTHALATE ND ug/L 525.2 0.05000 DIMETHYL PHTHALATE ND ug/L 525.2 0.05000 DI-BUTYL PHTHALATE ND ug/L 525.2 0.05000 ENDRIN ND ug/L 525.2 0.05000 ENDRIN ND ug/L 525.2 0.05000 ENDRIN ND ug/L 525.2 0.05000 FLUORENE ND ug/L 525.2 0.05000 FLUORENE ND ug/L 525.2 0.05000 HEPTACHLOR ND ug/L 525.2 0.05000 HEPTACHLOR ND ug/L 525.2 0.05000 HEPTACHLOR ND ug/L 525.2 0.05000 HEXACHLOROGENZENE ND ug/L 525.2 0.05000 ILIDANE (BHC - GAMMA) ND ug/L 525.2 0.02000 ILIDANE (BHC - GAMMA) ND ug/L 525.2 0.02000 IMETHUSICAL OR ND ug/L 525.2 0.05000 ILIDANE (BHC - GAMMA) ND ug/L 525.2 0.05000 ILIDANE (BHC - GAMMA) ND ug/L 525.2 0.05000 IMETRIBUZIN ND ug/L 525.2 0.05000 IMETRIBUZIN ND ug/L 525.2 0.05000 IMETRIBUZIN ND ug/L 525.2 0.05000		BENZO(B)FLUORANTHENE	ND		ug/L	525.2	c). 05 000		
BENZYL BUTYL PHTHALATE ND		BENZO(G,H,I)PERYLENE	ND		ug/L	525.2	ŧ	0.05000		
BROMACIL BUTACHLOR BUTACHLOR ND Ug/L 925.2 0.05000 CHLORDANE, TECHNICAL ND Ug/L 525.2 0.05000 CHRYSENE ND Ug/L 525.2 0.05000 CYANAZINE ND Ug/L 525.2 0.05000 DIETHYLHEXYL}-ADIPATE ND Ug/L 525.2 0.05000 DIETHYLHEXYL}-PHTHALATE 0.39 Ug/L 525.2 0.05000 DIEDRIN ND Ug/L 525.2 0.05000 DIEDRIN ND Ug/L 525.2 0.05000 DIETHYLPHTHALATE ND Ug/L 525.2 0.05000 DIETHYLPHTHALATE ND Ug/L 525.2 0.05000 DIETHYLPHTHALATE ND Ug/L 525.2 0.05000 DIMETHYLPHTHALATE ND		BENZO(K)FLUORANTHENE	ND		ug/L	525.2	C	0.05000		
BUTACHLOR ND Ug/L 525.2 0.10006 CHLORDANE, TECHNICAL ND Ug/L 525.2 0.05000 CHRYSENE ND Ug/L 525.2 0.05000 CYANAZINE ND Ug/L 525.2 0.05000 D(ETHYLHEXYL)-ADIPATE ND Ug/L 525.2 0.05000 D(ETHYLHEXYL)-PHTHALATE 0.39 Ug/L 525.2 0.05000 DIELDRIN ND Ug/L 525.2 0.05000 DIELDRIN ND Ug/L 525.2 0.05000 DIETHYL PHTHALATE ND Ug/L 525.2 0.05000 DIETHYL PHTHALATE ND Ug/L 525.2 0.05000 DIETHYL PHTHALATE ND Ug/L 525.2 0.05000 DIH-BUTYL PHTHALATE ND Ug/L 525.2 0.05000 EPTC ND Ug/L 525.2 0.05000 HEPTACHLOR POXIDE ND Ug/L 525.2 0.05000 HEPTACHLOR SENZENE ND Ug/L 525.2 0.05000 HEXACHLOROBENZENE ND Ug/L 525.2 0.05000 HEXACHLOROGYCLO-PENTADIENE ND Ug/L 525.2 0.05000 HEXACHLOROGYCLO-PENTADIENE ND Ug/L 525.2 0.05000 MALATHION ND Ug/L 525.2 0.05000 METHOLACHLOR ND Ug/L 525.2 0.05000		BENZYL BUTYL PHTHALATE	ND		ug/L	525.2	C	0.60000		•
CHLORDANE, TECHNICAL ND ug/L 525.2 0.02000 CHRYSENE ND ug/L 525.2 0.05000 CYANAZINE ND ug/L 525.2 0.05000 D(ETHYLHEXYL)-ADIPATE ND ug/L 525.2 0.02000 D(ETHYLHEXYL)-PHTHALATE 0.39 ug/L 525.2 0.06000 DIAZINON ND ug/L 525.2 0.05000 DIBENZO(A,H)ANTHRACENE ND ug/L 525.2 0.05000 DIBENZO(A,H)ANTHRACENE ND ug/L 525.2 0.05000 DIELDRIN ND ug/L 525.2 0.05000 DIELTHYL PHTHALATE ND ug/L 525.2 0.05000 DIELTHYL PHTHALATE ND ug/L 525.2 0.05000 DIMETHYL PHTHALATE ND ug/L 525.2 0.05000 DIA-BUTYL PHTHALATE ND ug/L 525.2 0.05000 DIA-BUTYL PHTHALATE ND ug/L 525.2 0.05000 DIA-BUTYL PHTHALATE ND ug/L 525.2 0.05000 ENDRIN ND ug/L 525.2 0.05000 EPTC ND ug/L 525.2 0.05000 PLUORANTHENE ND ug/L 525.2 0.05000 PLUORANTHENE ND ug/L 525.2 0.05000 HEPTACHLOR ND ug/L 525.2 0.02000 HEXACHLOROGYCLO-PENTADIENE ND ug/L 525.2 0.02000 INDENO(1,2,3-CD)PYRENE ND ug/L 525.2 0.02000 INDENO(1,2,3-CD)PYRENE ND ug/L 525.2 0.02000 METHOLYCHLOR ND ug/L 525.2 0.02000 MALATHION ND ug/L 525.2 0.02000 METHOLYCHLOR ND ug/L 525.2 0.02000		BROMACIL	ND		ug/L	525.2	c	0.05000		
CHRYSENE ND ug/L 525.2 0.05000 CYANAZINE ND ug/L 525.2 0.05000 D(ETHYLHEXYL)-ADIPATE ND ug/L 525.2 0.05000 D(ETHYLHEXYL)-PHTHALATE 0.39 ug/L 525.2 0.05000 D(EDIX) DIAZINON ND ug/L 525.2 0.05000 DIBENZO(A,H)ANTHRACENE ND ug/L 525.2 0.05000 DIBENZO(A,H)ANTHRACENE ND ug/L 525.2 0.05000 DIBENZO(A,H)ANTHRACENE ND ug/L 525.2 0.05000 DIELDRIN ND ug/L 525.2 0.05000 DIELDRIN ND ug/L 525.2 0.05000 DIMETHYL PHTHALATE ND ug/L 525.2 0.05000 DI-N-BUTYL PHTHALATE ND ug/L 525.2 0.05000 DI-N-BUTYL PHTHALATE ND ug/L 525.2 0.05000 ENDRIN ND ug/L 525.2 0.05000 EPTC ND ug/L 525.2 0.05000 FLUORANTHENE ND ug/L 525.2 0.05000 FLUORENE ND ug/L 525.2 0.05000 FLUORENE ND ug/L 525.2 0.05000 HEPTACHLOR POXIDE ND ug/L 525.2 0.05000 HEPTACHLOR POXIDE ND ug/L 525.2 0.05000 HEPTACHLOR POXIDE ND ug/L 525.2 0.02000 HEXCACHLOROGENZENE ND ug/L 525.2 0.02000 HEXACHLOROGENZENE ND ug/L 525.2 0.02000 INDENO(1,2,3-CD)PYRENE ND ug/L 525.2 0.02000 INDENO(1,2,3-CD)PYRENE ND ug/L 525.2 0.05000 METHORYCHLOR ND ug/L 525.2 0.05000		BUTACHLOR	ND		սց/Լ	525.2	C	0.10000		
CYANAZINE ND ug/L 525.2 0.05000 DKETHYLHEXYL)-ADIPATE ND ug/L 525.2 0.02000 DIETHYLHEXYL)-PHTHALATE 0.39 ug/L 525.2 0.05000 DIELDRIN ND ug/L 525.2 0.05000 DIELDRIN ND ug/L 525.2 0.05000 DIETHYL PHTHALATE ND ug/L 525.2 0.60000 DIMETHYL PHTHALATE ND ug/L 525.2 0.60000 DI-BUTYL PHTHALATE ND ug/L 525.2 0.60000 DI-N-BUTYL PHTHALATE ND ug/L 525.2 0.60000 DI-N-BUTYL PHTHALATE ND ug/L 525.2 0.60000 ENDRIN ND ug/L 525.2 0.00000 EPTC ND ug/L 525.2 0.02000 EPTC ND ug/L 525.2 0.05000 FLUGRANTHENE ND ug/L 525.2 0.05000 HEPTACHLOR ND ug/L </td <td></td> <td>CHLORDANE, TECHNICAL</td> <td>ND</td> <td></td> <td>ug/L</td> <td>525.2</td> <td>C</td> <td>0.02000</td> <td></td> <td></td>		CHLORDANE, TECHNICAL	ND		ug/L	525.2	C	0.02000		
DKETHYLHEXYL)-ADIPATE ND		CHRYSENE	ND		ug/L	525.2	C	0.05000		
DI(ETHYLHEXYL)-PHTHALATE		CYANAZINE	ND		ug/L	525.2	C	0.05000		
DIAZINON ND ug/L 525.2 0.05000 DIBENZO(A,H)ANTHRACENE ND ug/L 525.2 0.05000 DIELDRIN ND ug/L 525.2 0.05000 DIETHYL PHTHALATE ND ug/L 525.2 0.60000 DIMETHYL PHTHALATE ND ug/L 525.2 0.60000 DI-BUTYL PHTHALATE ND ug/L 525.2 0.60000 ENDRIN ND ug/L 525.2 0.02000 EPTC ND ug/L 525.2 0.07000 FLUORANTHENE ND ug/L 525.2 0.05000 FLUORENE ND ug/L 525.2 0.05000 HEPTACHLOR ND ug/L 525.2 0.02000 HEPTACHLOR EPOXIDE ND ug/L 525.2 0.02000 HEXACHLOROBENZENE ND ug/L 525.2 0.02000 HEXACHLOROGYCLO-PENTADIENE ND ug/L 525.2 0.02000 INDANIE (BHG - GAMMA) ND		DKETHYLHEXYL)-ADIPATE	ND		ug/L	525.2	C	0.02000		
DIBENZO(A,H)ANTHRACENE ND		DI(ETHYLHEXYL)-PHTHALATE	0.39		ug/L	525.2	C	0,00008.0		Lab BIG avg 0.2 ppb
DIELDRIN ND Ug/L 525.2 0.05000 DIETHYL PHTHALATE ND Ug/L 525.2 0.60000 DIMETHYL PHTHALATE ND Ug/L 525.2 0.60000 DI-N-BUTYL PHTHALATE ND Ug/L 525.2 0.50000 ENDRIN ND Ug/L 525.2 0.02000 ENDRIN ND Ug/L 525.2 0.07000 FLUORANTHENE ND Ug/L 525.2 0.07000 FLUORANTHENE ND Ug/L 525.2 0.05000 FLUORENE ND Ug/L 525.2 0.05000 HEPTACHLOR ND Ug/L 525.2 0.02000 HEPTACHLOR EPOXIDE ND Ug/L 525.2 0.02000 HEXACHLORGENZENE ND Ug/L 525.2 0.02000 HEXACHLORGOYCLO-PENTADIENE ND Ug/L 525.2 0.05000 INDANE (BHC - GAMMA) ND Ug/L 525.2 0.05000 METHOXYCHLOR ND <t< td=""><td></td><td>DIAZINON</td><td>ND</td><td></td><td>ug/L</td><td>525.2</td><td>C</td><td>0.05000</td><td></td><td></td></t<>		DIAZINON	ND		ug/L	525.2	C	0.05000		
DIETHYL PHTHALATE ND ug/L 525.2 0.60000 DIMETHYL PHTHALATE ND ug/L 525.2 0.60000 DI-N-BUTYL PHTHALATE ND ug/L 525.2 0.50000 ENDRIN ND ug/L 525.2 0.02000 EPTC ND ug/L 525.2 0.07000 FLUORANTHENE ND ug/L 525.2 0.05000 FLUORENE ND ug/L 525.2 0.05000 HEPTACHLOR ND ug/L 525.2 0.02000 HEPTACHLOR EPOXIDE ND ug/L 525.2 0.02000 HEXACHLOROGYCLO-PENTADIENE ND ug/L 525.2 0.02000 HEXACHLOROGYCLO-PENTADIENE ND ug/L 525.2 0.05000 INDENO(1,2,3-CD)PYRENE ND ug/L 525.2 0.05000 METHOXYCHLOR ND ug/L 525.2 0.05000 METHOXYCHLOR ND ug/L 525.2 0.02000 METOLACHLOR ND <td></td> <td>DIBENZO(A,H)ANTHRACENE</td> <td>ND</td> <td></td> <td>ug/L</td> <td>525.2</td> <td>C</td> <td>0.05000</td> <td></td> <td></td>		DIBENZO(A,H)ANTHRACENE	ND		ug/L	525.2	C	0.05000		
DIMETHYL PHTHALATE ND ug/L 525.2 0.60000 DIN-BUTYL PHTHALATE ND ug/L 525.2 0.60000 ENDRIN ND ug/L 525.2 0.02000 EPTC ND ug/L 525.2 0.07000 FLUGRANTHENE ND ug/L 525.2 0.05000 FLUGRENE ND ug/L 525.2 0.05000 HEPTACHLOR ND ug/L 525.2 0.02000 HEPTACHLOR EPOXIDE ND ug/L 525.2 0.02000 HEXACHLOROBENZENE ND ug/L 525.2 0.02000 HEXACHLOROCYCLO-PENTADIENE ND ug/L 525.2 0.05000 INDANE (BHC - GAMMA) ND ug/L 525.2 0.05000 MALATHION ND ug/L 525.2 0.05000 METHOXYCHLOR ND ug/L 525.2 0.05000 METOLACHLOR ND ug/L 525.2 0.05000 METRIBUZIN ND ug/L <td></td> <td>DIELDRIN</td> <td>ND</td> <td></td> <td>ug/L</td> <td>525.2</td> <td></td> <td>0.05000</td> <td></td> <td></td>		DIELDRIN	ND		ug/L	525.2		0.05000		
DI-N-BUTYL PHTHALATE ND ug/L 525.2 0.60000 ENDRIN ND ug/L 525.2 0.02000 EPTC ND ug/L 525.2 0.07000 FLUORANTHENE ND ug/L 525.2 0.05000 FLUORENE ND ug/L 525.2 0.05000 HEPTACHLOR ND ug/L 525.2 0.02000 HEPTACHLOR EPOXIDE ND ug/L 525.2 0.02000 HEXACHLOROBENZENE ND ug/L 525.2 0.02000 HEXACHLOROCYCLO-PENTADIENE ND ug/L 525.2 0.02000 INDENO(1,2,3-CD)PYRENE ND ug/L 525.2 0.05000 LINDANE (BHC - GAMMA) ND ug/L 525.2 0.05000 METHOXYCHLOR ND ug/L 525.2 0.02000 METOLACHLOR ND ug/L 525.2 0.02000 METRIBUZIN ND ug/L 525.2 0.05000		DIETHYL PHTHALATE	ND		ug/L	525.2	c	0.60000		
ENDRIN ND ug/L 525.2 0.02000 EPTC ND ug/L 525.2 0.07000 FLUORANTHENE ND ug/L 525.2 0.05000 FLUORENE ND ug/L 525.2 0.05000 HEPTACHLOR ND ug/L 525.2 0.02000 HEPTACHLOR EPOXIDE ND ug/L 525.2 0.02000 HEXACHLOROBENZENE ND ug/L 525.2 0.02000 HEXACHLOROCYCLO-PENTADIENE ND ug/L 525.2 0.02000 INDENO(1,2,3-CD)PYRENE ND ug/L 525.2 0.05000 LINDANE (BHC - GAMMA) ND ug/L 525.2 0.02000 MALATHION ND ug/L 525.2 0.05000 METHOXYCHLOR ND ug/L 525.2 0.05000 METHOXYCHLOR ND ug/L 525.2 0.02000 METOLACHLOR ND ug/L 525.2 0.05000 METOLACHLOR ND ug/L 525.2 0.05000 METOLACHLOR ND ug/L 525.2 0.05000		DIMETHYL PHTHALATE	NiD		ug/L	525.2	C	0.60000		•
EPTC ND ug/L 525.2 0.07000 FLUORANTHENE ND ug/L 525.2 0.05000 FLUORENE ND ug/L 525.2 0.05000 HEPTACHLOR ND ug/L 525.2 0.02000 HEPTACHLOR EPOXIDE ND ug/L 525.2 0.02000 HEXACHLOROBENZENE ND ug/L 525.2 0.02000 HEXACHLOROCYCLO-PENTADIENE ND ug/L 525.2 0.02000 INDENO(1,2,3-CD)PYRENE ND ug/L 525.2 0.05000 LINDANE (BHC - GAMMA) ND ug/L 525.2 0.05000 MALATHION ND ug/L 525.2 0.05000 METHOXYCHLOR ND ug/L 525.2 0.05000 METOLACHLOR ND ug/L 525.2 0.02000 METOLACHLOR ND ug/L 525.2 0.05000 METOLACHLOR ND ug/L 525.2 0.05000 METRIBUZIN ND ug/L 525.2 0.05000		DI-N-BUTYL PHTHALATE	ND		ug/L	525.2	C	0.60000		
FLUORANTHENE ND ug/L 525.2 0.05000 FLUORENE ND ug/L 525.2 0.05000 HEPTACHLOR ND ug/L 525.2 0.02000 HEPTACHLOR EPOXIDE ND ug/L 525.2 0.02000 HEXACHLOROBENZENE ND ug/L 525.2 0.02000 HEXACHLOROCYCLO-PENTADIENE ND ug/L 525.2 0.02000 INDENO(1,2,3-CD)PYRENE ND ug/L 525.2 0.05000 LINDANE (BHC - GAMMA) ND ug/L 525.2 0.05000 MALATHION ND ug/L 525.2 0.05000 METHOXYCHLOR ND ug/L 525.2 0.02000 METOLACHLOR ND ug/L 525.2 0.25000 METRIBUZIN ND ug/L 525.2 0.05000		ENDRIN	ND		ug/L	525.2	C	0.02000		
FLUORENE ND ug/L 525.2 0.05000 HEPTACHLOR ND ug/L 525.2 0.02000 HEPTACHLOR EPOXIDE ND ug/L 525.2 0.02000 HEXACHLOROGENZENE ND ug/L 525.2 0.02000 HEXACHLOROGYCLO-PENTADIENE ND ug/L 525.2 0.02000 INDENO(1,2,3-CD)PYRENE ND ug/L 525.2 0.05000 LINDANE (BHC - GAMMA) ND ug/L 525.2 0.05000 MALATHION ND ug/L 525.2 0.05000 METHOXYCHLOR ND ug/L 525.2 0.02000 METOLACHLOR ND ug/L 525.2 0.05000 METRIBUZIN ND ug/L 525.2 0.05000		EPTC	ND		ug/L	525.2	(0.07000		
HEPTACHLOR		FLUORANTHENE	ND		ug/L	525,2	C	0.05000		
HEPTACHLOR EPOXIDE ND ug/L 525.2 0.02000 HEXACHLOROBENZENE ND ug/L 525.2 0.02000 HEXACHLOROCYCLO-PENTADIENE ND ug/L 525.2 0.02000 INDENO(1,2,3-CD)PYRENE ND ug/L 525.2 0.05000 LINDANE (BHC - GAMMA) ND ug/L 525.2 0.02000 MALATHION ND ug/L 525.2 0.05000 METHOXYCHLOR ND ug/L 525.2 0.02000 METOLACHLOR ND ug/L 525.2 0.25000 METRIBUZIN ND ug/L 525.2 0.05000		FLUORENE	ND		ug/L	525.2	C	0.05000		
HEXACHLOROBENZENE ND ug/L 525.2 0.02000 HEXACHLOROCYCLO-PENTADIENE ND ug/L 525.2 0.02000 INDENO(1,2,3-CD)PYRENE ND ug/L 525.2 0.05000 LINDANE (BHC - GAMMA) ND ug/L 525.2 0.02000 MALATHION ND ug/L 525.2 0.05000 METHOXYCHLOR ND ug/L 525.2 0.02000 METOLACHLOR ND ug/L 525.2 0.02000 METRIBUZIN ND ug/L 525.2 0.05000		HEPTACHLOR	ND		ug/L	525.2	(0.02000		
HEXACHLOROGYCLO-PENTADIENE ND ug/L 525.2 0.02000 INDENO(1,2,3-CD)PYRENE ND ug/L 525.2 0.05000 LINDANE (BHC - GAMMA) ND ug/L 525.2 0.02000 MALATHION ND ug/L 525.2 0.05000 METHOXYCHLOR ND ug/L 525.2 0.02000 METOLACHLOR ND ug/L 525.2 0.05000 METRIBUZIN ND ug/L 525.2 0.05000		HEPTACHILOR EPOXIDE	ND		ug/L	525.2	(0.02000		
INDENO(1,2,3-CD)PYRENE ND ug/L 525.2 0.05000 LINDANE (BHC - GAMMA) ND ug/L 525.2 0.02000 MALATHION ND ug/L 525.2 0.05000 METHOXYCHLOR ND ug/L 525.2 0.02000 METCLACHLOR ND ug/L 525.2 0.25000 METRIBUZIN ND ug/L 525.2 0.05000		HEXACHLOROBENZENE	ND		ug/L	525.2	(0.02000		
LINDANE (BHC - GAMMA) ND ug/L 525.2 0.02000 MALATHION ND ug/L 525.2 0.05000 METHOXYCHLOR ND ug/L 525.2 0.02000 METOLACHLOR ND ug/L 525.2 0.25000 METRIBUZIN ND ug/L 525.2 0.05000		HEXACHLOROGYCLO-PENTADIENE	ND		ug∕L	525.2	C	0.02000		
MALATHION ND ug/L 525.2 0.05000 METHOXYCHLOR ND ug/L 525.2 0.02000 METOLACHLOR ND ug/L 525.2 0.25000 METRIBUZIN ND ug/L 525.2 0.05000		INDENO(1,2,3-CD)PYRENE	ND		ug/L	525.2	(0.05000		
METHOXYCHLOR ND ug/L 525.2 0.02000 METOLACHLOR ND ug/L 525.2 0.25000 METRIBUZIN ND ug/L 525.2 0.05000		LINDANE (BHC - GAMMA)	ND		ug/L	525.2	C	0.02000		
METCLACHLOR ND ug/L 525.2 0.25000 METRIBUZIN ND ug/L 525.2 0.05000		MALATHION	ND		ug/L	525.2	(0.05000		
METCLACHLOR ND ug/L 525.2 0.25000 METRIBUZIN ND ug/L 525.2 0.05000		METHOXYCHLOR	ND		ug/L	525.2	C	0.02000		
METRIBUZIN ND ug/L 525.2 0.05000		METOLACHLOR	ND		_	525.2	(0.25000		
		METRIBUZIN	ND		ug/L	525.2				
		NAPTHALENE								

[&]quot;Notation:

[%] Recovery = (Result of Analysis)/(True Value) * 100

NA = Indicates % Recovery could not be calculated.

QCS: Quality Control Sample, a solution containing known concentrations of method analytes which is used to fortify an aliquot of reagent matrix. The QCS is obtained from an external source and is used to check lab performance.

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ME or LRB. Method Blank or Laboratory Reagent Blank, an aliquot of reagent matrix is analyzed exactly like a sample, and its purpose is to determine if there is background contamination.



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SAMPLE INDEPENDENT QUALITY CONTROL REPORT

Method Blank

Reference Number: 09-02151

			True			%	QC	
Balch	Analyte	Result	Value	Units	Method	Recovery Limits	Qualifier Type*	Comment
525_090217	PARATHION	ND		ug/L	525.2	0.0500	O MB	
	PENTACHLOROPHENOL	ND		ug/L	525.2	0.0400	ю	
	PERYLENE-D12 (Surr)	106		%	525.2			
	PHENANTHRENE	ND		ug/L	525.2	0.0500	0	
	PROPACHLOR	ND		ug/L	525.2	0.0500	ю	
	PYRENE	ND		ug/L	525.2	0.0500	Ю	1
	PYRENE-D10 (Surr)	92		%	525.2			
	SIMAZINE	ND		ug/L	525.2	0.0200	10	
	TERBACIL	ND		ug/L	525.2	0.0500	10	
	TRIFLURALIN	ND		ug/L	525.2	0.0500	10	
	TRIPHENYLPHOSPHATE (Suit)	93		%	525.2			
531_090219	3-HYDROXYCARBOFURAN	ND		ug/L	531.2	0.5000	ю мв	
	ALDICARB	ND		ug/L	531.2	0.2500	10	
	ALDICARB SULFONE	ND		ug/L	531.2	0.4000	10.	
	ALDICARB SULFOXIDE	ND		ug/L	531.2	0.2500	10	
	BDMC (SURR)	114		%	531.2	0.0000	10	
	CARBARYL	ND		ug/L	531.2	0.5000	10	
	CARBOFURAN	ND		ug/L	531.2	0.4500	10	
	METHIOCARS	ND		ug/L	531.2	1,0000	00	
	METHOMYL	ND		ug/L	531.2	0.2500	ю	
	OXYMAL	ND		υg√L	531.2	1.0000	ю	
	PROPOXUR (BAYGON)	ND		ug/L	531.2	0.2500	x	÷
ec_090216	ELECTRICAL CONDUCTIVITY	ND		uS/cm	SM2510 B	2.5000	00 MB	
ec_090216	ELECTRICAL CONDUCTIVITY	ND		uS/cm	SM2510 B	2.5000	00 MB	
ec_090216	ELECTRICAL CONDUCTIVITY	ND		uS/cm	SM2510 B	2.5000	00 MB	
ec_090216	ELECTRICAL CONDUCTIVITY	ND		uS/cm	SM2510 B	2.5000	ID MB	-
OPHOS-090213	ORTHO-PHOSPHATE	ND		mg/L	SM4500-P F	0.100	00 MB	

^{*}Notation:

[%] Recovery = (Result of Analysis)/(True Value) * 100

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SAMPLE INDEPENDENT **QUALITY CONTROL REPORT**

Method Blank

Reference Number: 09-02151

			True			%		QC	
Batch	Analyte	Result	Value	Units	Method	Recovery Lin	mits	Qualifier Type*	Comment
TDS_090218	TOTAL DISSOLVED SOLIDS	ND		mg/L	SM2540 C	2.	50000	MB	
TDS_090218	TOTAL DISSOLVED SOLIDS	ND		mg/L	SM254D C	2.	50000	MB	
TDS_090218	TOTAL DISSOLVED SOLIDS	ND		mg/L	SM2540 C	2.9	50000	МВ	
turb_090213	TURBIDITY	ND .		NTU	180.1	. 0.0	02000	MB	

^{*}Notation:

[%] Recovery # (Result of Analysis)/(True Value) * 100

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SAMPLE INDEPENDENT QUALITY CONTROL REPORT

Quality Control Sample

Reference Number: 09-02151

.			True			%		QC	
Batch	Analyte	Result	Value	Units	Method	Recove	ery Limits	Qualifier Type'	Comment
200.7-090216A	HARDNESS	130	132.3	mg/L	200.7	98	80-120	ocs	
531_090219	3-HYDROXYCARBOFURAN	45.1	40	ug/L	531.2	113	70-130	ocs	
	ALDICARB .	37.4	37.3	ug/L	531.2	100	70-130		
	ALDICARB SULFONE	46.3	40.2	ug/L	531.2	115	70-130		
•	ALDICARB SULFOXIDE	50.3	44.9	ug/L	531.2	112	70-130		
	BDMC (SURR)	114		%	531.2		70-130		
	CARBARYL	43.4	46	ug/L	531.2	94	70-130		
	CARBOFURAN	68.2	60.9	ug/L	531.2	112	70-130		
	METHIOCARB	128	121	ug/L	531,2	106	70-130		
	METHOMYL	67.2	61.4	ug/L	531.2	109	70-130		
	OXYMAL .	65.6	59.9	ug/L	531.2	110	70-130		
	PROPOXUR (BAYGON)	100	96.7	ug/L	531.2	103	70-130		
COD_090216	CHEMICAL OXYGEN DEMAND	87	92	mg/L	SM5220 D	95	80-120	QCS	
D090304A	BROMATE	0.0166	0.0157	mg/L	300,1	106	75-125	QCS	
				11.5.4	000.1	100	70-123	QU3	
ec_090216	ELECTRICAL CONDUCTIVITY	16 1	150.5	uS/cm	SM2510 B	107	80-120	QCS	
on 000016	FI SCIPION CONDUCTOR						00 120	400	
ec_090216	ELECTRICAL CONDUCTIVITY	. 161	150.5	uS/cm	SM2510 B	107	80-120	QCS	
ec_090216	ELECTRICAL CONDUCTIVITY	162	150.5	uS/cm	SM2510 B	108	80-120	QCS	
ec_090216	ELECTRICAL CONDUCTIVITY	162	150.5	uS/cm	SM2510 B	108	80-120	QCS	
1090213A	CHLORIDE	30.7	30.0	mg/L	300.0	102	80-120	000	
	NITRATE-N	2.44	2.50	mg/L	300.0	98		QCS	
		4.	2.00	419te	300.0	90	80-120		
OPHOS-090213	ORTHO-PHOSPHATE	0.47			_				
UF1103*V30213	OMHOPPIUSPRAIC	0.47	0.49	mg/Ļ	SM4500-P F	96	70-130	QCS	

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SAMPLE INDEPENDENT QUALITY CONTROL REPORT

Quality Control Sample

Reference Number: 09-02151

Report Date: 03/05/09

			True			%	Q C	
Batch	Analyte	Result	Válue	Units	Method	Recovery Limits		Comment
turb_090213	TURBIDITY	 0.99	1.00	NTU	180.1	99 70-13	0 OCS	

*Notation:

% Recovery = (Result of Analysis)/(True Value) * 100

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Reference Number: 09-02151 Report Date: 3/5/2009 Page 1 of 5 Comments 물물 퉏 뎔 ᅙ 훒 훒 5 9 뎚 밁 집 Qualifier S %RPD LImits Å D 3 9 Ä i i £ £ Ï 1111 F F 545 Ž 9-9 ğ 545 0.0 28.83 21.1 5.3 ₹ 0 Ξ 0.0 10.5 3.7 0.0 Ξ 8 9.5 7. Ξ 0.3 9.6 Duplicate and Matrix Spike/Matrix Spike Duplicate Report **QUALITY CONTROL REPORT** Corporate Office 800.755.9295 • 360.757.1400 • 360.757.14027ax
Bellingham WA 805 Orchard Dr Sulta 4 - 98225
Mexokriden 360.571.0688 • 360.671.1577ax mg CaCO3/L uS/cm US/CH ₩2/Sh пgА mgA. 훒 ם 퉏 7<u>6</u>F J. J.GE 뼥 Burlington WA 1620 S Wahrut St - 98233 Corporate Office 800.755.9295 • 360.757.1 Dupficate Result 0.09 19.2 0.03 0.17 109 0.25 0.91 0.24 23 8 2 38 7.8 Result 0.21 19.3 0.04 0.29 0.24 50 6.0 2.3 등 4 86 얾 7: 7.6 4 4470 1,3-DIMETHYL-2-NITROBENZENE (Sur TRIPHENYLPHOSPHATE (Sum) CHEMICAL OXYGEN DEMAND 4313 ELECTRICAL CONDUCTIVITY ELECTRICAL CONDUCTIVITY ELECTRICAL CONDUCTIVITY PERYLENE-D12 (Sur) 4483 ORTHO-PHOSPHATE 4470 PYRENE-D10 (Surt) 4316 HARDNESS NITRATE-N NITRATEN CHLORIDE 4455 NITRATE-N CHLORIDE CHLORIDE 3913 ATRAZINE 4470 ATRAZINE BROMACIL Sample Analyte 4470 4483 4470 4506 4470 4506 4483 4483 4506 4471 4455 Duplicate OPHOS-090213 200.7-090216A COD_090216 525_090217 EC_090216 PH_090213 D090304A 1090213A

Batch

%RPD = Relative Percent Difference

NA = indicates %RPD could not be calculated

Matrix Spike (MS/Matrix Spike Duplicate (MSD) analyses are used to determine the accuracy (MS) and precision (MSD) of a analytical method in a given sample matrix. Therefore, the usefulness of this report is limited to samples of similar matrices analyzed in the same analytical batch.



Reference Number: 09-02151 Report Date: 3/5/2009

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Duplicate

Comments 민 줊 a Da 5 2 ö Limit 1 1 0-20 0-21 0-50 8 3 SRPD 5,0 7.0 0.0 2.2 8.0 pH Units PH Units 돌등품 Mg/L D TN Duplicate Resux 8.55 8.08 4.86 0.23 5.97 8 Result 6.50 4.53 8.02 0.23 5.94 8 4508 TOTAL DISSOLVED SOLIDS 4378 HYDROGEN ION (pH) 4482 HYDROGEN ION (pH) 4506 HYDROGEN ION (pH) 4482 TURBIDITY 4508 TURBIDITY Sample Analyte TURB_090213 TDS_090218 Batch

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Matrix Spike

Reference Number: 09-02151 Report Date: 3/5/2009 Page 3 of 5

Matrix Spike	oike				Dupficate									
				Spike	Spilke	Spike		Percent Recovery	ecovery				8	
Batch	Sample	Sample Analyte	Result	Result	Result	Conc	Units .	SE	MSD	Limits	KRPD	Limits	Qualifier	Comments
200.7-090216A	_													
	4237	HARDNESS	2	71.6	6.07	69.5	mg CaCO3/L 103	8	102	80-120	1:0	09-0	LFM	
	4316	4316 HARDNESS	19.3	69.3	88.7	69.5	mg CaCO3/L 101	5	100	80-120	6.0	09-0	LFM	
515.4_090219														
	4470	2,4-D	₽	4.45	4.4	_	g Jûr	2	88	70-130	₽	0-20	LFM	
	4470	2,4,5 - TP (SILVEX)	9	2.65		2.5	Jg/L 1	106	108	70-130	E :	0-20	LFM	
	4470	PENTACHLOROPHENOL	Q	2.8		2.5	ig/L 1	112	112	70-130	0.0	0-30	L'FM	
	4470	DALAPON	2	30.4		32.5	g. Jou	*	96	70-130	2.9	0-20	LFM	
	4470	DINOSEB	S	5.3	5.3	_		9	106	70-130	0.0	020	LFM	
	4470	PICLORAM	9	2.7		_	•	108	118	70-130	7.1	05-0	LFM	
	4470	DICAMBA		2.7				108	108	70-130	0.0	9	LFM	
	4470	DCPA (ACID METABOLITES)	2	2.3		2.5		92	4	70-130	9	8	LFM	
	4470	2,4 DB	2	19.6		8		98	66	70-130	0.5	0-20	LFM	
	4470	2,4,5 T	9	2.7		2.5		108	108	70-130	0.0	99	LFM	
	4470	BENTAZON	2	4.8	5.4	_	ug/L 9	96	108	70-130	1 .	05-0	LFM	
-	4470	DICHLORPROP	2	~		_	tour 9	76	68	70-130	7.	0-20	LFW	
	4470	ACIFLUORFEN	오	2.6	2.65	2.5	Jg/L 1	104	106	70-130	,	0-20	LFM	
	4470	2,4 - DCAA (SURR)	87	&	8	•	*		Ą.	70-130	Š	0-20	LFM	
525_090217														
I	3079	ENDRIN	9	0.96	_	_	6 T/Gr	96	ĄN	70-130	¥	0-60	LFM	
	3079	LINDANE (BHC - GAMMA)	9	1.0 1.0	_	_		101	ΑN	70-130	¥	0-80	LFM	
	3079	METHOXYCHLOR	2	1.12	-	_		112	A	70-130	¥	08-0	LFM	
	3079	ALACHLOR	Ð	2.07		_	ug/L 1	104	Ą.	70-130	¥	08-0	LFM	
	3079	ATRAZINE	9	2.28	.,	_	ug/L 1	114	NA A	70-130	¥	09-0	LFM	
	3079	BENZO(A)PYRENE	Q	1.19	_	_	ug/L 1	119	ď Z	70-130	¥	9	EP.	
	3079	CHLORDANE, TECHNICAL	2	0.85	_	_	ug/L 8	85	A N	70-130	×	90	LFIM	
	3079	DI(ETHYLHEXYL)-ADIPATE	9	1.27	•	_		127	NA	70-130	X A	0-60	LFM	
	3079	DI(ETHYLHEXYL)-PHTHALATE	2	1.14	•	_		114	42	70-130	Ϋ́	09-0	LFM	
	3079	HEPTACHLOR	2	1.08	-	_	ug/L 1	108	Y.	70-130	¥	0.60	LFM	
	3079	HEPTACHLOR EPOXIDE	ş	1.2	_	_		120	¥.	70-130	Ā	충	LFM	
	3079	HEXACH, OROBENZENE	Ş	0.96		_		96	¥Z.	70-130	¥	0-60	ÆĽ	
	3079	HEXACHLOROCYCLO-PENTADIENE	9	1.05	_	_	ug/l. 1	105	¥	70-130	¥	08-0	LFM	
	3079	SIMAZINE	2	1.13	•	_		113	٧Z	70-130	¥	9-6	LFR	
	3079	PENTACHLOROPHENOL	2	3.98	4	_	ug/l. 1	100	NA	70-130	¥	0-20	LFIM	
	3079	ALDRIN	2	1.04	•	_		104	ΑN	70-130	¥	09-0	LFM	
	3079	BUTACHLOR	9	1.18	•	_	,	116	VA	70-130	¥	09-0	LFM	

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Reference Number: 09-02151 Report Date: 3/5/2009

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Matrix Spike

Batch

Spike			Duplicate										
		Spike	Spike	Spike		Percen	Percent Recovery				ö		
Sample Analyte	Result	Result	Result	Conc	Units	W	MSD	Limits	%RPD	Limits	Cuekher	S	Comments
3079 DIELDRIN	Q.	1.09		-	ug/L	109	ΑN	70-130	Ā	09-0		LFR	
3079 METOLACHLOR	오	1.12		-	ug/L	112	Ą	70-130	₹ Z	8		F	
3079 METRIBUZIN	무	1,05		_	ng/l.	105	ν V	70-130	Ϋ́	용		LFM	
3079 PROPACHLOR	무	1.17		_	ug/L	111	AA	70-130	Ą.	8		L.F.M	
3079 BROMACIL	윷	1,1		_	ng/L	Ŧ	Y V	70-130	Ą	8		LFM	
3079 TERBACK	2	1.37		-	ug/L	137	Y Y	70-130	¥		몆	F	
3079 DIAZINON	Q	3.47		6	ug/L	118	¥	70-130	AN	080		LFM	
3079 EPTC	S	1.01		-	ug/L	흔	Ą	70-130	¥	240		LFM	
3079 4,4-DDD	2	1.09		-	ug/L	109	¥	70-130	¥	09-0		LFM	
3079 4,4-DDE	Q	-		•	ug/L	5	Ā	70-130	٨	0-60		LFM	
3079 4,4-DDT	2	1.09			ug/L	109	AN	70-130	٧	09-0		LFM	
3079 CYANAZINE	S	2.43		2	ug/L	72	Ą	70-130	Ą	0-60		FM	
3079 MALATHION	2	2.71		m	ug/L	8	¥	70-130	¥	8		LFM	
3079 PARATHION	9	2.3		2	ug/L	115	Ą	70-130	¥	8		LF.M	
3079 TRIFLURALIN	ð	1.07		_	ug/L	107	Ą	70-130	¥	09-0		LFM	
3079 FLUORENE	9	1.13		-	ng√.	=======================================	Ą	70-130	¥	09-0		LFM	•
3079 ACENAPHTHYLENE	9	1.08		_	ng/L	99	Ā	70-130	¥.	08-0		LFM	
3079 ANTHRACENE	N O	96.0		_	ug/L	86	¥	70-130	Ā	09-0		LFM	
3079 BENZ(A)ANTHRACENE	9	1.05		_	ug/L	5	¥	70-130	¥	8		LFM	
3079 BENZO(B)FLUORANTHENE	õ	1.05		_	ug/L	105	Ą	70-130	Ą	09-0		EFM.	
3079 BENZO(G,H,I)PERYLENE	ð	-		-	ng/L	100	Ą	70-130	Ž	09-0		LFM	
3079 BENZO(K)FLUORANTHENE	ð	1.01		-	ng/L	É	¥	70-130	¥	9-60		F	
3079 CHRYSENE	9	1.09		_	ug/t	109	¥	70-130	¥	0-60		N.	
3079 DIBENZO(A,H)ANTHRACENE	2	1.13		-	ug/L	13	¥	70-130	¥	090		LFM	
3079 INDENO(1,2,3-CD)PYRENE	2	1.21			ng/L	121	ž	70-130	¥	09-0		F	
3079 PHENANTHRENE	S	1.03		_	ug/L	103	NA	70-130	¥	09-0		LFM	
3079 PYRENE	2	1.01		_	Van	Ę	Ą	70-130	¥	9-80		LFM	
3079 BENZYL BUTYL PHTHALATE	Ş	1,14		_	ug/L	114	Ą.	70-130	¥	9		LF.	
3079 DIN-BUTYL PHTHALATE	ç	1.09			ng/¦	109	AA	70-130	ž	8		LFM	
3079 DIETHYL PHTHALATE	Q	1.16		_	ng/L	116	¥	70-130	ž	9		LFM	
3079 DIMETHYL PHTHALATE	S	=			ug/L	2	Ą	70-130	¥	040		LFIM	
3079 1,3-DIMETHYL-2-NITROBENZENE (Sur	83	26			*		ΑĀ	70-130	ž	09-0		M	
3079 PYRENE-D10 (Sum)	93	93			×		¥	70-130	ž	0-60		LFM	
3079 PERYLENE-D12 (Surt)	107	1 08			*		¥	70-130	ž	09-0		LFM	
3079 TRIPHENYLPHOSPHATE (Sum)	93	98			*		V	70-130	¥	090		LFW	

531_090219

%RPD = Relative Percent Difference

NA = Indicates %RPD could not be calculated

Matrix Spike (MSV/Matrix Spike Duplicate (MSD) analyses are used to determine the accuracy (MS) and precision (MSD) of a analytical metrice analytical metrices analytical batch.



Page 5 of 5 Reference Number: 09-02151

Report Date: 3/5/2009

Matrix Spike	sike				Duplicate									
:				Spike	Spilke	Spike		Percent	Percent Recovery				ac	
Batch	Sample	Sample Analyte	Result	Result	Result	Conc	Units	MS	MSD	Limits	%RPD	Limits	Quaditier	Comments
:	4470	4470 OXYMAL	Q	13.7	14.8	5	ng/L	9	8	70-130	7.7	- - - - -	LFM	
	4470	CARBOFURAN	2	12.8	13.4	1 5	ugvi.	85	88	70-130	4.6	0-20	LFN	
	4470	ALDICARB SULFOXIDE	2	5	14.6	5	ug/L	8 7	97	70-130	11.6	09-0	LFM	
	4470	ALDICARB SULFONE	2	12.4	14	15	ug/L	83	83	70-130	12.1	9	LFM	
	4470	METHOMYL	9	13	14.5	15	ug/L	87	26	70-130	10.9	9-50	LFM	
	4470	3-HYDROXYCARBOFURAN	Ş	12.8	14.2	15	пgЛ	85	98	70-130	10.4	0-50	LFM	
	4470	ALDICARB	Ş	12.6	13.3	15	ug/L	84	69	70-130	5,4	0-50	LFM	
	4470	CARBARYL	Ð	11.7	12.3	6	ηğη	82	82	70-130	5.0	0-20	LFM	
	4470	PROPOXUR (BAYGON)	Q	12.6	13	5	ug/L	\$	87	70-130	3.1	0-20	LFM	
	4470	METHIOCARB	2	12.3	13.2	1 5	ηūγ	83	88	70-130	7.1	0-20	LFM	
	4470	BDMC (SURR)	103	106	109		%		NA A	70-130	¥,	050	LFM	
COD_090216														
ı	4483	CHEMICAL OXYGEN DEMAND	4	63	65	8	mg/l.	86	102	80-120	4.0	080	LFM	-
D090304A														
	4480	BROMATE	2	0.011		0.010	mg/L	110	AN A	75-125	ž	09-0	LFW	
	5536	BROMATE	용	0.0094		0.010	mg/L	26	¥.	75-125	¥	0-80	LFM	
f090213A														
	4455	NITRATE-N	0.29	1.24		0.1	щgЛ	92	Y.	80-120	¥	99	LFM	
	4455	CHLORIDE	1.7	2.75		1.00	mg/L	105	A	80-120	¥	080	LFW	
	4483	NITRATEN	6.0	19.9		20.00	тgЛ	98	¥	80-120	۲	99	F	
	4483	CHLORIDE	2.3	21.8		20.00	mg/L	88	¥	80-120	¥	09-0	F	
	4506	NITRATE-N	1.0	1.08		1.00	mg/L	86	Ā	80-120	¥	09-0	F	
	4506	CHLORIDE	2.6	8.4		1.00	mg/L	8	¥	80-120	¥	0.80	S	
OPHOS-090213	60													
	4483	4483 ORTHO-PHOSPHATE	0.24	1.25	1.26	0.1	щвуг	퉏	102	70-130	1.0	0-20	LFM	

%RPD = Relative Percent Difference

NA = Indicates %RPD could not be calculated

Matrix Spike (MS)Matrix Spike Duplicate (MSD) analyses are used to determine the accuracy (MS) and precision (MSD) of a analytical method in a given sample matrix. Therefore, the usefulness of this report is limited to samples of similar matrices analytical batch.





Qualifier Definitions

Reference Number: 09-02151 Report Date: 03/05/09

Qualifier	Definition
B5	The compound was detected in the sample below the State Reporting Limit, result is biased high.
HQ	High QCS recovery due to increased detector response of the sample extract. The continuing calibration checks are within acceptance limits.
J	Indicates an estimated concentration. This occurs when an analyte concentration is below the calibration curve but is above the method detection limit.
11	The amount detected is below the Method's Reporting Level but greater then the lab's Practical Quantitation Level.
M	Matrix induced bias assumed.
Q1	Acceptance limits do not apply. This compound is only qualitatively identified by the method.
s	Spiking amount was lower than the 5:1 spike to background (sample amount) basis for performance criteria. The reported criteria does not apply due to increased errors in measurement of both sample and spike concentration.

Note: Some qualifier definitions found on this page may pertain to results or QC data which are not printed with this report.

Chain of Custody / Analysis Request FORM: COC 01-05-2009 Ship Address: 810 S Main Street Report to: Walla Walla Basin Watershed Cour Sampled by œ ທີ Phone 5 O Email: Sample Receipt Request (Must include FAX or Email) Relinquished by Instructions 1. Use one line per sample Location. 2. Be specific in analysis requests. Enter number of containers (NEW) List each metal individually (NEW) Check off analyses to be performed for each sample Loaction Field ID Ç I DAIL - I WHITHELING 541.938-2170 FAX Milton-Freews St Locher Road Recharge Sites Γroy Baker さららい からからう OR Zp: 97862 Location 2/12/69 Date Phone: Turn Around Time Required 11:00 Time Quickest (100% surcharge) Phone Call Rec Half-time (50% surcharge Standard Emergency (Phone Call Reg BIII to City Сотр. Grab/ ☐ Visa P.O.# Phone: Card# Address: 810 South Main Street Received by Sample Matrix* Walla Walla Basin Watershed Counc Milton-Freewe st OR ₹ ਨ W - water Q Q DW - drinking water Date (Please complete all applicable shaded sections) ٥ ชีวิก CAR Time 2 FAX A Bromate SW - surface water GW - Ground water Hardness Ž $\mathbf{X}\mathbf{A}\mathbf{x}$ to P 97862 Date NO3, COD Analyses Requested Email: Ref# SOC Package Check Regulatory Program Time WW - waste water S - soil Other RCRA/CERCLA Clean Water Act Safe Drinking Water Act For Lab Use Only TDS, CI, O-Phos.pH,Turb,Ec Custody seals intact Chain of custody & labels agree Samples received intact Sample temp Other_ C satisfactory 09-0215 S S 805 W. Orchard Dr. Suite 4 Bellingham, WA 98225 Number of Containers Burlington, WA 98233 1.800.755.9295 TY THE VERNE ABORATORIES 1620 S. Walnut St. Total Containers Page Special Instructions Conditions on Receipt CO007096

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Corporate Office 800.755.9295 • 360.757.1400 • 360.757.1402fax

Bellingham WA 805 Orchard Dr Suite 4 - 98225

Macrobiology 360.671.0688 • 360.671.1577fax

Page 1 of 3

Data Report

Client Name: Walla Walla Basin Watershed Council

810 S Main Street

Milton-Freewater, OR 97862

Reference Number: 09-03591

Project: Locher Road Recharge Sites

Report Date: 3/18/09

Date Received: 3/13/09

Peer Review:

Lab N	cription: L1 - Locher Road 1 lumber: 7335						-	le Date: 3 cted By: 1			
CAS ID#	Parameter	Result	PQL	MDL	Units	DF	Method	Analyzed	Analys	t Batch	Comment
14797-55-8	NITRATE-N	5.21	0.100	0.015	mg/L	1	300.0	3/13/09	ВЈ	1090313A	
E-10173	TOTAL DISSOLVED SOLIDS	225	10	6	mg/L	1	SM2540 C	3/16/09	CCN	TDS_090316	
16887-00-6	CHLORIDE	5.9	20	0.012	mg/L	1	300.0	3/13/09	BJ	1090313A	
14265-44-2	ORTHO-PHOSPHATE	0.22	0.01	0.002	mg/L	1	SM4500-P F	3/13/09	so	OPHOS-090313	
E-10139	HYDROGEN ION (pH)	6.72			pH Units	1	SM4500-H+ B	3/13/09	CCN	PH_090313	
E-10617	TURBIDITY	8.12	0.05	0.03	NTU	1	180.1	3/13/09	CCN	TURB_090313	
E-10184	ELECTRICAL CONDUCTIVITY	355	10		uS/cm	1	SM2510 B	3/16/09	CCN	EC_090316	
E-11778	HARDNESS	150	3.30	0.055	mg CaCO3/L	1	200.7	3/16/09	вл	200.7-090316A	
E-10117	CHEMICAL OXYGEN DEMAND	ND	8	2.47	mg/L	1	SM5220 D	3/16/09	MAK	COD_090316	
15541-45-4	BROMATE	ND	0.005	0.00118	mg/L	1	300.1	3/17/09	MVP	D090317A	
Lab N	ription: L2 - Locher Road 2 umber: 7336						Colle	ole Date:	Baker		
Lab N		Result	PQL	MDL	Units	DF			Baker	t Batch	Comment
Lab N CAS ID#	umber: 7336	Result	PQL 0.100	MDL 0.015	Units mg/L	DF 1	Colle	cted By:	Baker Analys		Comment
Lab N CAS ID# 14797-55-8	umber: 7336 Parameter						Colle	cted By: Analyzed	Baker Analys вл	t Batch	Comment
Lab N CAS ID# 14797-55-8 E-10173	umber: 7336 Parameter NITRATE-N	5.42	0.100	0.015	mg/L	1	Colle Method	Analyzed	Baker Analys BJ CCN	t Batch 1090313A	Comment
	umber: 7336 Parameter NITRATE-N TOTAL DISSOLVED SOLIDS	5.42 196	0.100 10	0.015 6	mg/L mg/L	1	Method 300.0 SM2540 C	Analyzed 3/13/09 3/16/09	Baker Analys BJ CCN BJ	t Batch 1090313A TDS_090316	Comment
Lab N CAS ID# 14797-55-8 E-10173 16887-00-6 14265-44-2	umber: 7336 Parameter NITRATE-N TOTAL DISSOLVED SOLIDS CHLORIDE	5.42 196 6.2	0.100 10 20	0.015 6 0.012	mg/L mg/L mg/L	1 1 1	Colle Method 300.0 SM2540 C 300.0	Analyzed 3/13/09 3/16/09 3/13/09	Analys BJ CCN BJ so	t Batch 1090313A 1DS_090316 1090313A	Comment
Lab N CAS ID# 14797-55-8 E-10173 16887-00-6	umber: 7336 Parameter NITRATE-N TOTAL DISSOLVED SOLIDS CHLORIDE ORTHO-PHOSPHATE	5.42 196 6.2 0.22	0.100 10 20	0.015 6 0.012	mg/L mg/L mg/L mg/L	1 1 1 1	Method 300.0 SM2540 C 300.0 SM4500-P F	Analyzed 3/13/09 3/16/09 3/13/09	Analys BJ CCN BJ SO CCN	t Batch 1090313A 1050316 1090313A 0PHOS-090313	Comment
Lab N CAS ID# 14797-55-8 E-10173 16887-00-6 14265-44-2 E-10139	umber: 7336 Parameter NITRATE-N TOTAL DISSOLVED SOLIDS CHLORIDE ORTHO-PHOSPHATE HYDROGEN ION (pH)	5.42 196 6.2 0.22 6.59	0.100 10 20 0.01	0.015 6 0.012 0.002	mg/L mg/L mg/L mg/L pH Units	1 1 1 1 1	Colle Method 300.0 SM2540 C 300.0 SM4500-P F SM4500-H+ B	Analyzed 3/13/09 3/16/09 3/13/09 3/13/09 3/13/09	Baker Analys BJ CCN BJ SO CCN CCN	t Batch 1090313A 110S_090316 1090313A 0PHOS-090313 PH_090313	Comment
Lab N CAS ID# 14797-55-8 E-10173 16887-00-6 14265-44-2 E-10139 E-10617	umber: 7336 Parameter NITRATE-N TOTAL DISSOLVED SOLIDS CHLORIDE ORTHO-PHOSPHATE HYDROGEN ION (pH) TURBIDITY	5.42 196 6.2 0.22 6.59 2.02	0.100 10 20 0.01	0.015 6 0.012 0.002	mg/L mg/L mg/L mg/L pH Units	1 1 1 1 1 1 1	Colle Method 300.0 SM2540 C 300.0 SM4500-P F SM4500-H+ B 180.1	Analyzed 3/13/09 3/13/09 3/13/09 3/13/09 3/13/09 3/13/09	Analys BJ CCN BJ SO CCN CCN CCN	t Batch 1090313A 1090313A 1090313A 0PH0S-090313 PH_090313 1URB_090313	Comment
Lab N CAS ID# 14797-55-8 E-10173 16887-00-6 14265-44-2 E-10139 E-10617 E-10184 E-11778	umber: 7336 Parameter NITRATE-N TOTAL DISSOLVED SOLIDS CHLORIDE ORTHO-PHOSPHATE HYDROGEN ION (pH) TURBIDITY ELECTRICAL CONDUCTIVITY	5.42 196 6.2 0.22 6.59 2.02	0.100 10 20 0.01 0.05	0.015 6 0.012 0.002	mg/L mg/L mg/L mg/L pH Units NTU uS/cm	1 1 1 1 1 1 1	Colle Method 300.0 SM2540 C 300.0 SM4500-P F SM4500-H+ B 180.1 SM2510 B	Analyzed 3/13/09 3/16/09 3/13/09 3/13/09 3/13/09 3/13/09 3/13/09 3/13/09	Baker Analys BJ CCN BJ SO CCN CCN CCN BJ	t Batch 1090313A 10S_090316 1090313A 0PHOS-090313 PH_090313 1URB_090313 EC_090316	Comment
Lab N CAS ID# 14797-55-8 E-10173 16887-00-6 14265-44-2 E-10139 E-10617 E-10184	umber: 7336 Parameter NITRATE-N TOTAL DISSOLVED SOLIDS CHLORIDE ORTHO-PHOSPHATE HYDROGEN ION (pH) TURBIDITY ELECTRICAL CONDUCTIVITY HARDNESS	5.42 196 6.2 0.22 6.59 2.02 295 118	0.100 10 20 0.01 0.05 10 3.30	0.015 6 0.012 0.002 0.03	mg/L mg/L mg/L pH Units NTU uS/cm mg CaCO3/L	1 1 1 1 1 1 1 1	Colle Method 300.0 SM2540 C 300.0 SM4500-P F SM4500-H+ B 180.1 SM2510 B 200.7	Analyzed 3/13/09 3/13/09 3/13/09 3/13/09 3/13/09 3/13/09 3/18/09	Baker Analys BJ CCN BJ SO CCN CCN CCN BJ MAK	t Batch 1090313A 1090313A 1090313A 1090313A 10PHOS-090313 1TURB_090313 1TURB_090313 1EC_090316 200.7-090316A	Comment
Lab N CAS ID# 14797-55-8 E-10173 16887-00-6 14265-44-2 E-10139 E-10617 E-10184 E-11778 E-10117	Parameter NITRATE-N TOTAL DISSOLVED SOLIDS CHLORIDE ORTHO-PHOSPHATE HYDROGEN ION (pH) TURBIDITY ELECTRICAL CONDUCTIVITY HARDNESS CHEMICAL OXYGEN DEMAND	5.42 196 6.2 0.22 6.59 2.02 295 118	0.100 10 20 0.01 0.05 10 3.30	0.015 6 0.012 0.002 0.03 0.055	mg/L mg/L mg/L pH Units NTU uS/cm mg CaCO3/L mg/L	1 1 1 1 1 1 1 1 1 1 1	Colle Method 300.0 SM2540 C 300.0 SM4500-P F SM4500-H+ B 180.1 SM2510 B 200.7 SM5220 D 300.1	Analyzed 3/13/09 3/16/09 3/13/09 3/13/09 3/13/09 3/13/09 3/16/09 3/16/09	Baker Analys BJ CCN BJ SO CCN CCN CCN BJ MAK MVP	t Batch 1090313A TDS_090316 1090313A OPHOS-090313 TURB_090313 EC_090316 200.7-090316A COD_090316 D090317A	Comment
Lab N CAS ID# 14797-55-8 E-10173 16887-00-6 14265-44-2 E-10139 E-10617 E-10184 E-11778 E-10117 15541-45-4 Sample Desc	Parameter NITRATE-N TOTAL DISSOLVED SOLIDS CHLORIDE ORTHO-PHOSPHATE HYDROGEN ION (pH) TURBIDITY ELECTRICAL CONDUCTIVITY HARDNESS CHEMICAL OXYGEN DEMAND BROMATE	5.42 196 6.2 0.22 6.59 2.02 295 118	0.100 10 20 0.01 0.05 10 3.30	0.015 6 0.012 0.002 0.03 0.055	mg/L mg/L mg/L pH Units NTU uS/cm mg CaCO3/L mg/L	1 1 1 1 1 1 1 1 1 1 1	Colle Method 300.0 SM2540 C 300.0 SM4500-P F SM4500-H+ B 180.1 SM2510 B 200.7 SM5220 D 300.1	Analyzed 3/13/09 3/16/09 3/13/09 3/13/09 3/13/09 3/13/09 3/15/09 3/16/09 3/17/09	Analys BJ CCN BJ SO CCN CCN CCN BJ MAK MVP	t Batch 1090313A TDS_090316 1090313A OPHOS-090313 TURB_090313 EC_090316 200.7-090316A COD_090316 D090317A	Comment

Notes:

D.F. - Dilution Factor

ND = Not detected above the listed practical quantitation limit (PQL) or not above the Method Detection Limit (MDL), if requested.

PQL = Practical Quantitation Limit is the lowest level that can be achieved within specified limits of precision and accuracy during routine laboratory operating conditions.



Page 2 of 3 Reference Number: 09-03591

Report Date:3/18/09

Data Report

14797-55-8	NITRATE-N	2.05	0.100	0.015	mg/L	1	300.0	3/13/09	вј	1090313A
E-10173	TOTAL DISSOLVED SOLIDS	91	10	6	mg/L	1	SM2540 C	3/16/09	CCN	TDS_090316
16887-00-6	CHLORIDE	2.2	20	0.012	mg/L	1	300.0	3/13/09	вл	1090313A
14265-44-2	ORTHO-PHOSPHATE	0.17	0.01	0.002	mg/L	1	SM4500-P F	3/13/09	so	OPHOS-090313
E-10139	HYDROGEN ION (pH)	6.61			pH Units	1	SM4500-H+ B	3/13/09	CCN	PH_090313
E-10617	TURBIDITY	2.01	0.05	0.03	NTU	1	180.1	3/13/09	CCN	TURB_090313
E-10184	ELECTRICAL CONDUCTIVITY	136	10		uS/cm	1	SM2510 B	3/16/09	CCN	EC_090316
E-11778	HARDNESS	50.5	3.30	0.055	mg CaCO3/L	1	200.7	3/16/09	BJ	200.7-090316A
E-10117	CHEMICAL OXYGEN DEMAND	ND	8	2.47	mg/L	1	SM5220 D	3/16/09	MAK	COD_090316
15541-45-4	BROMATE	ND	0.005	0.00118	mg/L	1	300.1	3/17/09	MVP	D090317A

•	ription: IN1 - Locher Road In umber: 7338						•	ole Date: 3 cted By: 1			. —
CAS ID#	Parameter	Result	PQL	MDL	Units	DF	Method	Analyzed	Analys	t Batch	Comment
14797-55-8	NITRATE-N	0.58	0.100	0.015	mg/L	1	300.0	3/13/09	BJ I	090313A	
E-10173	TOTAL DISSOLVED SOLIDS	826	10	6	mg/L	1	SM2540 C	3/16/09	CCN .	TDS_090316	
16887-00-6	CHLORIDE	22	20	0.012	mg/L	1	300.0	3/13/09	вл і	1090313A	
14265-44-2	ORTHO-PHOSPHATE	0.06	0.01	0.002	mg/L	1	SM4500-P F	3/13/09	so ·	OPHOS-090313	
E-10139	HYDROGEN ION (pH)	11.94	•		pH Units	1	SM4500-H+ B	3/13/09	CCN	PH_090313	
E-10617	TURBIDITY	21.7	0.05	0.03	NTU	1	180.1	3/13/09	CCN	TURB_090313	
E-10184	ELECTRICAL CONDUCTIVITY	2560	100		uS/cm	10	SM2510 B	3/16/09	CCN	EC_090316	
E-11778	HARDNESS	37.5	3.30	0.055	mg CaCO3/L	1	200.7	3/16/09	в	200.7- 090316 A	
E-10117	CHEMICAL OXYGEN DEMAND	ND	8	2.47	mg/L	1	SM5220 D	3/16/09	MAK	COD_090316	
15541-45-4	BROMATE	ND	0.005	0.00118	mg/L	1	300.1	3/17/09	MVP	D090317A	

•	ription: MC1 - Mudcreek 1 umber: . 7339						,	ole Date: 3 cted By: 1			
CAS ID#	Parameter	Result	PQL	MDL	Units	DF	Method	Analyzed	Analys	t Batch	Comment
14797-55-8	NITRATE-N	1.82	0.100	0.015	mg/L	1	300.0	3/13/09	BJ	1090313A	
E-10173	TOTAL DISSOLVED SOLIDS	179	10	6	mg/L	1	SM2540 C	3/16/09	CCN	TDS_090316	
16887-00-6	CHLORIDE	5.5	20	0.012	mg/L	1	300.0	3/13/09	ВЈ	1090313A	
14265-44-2	ORTHO-PHOSPHATE	0.15	0.01	0.002	mg/L	1	SM4500-P F	3/13/09	so	OPHOS-090313	
E-10139	HYDROGEN ION (pH)	7.69			pH Units	1	SM4500-H+ B	3/13/09	CCN	PH_090313	
E-10617	TURBIDITY	0.64	0.05	0.03	NTU	1	180.1	3/13/09	CCN	TURB_090313	
E-10184	ELECTRICAL CONDUCTIVITY	272	10		u\$/cm	ı	SM2510 B	3/16/09	CCN	EC_090316	
E-11778	HARDNESS	112	3.30	0.055	mg CaCO3/L	1	200.7	3/16/09	ผ	200.7-090316A	
E-10117	CHEMICAL OXYGEN DEMAND	ND	8	2.47	mg/L	1	SM5220 D	3/16/09	MAK	COD_090316	
15541-45-4	BROMATE	ND	0.005	0.00118	mg/L	1	300.1	3/17/09	MVP	D090317A	

ND = Not detected above the listed practical quantitation limit (PQL) or not above the Method Detection Limit (MDL), if requested.
PQL = Practical Quantitation Limit is the lowest level that can be acheived within specified limits of precision and accuracy during routine laboratory operating conditions.



Page 3 of 3 Reference Number: 09-03591 Report Date:3/18/09

Data Report

•	ription: MC2 - Mudcreek 2 umber: 7340						•	ole Date: 3 cted By: 1		-	
CAS ID#	Parameter	Result	PQL	MDL	Units	DF	Method	Analyzed	Analysi	Batch	Comment
14797-55-8	NITRATE-N	2.59	0.100	0.015	mg/L	1	300.0	3/14/09	ву і	090313A	
E-10173	TOTAL DISSOLVED SOLIDS	216	10	6	mg/L	1	SM2540 C	3/16/09	ÇCN :	TDS_090316	
16887-00-6	CHLORIDE	7.5	20	0.012	mg/L	1	300.0	3/13/09	BJ I	090313A	
14265-44-2	ORTHO-PHOSPHATE	0.15	0.01	0.002	mg/L	1	SM4500-P F	3/13/09	so (OPHOS-090313	
E-10139	HYDROGEN ION (pH)	7.62			pH Units	1	SM4500-H+ B	3/13/09	CCN	PH_090313	
-10617	TURBIDITY	2.68	0.05	0.03	NTU	1	180.1	3/13/09	CCN .	TURB_090313	
-10184	ELECTRICAL CONDUCTIVITY	373	10		uS/cm	ı	SM2510 B	3/16/09	CCN	EC_090316	
E-11778	HARDNESS	142	3.30	0.055	mg CaCO3/L	ı	200.7	3/16/09	вл :	200.7-090316A	
-10117	CHEMICAL OXYGEN DEMAND	ND	8	2.47	mg/L	1	SM5220 D	3/16/09	MAK	COD_090316	
15541-45-4	BROMATE	ND	0.005	0.00118	mg/L	1	300.1	3/17/09	MVP	D090317A	

ND = Not detected above the listed practical quantitation limit (PQL) or not above the Method Detection Limit (MDL), if requested.

PQL = Practical Quantitation Limit is the lowest level that can be acheived within specified limits of precision and accuracy during routine laboratory operating conditions.



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HERBICIDES IN DRINKING WATER

Client Name: Walla Walla Basin Watershed Council

810 S Main Street

Milton-Freewater, OR 97862

Project::

Field ID: L1

Sample Description: Locher Road 1

Sampled By: Baker

Sample Date: 3/12/09

Source Type: Sampler Phone:

Reference Number: 09-03591

Project: Locher Road Recharge Sites

Lab Number: 07335

Report Date: 4/15/09

Date Analyzed: 03/14/09 Date Extracted: 515.4_090313

Analyst: CO

Peer Review:

Analytical Method: 515.4

Herbicides in Drinking Water

CAS	COMPOUND	RESULTS	UNITS	PQL	MDL	MCL	COMMENT
EPA R	egulated			· ·			
4-75-7	2,4 - D	ND	ug/L	0.5	0.2	70	
3-72-1	2,4,5 - TP (SILVEX)	ND	. ug/L	0.25	0.04	50	
37-86-5	PENTACHLOROPHENOL	ND	ug/L	0.25	0.05	1	
5-99-0	DALAPON	ND	ug/L	3.25	0.9	200	
8-85-7	DINOSEB	ND	ug/L	0.5	0.2	7.	
918-02-1	PICLORAM	ND	ug/L	0.25	0.07	500	
Other							
-14028	DCPA (ACID METABOLITES)	ND	ug/L	0.25	0.1		
918-00-9	DICAMBA	ND	ug/L	0.25	0.05		
4-82-6	2,4 DB	ND	ug/L	2			
3-76 - 5	2,4,5 T	ND	ug/L	0.25	0.04		
5057-89-0	BENTAZON	ND	ug/L	0.5	0.2		
20-36-5	DICHLORPROP	ND	ug/L	0.75	0.5		
0594-66-6	ACIFLUORFEN	ND	ug/L	0.25	0.1		
1-36-5	3,5 - DICHLOROBENZOIC ACID	ND	ug/L	0.25	0.2		

NOTES:
If a compound is detected > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.

MCL (Maximum Contaminant Level) maximum permissible level of a contaminant in water established by EPA; Federal Action Levels are 0.015 mg/L for Lead and 1.3 mg/L for Copper. Sodium has a recommended limit of 20 mg/L. A blank MCL value indicates a level is not currently established.

Trigger Level: DOH Drinking Water Response level. Systems with compounds detected in excess of this level are required to take additional samples. Contact your regional DOH office.

ND (Not Detected): indicates that the parameter was not detected above the State Reporting Limit (SRL).



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HERBICIDES IN DRINKING WATER

Client Name: Walla Walla Basin Watershed Council

810 S Main Street

Milton-Freewater, OR 97862

Project::

Field ID: L2

Sample Description: Locher Road 2

Sampled By: Baker

Sample Date: 3/12/09 Source Type:

Sampler Phone:

Reference Number: 09-03591

Project: Locher Road Recharge Sites

Lab Number: 07336

Report Date: 4/15/09 Date Analyzed: 03/14/09

Date Extracted: 515.4_090313

Analyst: CO

Peer Review:

Analytical Method: 515.4

Herbicides in Drinking Water

CAS	COMPOUND	RESULTS	UNITS	PQL	MDL	MCL	COMMENT	
EPA R	egulated			<u>.</u>				,
4-75-7	2,4 - D	ND	ug/L	0.5	0.2	70		
3-72-1	2,4,5 - TP (\$ILVEX)	ND	ug/L	0.25	0.04	50		
7-86-5	PENTACHLOROPHENOL	ND	ug/L	0.25	0.05	1		
5-99-0	DALAPON	ND	ug/L	3.25	0.9	200		
8-85-7	DINOSEB	ND	ug/L	0.5	0.2	7		
918-02-1	PICLORAM	ND	ug/L	0.25	0.07	500		
Other								
-14028	DCPA (ACID METABOLITES)	ND	ug/L	0.25	0.1			
918-00-9	DICAMBA	ND	ug/L	0.25	0.05			
4-82-6	2,4 DB	ND	u g/ L	2				
3-76-5	2,4,5 T	ND	ug/L	0.25	0.04		•	
5057-89-0	BENTAZON	ND	ug/L	0.5	0.2			
20-36-5	DICHLORPROP	ND	ug/L	0.75	0.5			
0594-66-6	ACIFLUORFEN	ND	ug/L	0.25	0.1			
1-36-5	3,5 - DICHLOROBENZOIC ACID	ND	ug/L	0.25	0.2		•	

ND (Not Detected): indicates that the parameter was not detected above the State Reporting Limit (SRL).

It a compound is detected > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.

MCL (Maximum Contaminant Level) maximum permissible level of a contaminant in water established by EPA; Federal Action Levels are 0.015 mg/L for Lead and 1.3 mg/L for Copper. Sodium has a recommended limit of 20 mg/L. A blank MCL value indicates a level is not currently established.

Trigger Level: DOH Drinking Water Response level. Systems with compounds detected in excess of this level are required to take additional samples. Contact your regional DOH office.



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HERBICIDES IN DRINKING WATER

Client Name: Walla Walla Basin Watershed Council

810 S Main Street

Milton-Freewater, OR 97862

Project::

Field ID: L3

Sample Description: Locher Road 3

Sampled By: Baker

Sample Date: 3/12/09

Source Type: Sampler Phone:

Reference Number: 09-03591

Project: Locher Road Recharge Sites

Lab Number: 07337

Report Date: 4/15/09 Date Analyzed: 03/14/09

Date Extracted: 515.4_090313

Analyst: CO

Peer Review:

Analytical Method: 515.4

Herbicides in Drinking Water

CAS	COMPOUND	RESULTS	UNITS	PQL	MDL	MCL	COMMENT	
EPA R	egulated							
4-75-7	2,4 - D	ND	ug/L	0.5	0.2	70		
3-72-1	2,4,5 - TP (SILVEX)	ND	ug/L	0.25	0.04	50		
7-86-5	PENTACHLOROPHENOL	ND	ug/L	0.25	0.05	1		
5-99-0	DALAPON	ND	ug/L	3.25	0.9	200		
8-85-7	DINOSEB	ND.	ug/L	0.5	0.2	7		
918-02-1	PICLORAM	ND	ug/L	0.25	0.07	500		
Other								
-14028	DCPA (ACID METABOLITES)	ND	ug/L	0.25	0.1			
918- 00-9	DICAMBA	ND	ug/L	0.25	0.05			
4-82-6	2,4 DB	ND	ug/L	2				
3-76-5	2,4,5 T	ND	ug/L	0.25	0.04			
5057-89-0	BENTAZON	ND	ug/L	0.5	0.2			
20-36-5	DICHLORPROP	ND	ug/L	0.75	0.5			
0594-66-6	ACIFLUORFEN	ND	ug/L	0.25	0.1			
1-36-5	3,5 - DICHLOROBENZOIC ACID	ND	ug/L	0.25	0.2			

NOTES:

If a compound is detected > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.

MCL (Maximum Contaminant Level) maximum permissible level of a contaminant in water established by EPA; Federal Action Levels are 0.015 mg/L for Lead and 1.3 mg/L for Copper. Sodium has a recommended limit of 20 mg/L. A blank MCL value indicates a level is not currently established.

Trigger Level: DOH Drinking Water Response level. Systems with compounds detected in excess of this level are required to take additional samples. Contact your regional DOH office.

NO (Not Detected): indicates that the parameter was not detected above the State Reporting Limit (SRL).



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HERBICIDES IN DRINKING WATER

Client Name: Walla Walla Basin Watershed Council

810 S Main Street

Milton-Freewater, OR 97862

Project::

Field ID: IN1

Sample Description: Locher Road In

Sampled By: Baker Sample Date: 3/12/09

Source Type:

Sampler Phone:

Reference Number: 09-03591

Project: Locher Road Recharge Sites

Lab Number: 07338

Report Date: 4/15/09 Date Analyzed: 03/14/09

Date Extracted: 515.4_090313

Analyst: CO

Peer Review:

Analytical Method: 515.4

Herbicides in Drinking Water

CAS	COMPOUND	RESULTS	UNITS	PQL	MDL	MCL	COMMENT
EPA R	egulated						-
4-75-7	2,4 - D	ND	ug/L	0.5	0.2	70	
3-72-1	2,4,5 - TP (\$ILVEX)	ND	ug/L	0.25	0.04	50	
-86-5	PENTACHLOROPHENOL	ND	ug/L	0.25	0.05	1	
-99-0	DALAPON	ND	ug/L	3.25	0.9	200	
3-85-7	DINOSEB	ND	ug/L	0.5	0.2	7	
918-02-1	PICLORAM	ND	ug/L	0.25	0.07	500	
Other	•						
14028	DCPA (ACID METABOLITES)	0.34	ug/L	0.25	0.1		
18-00-9	DICAMBA	ND	ug/L	0.25	0.05		
-82-6	2,4 DB	ND	ug/L	2			qualitative analysis
-76-5	2,4,5 T	ND	ug/L	0.25	0.04		
057-89-0	BENTAZON	ND	ug/L	0.5	0.2		
0-36-5	DICHLORPROP	ND	ug/L	0.75	0.5		
594-66-6	ACIFLUORFEN	ND	ug/L	0.25	0.1		
-36-5	3,5 - DICHLOROBENZOIC ACID	ND	ug/L	0.25	0.2		

INCLES.
If a compound is detected > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.
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If a compound is detected > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.
If a compound is detected > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.
If a compound is detected > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.
If a compound is detected > or = to the State Reporting Level, SRL, specified in the State Reporting Reporting Reporting Reporting Reporting Reporting Report Reporting
ND (Not Detected): indicates that the parameter was not detected above the State Reporting Limit (SRL).



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HERBICIDES IN DRINKING WATER

Client Name: Walla Walla Basin Watershed Council

810 S Main Street

Milton-Freewater, OR 97862

System Name:

System ID Number: DOH Source Number:

Multiple Sources:

Sample Type:

Sample Purpose: Investigative or Other

Sample Location: Locher Road 1

County:

Sampled By: Baker

Sampler Phone:

Reference Number: 09-03591

Project: Locher Road Recharge Sites

Field ID: L1

Lab Number: 046-07335

Date Collected: 3/12/09 00:00 Date Extracted: 515.4_090313

Date Analyzed: 03/14/09

Report Date: 3/19/09

Analyst: CO

Peer Review:

OH#	COMPOUNDS	RESULTS	UNITS	SRL	Trigger	MCL	COMMENT
	EPA Regulated					-	
37	2,4 - D	ND	ug/L	0.5	0.2	70	
38	2,4,5 - TP (SILVEX)	ND	ug/L	1.0	0.4	50	
134	PENTACHLOROPHENOL	ND	ug/L	0.2	0.08	1	
137	DALAPON	ND	ug/L	5	2	200	
139	DINOSEB	ND	ug/L	1.0	0.4	7	
140	PICLORAM	ND	ug/L	0.5	0.2	500	
	Other					-	
138	DICAMBA	ND	ug/L	0.2	0.2		
225	DCPA (ACID METABOLITES)	ND	ug/L	0.1	0.1		
135	2,4 DB	ND	ug/L	1.0	1.0	I	
136	2,4,5 T	ND	ug/L	0.4	0.4		
220	BENTAZON	ND	ug/L	0.5	0.5	İ	
221	DICHLORPROP	ND	ug/L	0.5	0.5		
223	ACIFLUORFEN	ND	ug/L	2.0	2.0		
226	3,5 - DICHLOROBENZOIC ACID	ND	ug/L	0.5	0.5		
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NOTES:

16 a compound is detected > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.

MCL (Maximum Contaminant Level) maximum permissible level of a contaminant in water established by EPA; Federal Action Levels are 0.015 mg/L for Lead and 1.3 mg/L for Copper. Sodium has a recommended limit of 20 mg/L. A blank MCL value indicates a level is not currently established.

Trigger Level; DOH Drinking Water Response level. Systems with compounds detected in excess of this level are required to take additional samples. Contact your regional DOH office.

ND (Not Detected): indicates that the parameter was not detected above the State Reporting Limit (SRL).



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HERBICIDES IN DRINKING WATER

Client Name: Walla Walla Basin Watershed Council

810 S Main Street

Milton-Freewater, OR 97862

System Name:

System ID Number:

DOH Source Number:

Multiple Sources:

Sample Type:

Sample Purpose: Investigative or Other

Sample Location: Locher Road 2

County:

Sampled By: Baker

Sampler Phone:

Reference Number: 09-03591

Project: Locher Road Recharge Sites

Field ID: L2

Lab Number: 046-07336

Date Collected: 3/12/09 00:00

Date Extracted: 515.4 090313

Date Analyzed: 03/14/09

Report Date: 3/19/09

Analyst: CO

Peer Review:

Mathed 545 4 For State Drinking Water Compliance

	COMPOUNDS	RESULTS	UNITS	SRL	Trigger	MCL	COMMENT
	EPA Regulated					<u> </u>	
37	2,4 - D	ND	ug/L	0.5	0.2	70	
38	2,4,5 - TP (SILVEX)	ND	ug/L	1.0	0.4	50	
134	PENTACHLOROPHENOL	ND	ug/L	0.2	0.08	1	
137	DALAPON	ND	ug/L	5	2	200	
139	DINOSEB	ND	ug/L	1.0	0.4	7	
140	PICLORAM	ND	ug/L	0.5	0.2	500	
,	Other						
138	DICAMBA	ND	ug/L	0.2	0.2		
225	DCPA (ACID METABOLITES)	ND	ug/L	0.1	0.1		
135	2,4 DB	ND	ug/L	1.0	1.0		
136	2,4,5 ⊤	ND	ug/L	0.4	0.4	ĺ	
220	BENTAZON	ND	ug/L	0.5	0.5		
221	DICHLORPROP	ND	ug/L	0.5	0.5		
223	ACIFLUORFEN	ND	ug/L	2.0	2.0		
226	3,5 - DICHLOROBENZOIC ACID	ND	ug/L	0.5	0.5		
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NOTES:
If a compound is detected > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.

MCL (Maximum Contaminant Level) maximum permissible level of a contaminant in water established by EPA; Federal Action Levels are 0.015 mg/L for Lead and 1.3 mg/L for Copper. Sodium has a recommended limit of 20 mg/L. A blank MCL value indicates a level is not currently established.

Trigger Level: DOH Drinking Water Response level. Systems with compounds detected in excess of this level are required to take additional samples. Contact your regional DOH office.

ND (Not Detected): indicates that the parameter was not detected above the State Reporting Limit (SRL).



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HERBICIDES IN DRINKING WATER

Client Name: Walla Walla Basin Watershed Council

810 S Main Street

Milton-Freewater, OR 97862

System Name:

System ID Number: DOH Source Number.

Multiple Sources:

Sample Type:

Sample Purpose: Investigative or Other

Sample Location: Locher Road 3

County:

Sampled By: Baker

Sampler Phone:

Reference Number: 09-03591

Project: Locher Road Recharge Sites

Field ID: L3

Lab Number: 046-07337

Date Collected: 3/12/09 00:00 Date Extracted: 515.4_090313

Date Analyzed: 03/14/09 Report Date: 3/19/09

Analyst: CO

Peer Review:

EPA Method 515.4 For State Drinking Water Compliance

DOH#	COMPOUNDS	RESULTS	UNITS	SRL	Trigger	MCL	COMMENT
	EPA Regulated			<u> </u>			
37	2,4 - D	ND	ug/L	0.5	0.2	70	·
38	2,4,5 - TP (SILVEX)	ND.	ug/L	1.0	0.4	50	
134	PENTACHLOROPHENOL	ND	ug/L	0.2	0.08	1	
137	DALAPON	ND	ug/L	5	2	200	
139	DINOSEB	ND	ug/L	1.0	0.4	7	
140	PICLORAM	ND	ug/L	0.5	0.2	500	
	Other						
138	DICAMBA	ND	ug/L	0.2	0.2		
225	DCPA (ACID METABOLITES)	ND	ug/L	0.1	0.1		
135	2,4 DB	ND	ug/L	1.0	1.0		
136	2,4,5 T	ND	ug/L	0.4	0.4		
220	BENTAZON	ND	ug/L	0.5	0.5		
221	DICHLORPROP	ND	ug/L	0.5	0.5		
223	ACIFLUORFEN	ND	ug/L	2.0	2.0	1	
226	3,5 - DICHLOROBENZOIC ACID	ND	ug/L	0.5	0.5		
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NOTES.
If a compound is detected > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.
If a compound is detected > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.
If a compound is detected > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.
If a compound is detected > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.
If a compound is detected > or = to the State Reporting Level, Sequence in t

ND (Not Detected): indicates that the parameter was not detected above the State Reporting Limit (SRL).



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CARBAMATES IN DRINKING WATER

Client Name: Walla Walla Basin Watershed Council

810 S Main Street

Milton-Freewater, OR 97862

System Name:

System ID Number:

DOH Source Number:

Multiple Sources:

Sample Type:

Sample Purpose: Investigative or Other

Sample Location: Locher Road 2

County:

Sampled By: Baker

Sampler Phone:

Reference Number: 09-03591

Project: Locher Road Recharge Sites

Field ID: L2

Lab Number: -07336

Date Collected: 3/12/09 00:00

Date Extracted: 531_090313

Date Analyzed: 03/13/09

Report Date: 3/18/09

Analyst: CO

Peer Review:

EPA Method 531 2 For State Drinking Water Compliance

DOH#	COMPOUNDS	RESULTS	UNITS	SRL	Trigger	MCL	COMMENT
	EPA Regulated		1				
148	OXYMAL.	ND	ug/L	4.0	4.0	200	
146	CARBOFURAN	ND	ug/L	1.8	1.8	40	
	EPA Unregulated		İ				
144	ALDICARB SULFOXIDE	ND	ug/L	1.0	1.0		
143	ALDICARB SULFONE	ND	ug/L	1.6	1.6		
147	METHOMYL	ND	ug/L	1.0	1.0		
141	3-HYDROXYCARBOFURAN	ND	ug/L	2.0	2.0		
142	ALDICARB	ND	ug/L	1.0	1.0		
145	CARBARYL	ND	ug/L	2.0	2.0		
	State Unregulated - Other						
326	PROPOXUR (BAYGON)	ND	ug/L	1.0			
327	METHIOCARB	ND	ug/L	4.0			
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NOTES:
If a compound is detected > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DOH.

MCL (Maximum Contaminant Level) maximum permissible level of a contaminant in water established by EPA; Federal Action Levels are 0.015 mg/L for Lead and 1.3 mg/L for Copper. Sodium has a recommended limit of 20 mg/L. A blank MCL value indicates a level is not currently established.

Trigger Level: DOH Drinking Water Response level. Systems with compounds detected in excess of this level are required to take additional samples. Contact your regional DOH office.

ND (Not Detected): indicates that the parameter was not detected above the State Reporting Limit (SRL).



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CARBAMATES IN DRINKING WATER

Client Name: Walla Walla Basin Watershed Council

810 S Main Street

Milton-Freewater, OR 97862

System Name:

System ID Number:

DOH Source Number:

Multiple Sources:

Sample Type:

Sample Purpose: Investigative or Other

Sample Location: Locher Road 3

County:

Sampled By: Baker

Sampler Phone:

Reference Number: 09-03591

Project: Locher Road Recharge Sites

Field ID: L3

Lab Number: -07337

Date Collected: 3/12/09 00:00

Date Extracted: 531_090313

Date Analyzed: 03/13/09

Report Date: 3/18/09

Analyst: CO

Peer Review:

EPA Method 531 2 For State Drinking Water Compliance

DOH#	COMPOUNDS	RESULTS	UNITS	SRL	Trigger	MCL	COMMENT
····- · · - ·	EPA Regulated			<u> </u>		†	
148	OXYMAL	ND	ug/L	4.0	4.0	200	
146	CARBOFURAN	ND	ug/L	1.8	1.8	40	
	EPA Unregulated						
144	ALDICARB SULFOXIDE	ND	ug/L	1.0	1.0		
143	ALDICARB SULFONE	ND	ug/L	1.6	1.6		
147	METHOMYL.	ND	ug/L	1.0	1.0	l	
141	3-HYDROXYCARBOFURAN	ND	ug/L	2.0	2.0		·
142	ALDICARB	ND	ug/L	1.0	1.0		
145	CARBARYL	ND	ug/L	2.0	2.0		
	State Unregulated - Other	E					
326	PROPOXUR (BAYGON)	ND	ug/L	1.0			
327	METHIOCARB	ND	ug/L	4.0			
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NOTES:

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If a compound is detected > or = to the State Reporting Level, SRL, specified increased monitoring frequencies may occur per DCH.

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If a compound is detected > or = to the State Reporting Level, SRL, specified in excess of the State Reporting Reporting Frequencies may occur per DCH.

If a compound is detected > or = to the State Report PCH.

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If a



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May 8, 2009

Mr. Troy Baker

Walla Walla Basin Watershed Council
810 S Main Street

Milton-Freewater, OR 97862

Dear Mr. Troy Baker,

Your project: Locher Road and Hall-Wei, and Recharge Sites, was received on Friday April 24, 2009.

All samples were analyzed within the accepted holding times, were appropriately preserved and were analyzed according to approved analytical protocols. The quality control data was within laboratory acceptance limits, unless specified in the QA reports.

If you have questions phone me at 800 755-9295.

RE: 09-05772 - Locher Road and Hall-V

Respectfully Submitted,

Lawrence J Henderson, PhD Director of Laboratories Enclosures Data Report



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Data Report

Client Name: Walla Walla Basin Watershed Council

810 S Main Street

Milton-Freewater, OR 97862

Reference Number: 09-05772

Project: Locher Road and Hall-Wentland Red

Report Date: 5/8/09

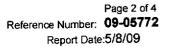
Date Received: 4/24/09

Peer Review:

Lab N	ription: MC1 - Mud Creek 1 umber: 11902							le Date: 4 cted By: \		m	
CAS ID#	Parameter	Result	PQL	MDL	Units	DF	Method	Analyzed	Analysi	t Batch	Comment
E-10139	HYDROGEN ION (pH)	7.07		<u>-</u>	pH Units	1	SM4500-H+ B	4/24/09	MAK I	PH_090424	
E-10617	TURBIDITY	1.04	0.05	0.02	NTU	1	180.1	4/24/09	MAK '	TURB_090424	
14797-55-8	NITRATE-N	1.56	0.100	0.015	mg/L	1	300.0	4/25/09	ВЈ	1098424A	
16887-00-6	CHLORIDE	7.4	20	0.012	mg/L	1	300.0	4/24/09	ВЈ	1090424A	
E-10173	TOTAL DISSOLVED SOLIDS	220	10	6	mg/L	1	SM2540 C	4/28/09	CCN	TDS_090428	
14265-44-2	ORTHO-PHOSPHATE	0.16	0.01	0.002	mg/L	1	SM4500-P F	4/24/09	so	OPHOS-090424	
E-10184	ELECTRICAL CONDUCTIVITY	363	10		uS/cm	1	SM2510 B	4/24/09	CCN	EC_090424	
15541-45-4	BROMATE	ND	0.005	0.00046	mg/L	1	300.1	4/29/09	MVP	D090429A	
E-11778	HARDNESS	149	3.30	0.055	mg CaCO3/L	1	200.7	4/27/09	BJ	200,7-090427A	
E-10117	CHEMICAL OXYGEN DEMAND	ND ·	8.0	2.47	mg/L	1	SM5220 D	5/7/09	MAK	COD_090507	
•	cription: MC2 - Mud Creek 2 lumber: 11903						Colle	ole Date:	Unknov	vr n	
Lab N	·	Result	PQL	MDL	Units	DF			Unknov	vr n	Comment
Lab N CAS ID#	lumber: 11903 Parameter	Result	PQL	MDL	Units pH Units	DF 1	Colle	cted By:	Unknov	vr n	Comment
Lab N CAS ID# E-10139	lumber: 11903		PQL 0.05	MDL 0.02	-		Colle	cted By: Analyzed	Unknov Analys	vn st Batch	Comment
Lab N CAS ID# E-10139 E-10617	Parameter HYDROGEN ION (pH)	6.94			pH Units	1	Method SM4500-H+ B	Analyzed	Unknov Analys MAK	st Batch	Comment
Lab N CAS ID# E-10139	Parameter HYDROGEN ION (pH) TURBIDITY	6.94 0.50	0.05	0.02	pH Units NTU	1	Method SM4500-H+ B	Analyzed 4/24/09 4/24/09	Unknov Analys MAK MAK	vn st Batch PH_090424 TURB_090424	Comment
Lab N CAS ID# E-10139 E-10617 14797-55-8	Parameter HYDROGEN ION (pH) TURBIDITY NITRATE-N	6.94 0.50 0.98	0.05 0.100	0.02 0.015	pH Units NTU mg/L	1 1	Method SM4500-H+ B 180.1 300.0	Analyzed 4/24/09 4/25/09	Analys MAK MAK BJ	t Batch PH_090424 TURB_090424	Comment
Lab N CAS ID# E-10139 E-10617 14797-55-8 16887-00-6	Parameter HYDROGEN ION (pH) TURBIDITY NITRATE-N CHLORIDE	6.94 0.50 0.98 5.3	0.05 0.100 20	0.02 0.015 0.012	pH Units NTU mg/L mg/L	1 1 1 1	Method SM4500-H+ B 180.1 300.0 300.0	Analyzed 4/24/09 4/25/09 4/24/09	Analys MAK MAK BJ BJ	st Batch PH_090424 TURB_090424 1090424A	Comment
Lab N CAS ID# E-10139 E-10617 14797-55-8 16887-00-6 E-10173	Parameter HYDROGEN ION (pH) TURBIDITY NITRATE-N CHLORIDE TOTAL DISSOLVED SOLIDS	6.94 0.50 0.98 5.3 163	0.05 0.100 20 10	0.02 0.015 0.012 6	pH Units NTU mg/L mg/L mg/L	1 1 1 1	Method SM4500-H+ B 180.1 300.0 300.0 SM2540 C	Analyzed 4/24/09 4/24/09 4/25/09 4/28/09	Analys MAK MAK BJ CCN	PH_090424 TURB_090424 1090424A TDS_090428	Comment
Lab N CAS ID# E-10139 E-10617 14797-55-8 16887-00-6 E-10173 14265-44-2	Parameter HYDROGEN ION (pH) TURBIDITY NITRATE-N CHLORIDE TOTAL DISSOLVED SOLIDS ORTHO-PHOSPHATE	6.94 0.50 0.98 5.3 163 0.13	0.05 0.100 20 10 0.01	0.02 0.015 0.012 6	pH Units NTU mg/L mg/L mg/L mg/L	1 1 1 1 1 1 1	Method SM4500-H+ B 180.1 300.0 300.0 SM2540 C SM4500-P F	Analyzed 4/24/09 4/24/09 4/25/09 4/28/09 4/28/09 4/28/09	Analys MAK MAK BJ CCN SO	t Batch PH_090424 TURB_090424 1090424A 1090424A TDS_090428 OPHOS-090424	Comment
Lab N CAS ID# E-10139 E-10617 14797-55-8 16887-00-6 E-10173 14265-44-2 E-10184	Parameter HYDROGEN ION (pH) TURBIDITY NITRATE-N CHLORIDE TOTAL DISSOLVED SOLIDS ORTHO-PHOSPHATE ELECTRICAL CONDUCTIVITY	6.94 0.50 0.98 5.3 163 0.13 275	0.05 0.100 20 10 0.01	0.02 0.015 0.012 6 0.002	pH Units NTU mg/L mg/L mg/L mg/L uS/cm	1 1 1 1 1	Method SM4500-H+ B 180.1 300.0 300.0 SM2540 C SM4500-P F SM2510 B	Analyzed 4/24/09 4/24/09 4/24/09 4/24/09 4/24/09 4/24/09 4/24/09	Analys MAK MAK BJ CCN SO CCN	t Batch PH_090424 TURB_090424 1090424A 1090424A TDS_090428 OPHOS-090424 EC_090424	Comment
Lab N CAS ID# E-10139 E-10617 14797-55-8 16887-00-6 E-10173 14265-44-2 E-10184 15541-45-4	Parameter HYDROGEN ION (pH) TURBIDITY NITRATE-N CHLORIDE TOTAL DISSOLVED SOLIDS ORTHO-PHOSPHATE ELECTRICAL CONDUCTIVITY BROMATE	6.94 0.50 0.98 5.3 163 0.13 275 ND	0.05 0.100 20 10 0.01 10 0.005	0.02 0.015 0.012 6 0.002	pH Units NTU mg/L mg/L mg/L mg/L uS/cm mg/L	1 1 1 1 1 1 1	Collee Method SM4500-H+ B 180.1 300.0 300.0 SM2540 C SM4500-P F SM2510 B 300.1	Analyzed 4/24/09 4/24/09 4/24/09 4/24/09 4/24/09 4/24/09 4/24/09 4/24/09	Analys MAK MAK BJ CCN SO CCN MVP	t Batch PH_090424 TURB_090424 1090424A TDS_090424 OPHOS-090424 EC_090424 D090429A	Comment
Lab N CAS ID# E-10139 E-10617 14797-55-8 16887-00-6 E-10173 14265-44-2 E-10184 15541-45-4 E-11778 E-10117	Parameter HYDROGEN ION (pH) TURBIDITY NITRATE-N CHLORIDE TOTAL DISSOLVED SOLIDS ORTHO-PHOSPHATE ELECTRICAL CONDUCTIVITY BROMATE HARDNESS	6.94 0.50 0.98 5.3 163 0.13 275 ND	0.05 0.100 20 10 0.01 10 0.005 3.30	0.02 0.015 0.012 6 0.002 0.00046 0.055	pH Units NTU mg/L mg/L mg/L uS/cm mg/L uS/cm	1 1 1 1 1 1 1	Collee Method SM4500-H+ B 180.1 300.0 300.0 SM2540 C SM4500-P F SM2510 B 300.1 200.7 SM5220 D	Analyzed 4/24/09 4/24/09 4/25/09 4/24/09 4/24/09 4/24/09 4/24/09 4/24/09 4/29/09 4/27/09	Analys MAK MAK BJ CCN SO CCN MVP BJ MAK	MT Batch PH_090424 TURB_090424 1090424A 1090424A TDS_090428 OPHOS-090424 EC_090424 D090429A 200.7-090427A	Comment

ND = Not detected above the listed practical quantitation limit (PQL) or not above the Method Detection Limit (MDL), if requested.

PQL = Practical Quantitation Limit is the lowest level that can be acheived within specified limits of precision and accuracy during routine laboratory operating conditions.





Data Report

E-10139	HYDROGEN ION (pH)	6.52			pH Units	1	SM4500-H+ B	4/24/09	MAK	PH_090424
E-10617	TURBIDITY	61.8	0.50	0.02	NTU	10	180.1	4/24/09	MAK	TURB_090424
14797-55-8	NITRATE-N	0.22	0.100	0.015	mg/L	1	300.0	4/25/09	BJ	1090424A
16887-00-6	CHLORIDE	0.7	20	0.012	mg/L	1	300.0	4/24/09	BJ	1090424A
E-10173	TOTAL DISSOLVED SOLIDS	52	10	6	mg/L	1	SM2540 C	4/28/09	CCN	TDS_090428
14265-44-2	ORTHO-PHOSPHATE	0.12	0.01	0.005	mg/L	1	SM4500-P F	4/24/09	so	OPHOS-090424
E-10184	ELECTRICAL CONDUCTIVITY	52.9	10		uS/cm	1	SM2510 B	4/24/09	CCN	EC_090424
15541-45-4	BROMATE '	ND	0.005	0.00046	mg/L	1	300.1	4/29/09	MVP	D090429A
E-11778	HARDNESS	20.8	3.30	0.055	mg	1	200.7	4/27/09	ВЈ	200.7-090427A
E-10117	CHEMICAL OXYGEN DEMAND	10	8.0	2.47	CaCO3/L mg/L	1	SM5220 D	5/7/09	MAK	COD_090507

	ription: L2 - Locher Road 2 umber: 11905						Sample Date: 4/23/09 Collected By: Unknown					
CAS ID#	Parameter	Result	PQL	MDL	Units	DF	Method	Analyzed	Analyst	Batch	Comment	
E-10139	HYDROGEN ION (pH)	6.45		-	pH Units	1	SM4500-H+ B	4/24/09	CCN F	PH_090424		
E-10617	TURBIDITY	10.9	0.05	0.02	NTU	1	180.1	4/24/09	MAK "	TURB_090424		
14797-55-8	NITRATE-N	17.3	0.100	0.015	mg/L	1	300.0	4/25/09	вл і	090424A		
16887-00-6	CHLORIDE	9.8	20	0,012	mg/L	1	300.0	4/24/09	BJ I	090424A		
E-10173	TOTAL DISSOLVED SOLIDS	302	10	6	mg/L	1	SM2540 C	4/28/09	CCN	TDS_090428		
14265-44-2	ORTHO-PHOSPHATE	0.20	0.01	0.002	mg/L	1	SM4500-P F	4/24/09	so	OPHOS-090424		
E-10184	ELECTRICAL CONDUCTIVITY	440	10		uS/cm	1	SM2510 B	4/24/09	CCN	EC_090424		
15541-45-4	BROMATE	ND	0.005	0.00046	mg/L	1	300.1	4/29/09	MVP	D090429A		
E-11778	HARDNESS	118	3.30	0.055	mg CaCO3/L	1	200.7	4/27/09	BJ	200.7-090427A		
E-10117	CHEMICAL OXYGEN DEMAND	NĎ	8.0	2.47	mg/L	1	SM5220 D	5/7/09	MAK	COD_090507		

•	ription: L1 - Locher Road 1 umber: 11906							le Date: 4 cted By: U		1	
AS ID#	Parameter	Result	PQL	MDL	Units	DF	Method	Analyzed	Analyst	Batch	Comment
-10139	HYDROGEN ION (pH)	6.71			pH Units	1	SM4500-H+ B	4/24/09	CCN P	H_090424	
-10617	TURBIDITY	15.5	0.05	0.02	NTU	1	180.1	4/24/09	MAK T	URB_090424	
4797-55-8	NITRATE-N	16.2	0.100	0.015	mg/L	1	300.0	4/25/09	BJ I	90424A	
6887-00-6	CHLORIDE	9.7	20	0.012	mg/L	1	300.0	4/24/09	BJ 1	90424A	
-10173	TOTAL DISSOLVED SOLIDS	279	10	6	mg/L	1	SM2540 C	4/28/09	CCN 1	DS_090428	
4265-44-2	ORTHO-PHOSPHATE	0.21	0.01	0.002	mg/L	1	SM4500-P F	4/24/09	\$0 C	PHOS-090424	
-10184	ELECTRICAL CONDUCTIVITY	434	10		uS/cm	1	SM2510 B	4/24/09	CCN E	C_090424	
5541-45-4	BROMATE	ND	0.005	0.00046	mg/L	1	300.1	4/29/09	MVP I	0090429A	
-11778	HARDNESS	167	3.30	0.055	mg CaCO3/L	1	200.7	4/27/09	8J 2	100.7-090427A	
-10117	CHEMICAL OXYGEN DEMAND	ND	8.0	2.47	mg/L	1	SM5220 D	5/7/09	MAK	COD_090507	

Notes:

ND = Not detected above the listed practical quantitation limit (PQL) or not above the Method Detection Limit (MDL), if requested.

PQL = Practical Quantitation Limit is the lowest level that can be acheived within specified limits of precision and accuracy during routine laboratory operating conditions.

D.F. - Dilution Factor



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Page 1 of 4



SAMPLE INDEPENDENT QUALITY CONTROL REPORT

Laboratory Fortified Blank

Reference Number: 09-05772

			True			%		QC	
Batch	Analyte	Result	Value	Units	Method	Recovery	Limits	Qualifier Type*	Comment
200.7-090427A	HARDNESS	73.3	69.5	mg/L	200.7	105	80-120	LFB	
COD_090507	CHEMICAL OXYGEN DEMAND	55	50	mg/L	SM5220 D	110	80-120	LFB	
OPHOS-090424	ORTHO-PHOSPHATE	1.01	1.00	mg/L	SM4500-P F	101	70-130	LFB	
tds_090428	TOTAL DISSOLVED SOLIDS	504	500	mg/L	SM2540 C	101	80-120	LFB	
tds_090428	TOTAL DISSOLVED SOLIDS	512	500	mg/L	SM2540 C	102	80-120	LFB	-

[%] Recovery = (Result of Analysis)/(True Value) * 100

NA = Indicates % Recovery could not be calculated.

QCS: Quality Control Sample, a solution containing known concentrations of method analytes which is used to fortify an aliquot of reagent matrix. The QCS is obtained from an external source and is used to check lab performance.

LFB: Laboratory Fortified Blank, an aliquot of reagent matrix to which known quantities of method analytes are added in the lab. The LFB is analyzed exactly like a sample, and its purpose is to determine whether method performance is within accepted control limits.

MB or LRB. Method Blank or Laboratory Reagent Blank, an aliquot of reagent matrix is analyzed exactly like a sample, and its purpose is to determine if there is background contamination.



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SAMPLE INDEPENDENT QUALITY CONTROL REPORT

Laboratory Reagent Blank

Reference Number: 09-05772

			True			%		QC	
Batch	Analyte	Result	Value	Units	Method	Recovery	Limits	Qualifier Type*	Comment
200.7-090427A	HARDNESS	ND		mg/L	200.7	***	10.0000	(LRB	
COD_090507	CHEMICAL OXYGEN DEMAND	ND		mg/L	SM5220 D		4.00000	LRB	
D090429A	BROMATE	ND.		mg/L	300.1		0.00500	LRB	
109042 4 A	CHLORIDE NITRATE-N	ND ND		mg/L mg/L	300.0 300.0		0.10000 0.10000		
OPHOS-090424	ORTHO-PHOSPHATE	`ND		mg/L	SM4500-P F		0.10000) LRB	
TURB_090424	TURBIDITY	ND		NTU	180.1		0.02000) LRB	

^{*}Notation:

[%] Recovery = (Result of Analysis)/(True Value) * 100

NA = Indicates % Recovery could not be calculated.

QCS: Quality Control Sample, a solution containing known concentrations of method analytes which is used to fortify an aliquot of reagent matrix. The QCS is obtained from an external source and is used to check lab performance.

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MB or LRB: Method Blank or Laboratory Reagent Blank, an aliquot of reagent matrix is analyzed exactly like a sample, and its purpose is to determine if there is background contamination.



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SAMPLE INDEPENDENT QUALITY CONTROL REPORT

Method Blank

Reference Number: 09-05772

			True			%		QC	
Batch	Analyte	Result	Value	Units	Method	Recovery	Limits	Qualifier Type*	Comment
200.7-090427A	HARDNESS	ND		mg/L	200.7		0.82000	МВ	
ec_090424	ELECTRICAL CONDUCTIVITY	ND		uS/cm	SM2510 B		2.50000	MB	•
ec_090424	ELECTRICAL CONDUCTIVITY	ND		u\$/cm	\$M2510 B		2.50000	МВ	
ec_090424	ELECTRICAL CONDUCTIVITY	.ND		uS/cm	SM2510 B		2.50000	МВ	
ec_090424	ELECTRICAL CONDUCTIVITY	ND		uS/cm	SM2510 B		2.50000	MB	
ec_090424	ELECTRICAL CONDUCTIVITY	ND .		uS/cm	SM2510 B		2.50000) MB	
ec_090424	ELECTRICAL CONDUCTIVITY	ND		ц\$/cm	SM2510 B		2.50000) MB	
OPHOS-090424	ORTHO-PHOSPHATE	ND		mg/L	SM4500-P F		0.10000) MB	
tds_090428	TOTAL DISSOLVED SOLIDS	ND		mg/L	SM2540 C		2.50000	мв.	
tds_090428	TOTAL DISSOLVED SOLIDS	ND		mg/L	SM2540 C		2.5000) МВ	
	•								

^{*}Notation:

[%] Recovery = (Result of Analysis)/(True Value) * 100

NA = Indicates % Recovery could not be calculated.

QCS: Quality Control Sample, a solution containing known concentrations of method analytes which is used to fortify an aliquot of reagent matrix. The QCS is obtained from an external source and is used to check lab performance.

LFB: Laboratory Fortified Blank, an aliquot of reagent matrix to which known quantities of method analytes are added in the lab. The LFB is analyzed exactly like a sample, and its purpose is to determine whether method performance is within accepted control limits.

MB or LRB: Method Blank or Laboratory Reagent Blank, an aliquot of reagent matrix is analyzed exactly like a sample, and its purpose is to determine if there is background contamination.



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SAMPLE INDEPENDENT QUALITY CONTROL REPORT

Quality Control Sample

Reference Number: 09-05772

·									
			True			%		QC	
Batch	Analyte	Result	Value	Units	Method	Recovery		QualifierType*	Comment
200.7-090427A	HARDNESS	135	132.3	mg/L	200.7	102	80-120	QCS	
COD_090507	CHEMICAL OXYGEN DEMAND	94	92	mg/L	SM5220 D	102	80-120	QCS	
D090429A	BROMATE	-0.0155	0.0157	mg/L	300.1	99	75-125	qcs	
ec_090424	ELECTRICAL CONDUCTIVITY	155	150.1	uS/cm	SM2510 B	103	80-120	QCS	
ec_090424	ELECTRICAL CONDUCTIVITY	154	150.1	uS/cm	SM2510 B	103	80-120	qcs	•
ec_090424	ELECTRICAL CONDUCTIVITY	154	150.1	u\$/cm	SM2510 B	103	80-120	QCS	
ec_090424	ELECTRICAL CONDUCTIVITY	150	150.1	uS/cm	SM2510 B	100	80-120	QCS	
ec_090424	ELECTRICAL CONDUCTIVITY	152	150.1	uS/cm	SM2510 B	101	80-120		
ec_090424	ELECTRICAL CONDUCTIVITY	154	. 150.1	u\$/cm	SM2510 B	103	80-120	QCS	
1090 424 A	CHLORIDE	·29	30.0	mg/L	300.0	97	80-120		
	NITRATE-N	2.42	2.50	mg/L	300.0	97	80-120)	
OPHQS-090424	ORTHO-PHOSPHATE	0.45	0.49	mg/L	SM4500-P F	92	70-130	g QCS	
TURB_090424	TURBIDITY	0.95	1.00	NTU	180.1	95	70-13	o qcs	

[%] Recovery = (Result of Analysis)/(True Value) * 100

NA = Indicates % Recovery could not be calculated.

QCS: Quality Control Sample, a solution containing known concentrations of method analytes which is used to fortify an aliquot of reagent matrix. The QCS is obtained from an external source and is used to check lab performance.

LFB: Laboratory Fortified Blank, an aliquot of reagent matrix to which known quantities of method analytes are added in the lab. The LFB is analyzed exactly like a sample, and its purpose is to determine whether method performance is within accepted control limits.

MB or LRB: Method Blank or Laboratory Reagent Blank, an aliquot of reagent matrix is analyzed exactly like a sample, and its purpose is to determine if there is background contamination.



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QUALITY CONTROL REPORT

Duplicate and Matrix Spike/Matrix Spike Duplicate Report

Reference Number: 09-05772 Report Date: 5/8/2009

Duplicate	ø								
				Duplicate	:	9	į	OC Ouslifer	Comments
Batch	Sample Analyte	nalyte	Result	Result	Units		١		
200.7-090427A	L	11461 HARDNESS 11780 HARDNESS	142	143 155	mg CaCO3/L mg CaCO3/L	0.0	0.45	DUP	
COD_090507	12748 CH	12748 CHEMICAL OXYGEN DEMAND	3600	3500	πg/L	2.8	9-45	OUP	
D090429A									
EC_090424		VENTALISMOOT A CHERCA	536	367	uS/em	0.3	045	DUP	
	1146U EL	ELECTRICAL CONDUCTIVITY	622	780	uS/cm	0.4	0-45	DUP	
		ELECTRICAL CONDUCTIVITY	386	401	m2/cm	3.8	045	DUP	
		ELECTRICAL CONDUCTIVITY	275	272	uS/cm	77	0-45	DUP	
		ELECTRICAL CONDUCTIVITY	367	362	uS/cm	<u>4</u>	54.	DUP	
1090424A					:	Š	97.0	2	
	11797 C	11797 CHLORIDE	9.0	9:0	mg/L	2 6	n (S =	
	11864 C	CHLORIDE	89	18	mg/L	0.0	£ ;	5 2	
	11934 C	CHLORIDE	15	15	mg/L	0.0	0-45	900 1111	
		CHLORIDE	7.4	7.4	mg/L	0.0	0.45	400	
OPHOS-090424		11910 ORTHO-PHOSPHATE	0.16	0.16	T/BW	0.0	0-20	DUP	
PH_090424	11911 H	11911 HYDROGEN ION (pH)	6.05	6.03	pH Units	0.3	0.45	DUP	
TDS_090428 TURB_090424	4	11911 TURBIDITY 11934 TURBIDITY	284 0.68	280 0.71	UTN	4.1	0-50	and Pud	

%RPD = Relative Percent Difference

NA = Indicates %RPD could not be calculated

Matrix Spike (MS)/Matrix Spike Duplicate (MSD) analyses are used to determine the accuracy (MS) and precision (MSD) of a analytical method in a given sample matrix. Therefore, the usefulness of this report is limited to samples of similar matrices analyzed in the same analytical batch.

Only Duplicate sample with detections are listed in this report



Reference Number: 09-05772 Report Date: 5/8/2009 Page 2 of 2

Matrix Spike	oike			Duplicate										-
•			Spike	Spike	Spike		Percent Recovery	COVETY				သွ		
Batch	Sample Analyte	Result	Result	Result	Conc	Units	MS N	MSD	Limits	%RPD	Limits	Qualifier	Comments	
200.7-090427A				;	;	0		6	20	4	08.0	2	5	
	11461 HARDNESS	142	210	211	69.5	mg cacos/r sa			021-06	<u> </u>	3	<u>.</u>		
	11780 HARDNESS	155	225	223	69.5	mg CaCO3/L 101		86	80-120	2.9	09-0	LFM	5	
COD 090507														
	11607 CHEMICAL OXYGEN DEMAND	Q X	45	48	20	mg/L	6 06	96	80-120	6.5	0-0	E F	Σ	
	42242 CHEMICAL OXYGEN DEMAND	9	64	49	20	mg/L	86	86	80-120	0.0	09-0	LFM	Σ	
	12748 CHEMICAL OXYGEN DEMAND	3600	2900	2900	2500	mg/L	92 6	95	80-120	0.0	09-0	LFIM	Σ	
D090429A		,						•	75.175	4	0-60	FM	5	
	11629 BROMATE	9	0.0087		0.070	J/6E			2	•	2			
	11909 BROMATE	2	0.0107		0.010	mg/L	107	¥ Z	75-125	₹ Z	0-0	F.	⋝	,
1090424A		,				•			9	•	٠ .	710		
	11797 CHLORIDE	9.0	20.1		20.00	mg/L	88	۲ ۲	80-120	ď Z	0-00	5	-	
	44064 NITDATEN	Q	1.08		1.00	mg/L	108	Y.	80-120	¥	0-60	LFM	Σ	
	N-H-CALL TOOK	S	1 05		1,00	mo/L	105	¥	80-120	Ą	09-0	LFM	2	
	1.954 NITANIE-N	9	18.8		20.00	mg/L	94	¥	80-120	¥	09-0	LFM	₽	
	N-U-XXIIN Z#BLE	1				,			0,00	1	6	Nu -	5	
	11942 CHLORIDE	7,4	26.3		20.00	mg/L	96		80-120	⊈ Z	ng-n	5		
OPHOS-090424	24								:	,		_		
1	11910 ORTHO-PHOSPHATE	0.16	1.17	1.19	1.00	mg/L	<u>1</u>	103	70-130	2.0	0-20	E L		

%RPD = Relative Percent Difference

NA = Indicates %RPD could not be calculated
Matrix Spike (MS)/Matrix Spike Duplicate (MSD) analyses are used to determine the accuracy (MS) and precision (MSD) of a analytical method in a given sample matrix. Therefore, the usefulness of this report is limited to samples of similar matrices analyzed in the same analytical batch.

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Data Report

Client Name: Walla Walla Basin Watershed Council

810 S Main Street

Milton-Freewater, OR 97862

Reference Number: 09-07745

Project: Locher Road Recharge Sites and Ha

Report Date: 6/11/09

Date Received: 5/29/09

Collected By: Unknown

Analyzed Analyst Batch

Comment

Method

Peer Review:

CAS ID#	Parameter	Result	PQL	MDL	Units	DF	Method	Analyzed	Analys	st Batch	Comment
E-10139	HYDROGEN ION (pH)	6.89			pH Units	1	\$M4500-H+ B	5/29/09	MAK	PH_090529	
E-10617	TURBIDITY	1.30	0.05	0.02	NTU	1	180.1	5/29/09	MAK	TURB_090529	
14797-55-8	NITRATE-N	7.07	0.100	0.015	mg/L	1	300.0	5/29/09	ВЈ	1090529A	
16887-00-6	CHLORIDE	7.26	0.1	0.012	mg/L	1	300.0	5/29/09	вј	1090529A	
E-10173	TOTAL DISSOLVED SOLIDS	266	10		mg/L	1	SM2540 C	6/3/09	CCN	TDS_090603	
14265-44-2	ORTHO-PHOSPHATE	0.27	0.01	0.002	mg/L	1	SM4500-P F	5/29/09	so	OPHOS-090529B	
E-10184	ELECTRICAL CONDUCTIVITY	431	10		uS/cm	1	SM2510 B	6/3/09	CCN	EC_090603	
15541-45-4	BROMATE	ND	0.005	0.0016	mg/L	ı	300.1	6/1/09	MVP	D090601A	
E-11778	HARDNESS	172	3.30	0.055	mg CaCO3/L	1	200.7	6/1/09	BJ	200.7-090601B	. :-
					CGCCGE						
Sample Desc	chemical oxygen demand ription: L-2 - Locher Rd. 2 Well umber: 16096	ND	8.0	2.47	mg/L	1	-	6/2/09 ble Date: ! cted By:			
Sample Desc Lab N	ription: L-2 - Locher Rd. 2 Well	ND Result	8.0	2.47 MDL		1 DF	Samp	ole Date:	5/28/09 Unknov) wn	Comment
Sample Desc Lab N CAS ID#	ription: L-2 - Locher Rd. 2 Well umber: 16096				mg/L		Samp Colle	ole Date:	5/28/09 Unknov) wn	Comment
Sample Desc Lab N CAS ID# E-10139	ription: L-2 - Locher Rd. 2 Well umber: 16096 Parameter	Result			mg/L Units	DF	Samp Colle Method	ole Date: ! cted By: ! Analyzed	5/28/09 Unknov Analys	wn st Batch	Comment
Sample Desc Lab N CAS ID# E-10139 E-10617	ription: L-2 - Locher Rd. 2 Well umber: 16096 Parameter HYDROGEN ION (pH)	Result 6.69	PQL	MDL	mg/L Units pH Units	DF 1	Samp Colle Method SM4500-H+ B	ole Date: steed By: Analyzed	5/28/09 Unknov Analys	wn st Batch PH_090529	Comment
Sample Desc Lab N CAS ID# E-10139 E-10617 14797-55-8	ription: L-2 - Locher Rd. 2 Well umber: 16096 Parameter HYDROGEN ION (pH) TURBIDITY	Result 6.69 9.13	PQL 0.05	MDL 0.02	Units PH Units NTU	DF 1	Samp Colle Method SM4500-H+ B 180.1	ole Date: 4 cted By: 4 Analyzed 5/29/09	5/28/09 Unknov Arialys MAK MAK	Nn st Batch PH_090529 TURB_090529	Comment
Sample Desc Lab N CAS ID# E-10139 E-10617 14797-55-8 16887-00-6	ription: L-2 - Locher Rd. 2 Well umber: 16096 Parameter HYDROGEN ION (pH) TURBIDITY NITRATE-N	Result 6.69 9.13 15.4	PQL 0.05 0.100	MDL 0.02 0.015	Units PH Units NTU mg/L	DF 1 1 1	Samp Colle Method SM4500-H+ B 180.1 300.0	Analyzed 5/29/09 5/30/09	5/28/09 Unknov Analys MAK MAK BJ	MT St Batch PH_090529 TURB_090529 K090529A	Comment
Sample Desc Lab N CAS ID# E-10139 E-10617 14797-55-8 16887-00-6 E-10173	ription: L-2 - Locher Rd. 2 Well umber: 16096 Parameter HYDROGEN ION (pH) TURBIDITY NITRATE-N CHLORIDE	Result 6.69 9.13 15.4 8.9	PQL 0.05 0.100 0.1	MDL 0.02 0.015	Units PH Units NTU mg/L mg/L	DF 1 1 1 1	Samp Colle Method SM4500-H+ B 180.1 300.0 300.0	Analyzed 5/29/09 5/29/09	5/28/09 Unknov Analys MAK MAK BJ BJ	NT st Batch PH_090529 TURB_090529 1090529A	
Sample Desc Lab N CAS ID# E-10139 E-10617 14797-55-8 16887-00-6 E-10173 14265-44-2	ription: L-2 - Locher Rd. 2 Well umber: 16096 Parameter HYDROGEN ION (pH) TURBIDITY NITRATE-N CHLORIDE TOTAL DISSOLVED SOLIDS	Result 6.69 9.13 15.4 8.9 286	PQL 0.05 0.100 0.1	MDL 0.02 0.015 0.012	Units PH Units NTU mg/L mg/L	DF 1 1 1 1 1	Samp Colle Method SM4500-H+ B 180.1 300.0 300.0 SM2540 C	Analyzed 5/29/09 5/29/09 5/29/09 5/29/09 5/29/09 5/29/09	5/28/09 Unknov Analys MAK MAK BJ BJ CCN	Nn St Batch PH_090529 TURB_090529 1090529A TDS_090603	
Sample Desc Lab N CAS ID# E-10139 E-10617 14797-55-8 16687-00-6 E-10173 14265-44-2 E-10184	ription: L-2 - Locher Rd. 2 Well umber: 16096 Parameter HYDROGEN ION (pH) TURBIDITY NITRATE-N CHLORIDE TOTAL DISSOLVED SOLIDS ORTHO-PHOSPHATE	Result 6.69 9.13 15.4 8.9 286 0.26	PQL 0.05 0.100 0.1 10 0.01	MDL 0.02 0.015 0.012	Units PH Units NTU mg/L mg/L mg/L	DF 1 1 1 1 1 1 1	Samp Colle Method SM4500-H+ B 180.1 300.0 300.0 SM2540 C SM4500-P F	Analyzed 5/29/09 5/29/09 5/29/09 5/29/09 5/29/09	Arialys MAK MAK BJ GCN SO	NT St Batch PH_090529 TURB_090529 1090529A 1090529A TDS_090603 OPHOS-090529B	
•	ription: L-2 - Locher Rd. 2 Well umber: 16096 Parameter HYDROGEN ION (pH) TURBIDITY NITRATE-N CHLORIDE TOTAL DISSOLVED SOLIDS ORTHO-PHOSPHATE ELECTRICAL CONDUCTIVITY	Result 6.69 9.13 15.4 8.9 286 0.26 422	PQL 0.05 0.100 0.1 10 0.01	MDL 0.02 0.015 0.012 0.002	Units PH Units NTU mg/L mg/L mg/L mg/L mg/L mg/L	DF 1 1 1 1 1 1 1 1 1	Samp Colle Method SM4500-H+ B 180.1 300.0 300.0 SM2540 C SM4500-P F SM2510 B	Analyzed 5/29/09 5/29/09 5/29/09 5/29/09 5/29/09 6/3/09	S/28/09 Arialys MAK MAK BJ GGN SO GCN	MTN st Batch PH_090529 TURB_090529 1090529A TDS_090603 OPHOS-090529B EC_090603	

Notes:

CAS ID#

ND = Not detected above the listed practical quantitation limit (PQL) or not above the Method Detection Limit (MDL), if requested.

PQL = Practical Quantitation Limit is the lowest level that can be acheived within specified limits of precision and accuracy during routine laboratory operating conditions.

Result

PQL

MDL

Units

D.F. - Dilution Factor

Lab Number: 16097

Parameter





HARDNESS

CHEMICAL OXYGEN DEMAND

Data Report

E-10139	HYDROGEN ION (pH)	6.80			ρH Units	1	SM4500-H+ B	5/29/09	MAK	PH_090529
E-10617	TURBIDITY	6.88	0.05	0.02	NTU	1	180.1	5/29/09	MAK	TUR8_090529
14797-55-8	NITRATE-N	1.84	0.100	0.015	mg/L	1	300.0	5/30/09	BJ	1090529A
16887-00-6	CHLORIDE	1.42	0.1	0.012	mg/L	1	300.0	5/29/09	вј	1090529A
E-10173	TOTAL DISSOLVED SOLIDS	96	10		mg/L	1	SM2540 C	6/3/09	CCN	TD\$_090603
14265-44-2	ORTHO-PHOSPHATE	ND	0.01	0.002	mg/L	1	SM4500-P F	5/29/09	so	OPHO\$-090529B
E-10184	ELECTRICAL CONDUCTIVITY	121	10		uS/cm	1	SM2510 B	6/3/09	CCN	EC_090603
15541-45-4	BROMATE	ND	0.005	0.0016	mg/L	1	300.1	6/1/09	MVP	D090601A
E-11778	HARDNESS	44.0	3.30	0.055	mg CaCO3/L	1	200.7	6/1/09	BJ	200.7-0906018
E-10117	CHEMICAL OXYGEN DEMAND	ND	8.0	2.47	mg/L	1	SM5220 D	6/2/09	MAK	COD_090602

•	cription: SW-1 - Locher Rd. SW 1 cumber: 16098						•	ole Date: 5 cted By: \			
CAS ID#	Parameter	Result	PQL	MDL	Units	DF	Method	Analyzed	Analys	st Batch	Comment
E-10139	HYDROGEN ION (pH)	7.18			pH Units	1	SM4500-H+ B	5/29/09	MAK	PH_090529	
E-10617	TURBIDITY	3.97	0.05	0.02	NTU	1	180.1	5/29/09	MAK	TURB_090529	
14797-55-8	NITRATE-N	1.37	0.100	0.015	mg/L	1	300.0	5/30/09	BJ	1090529A	
16887-00-6	CHLORIDE	6.8	0,1	0.012	mg/L	1	300.0	5/29/09	ВЈ	1090529A	
E-10173	TOTAL DISSOLVED SOLIDS	226	10		mg/L	1	SM2540 C	6/3/09	CCN	TD\$_090603	
14265-44-2	ORTHO-PHOSPHATE	0.26	0.01	0.002	mg/L	1	SM4500-P F	5/29/09	50	OPHOS-090529B	
E-101 84	ELECTRICAL CONDUCTIVITY	374	10		u\$/cm	1	SM2510 B	6/3/09	CCN	EC_090603	
15541-45-4	BROMATE	ND	0.005	0.0016	mg/L	1	300.1	6/1/09	MVP	D090601A	

0.055

2.47

mg CaCO3/L

SM5220 D

COD_090602

mg/L

•	ription: SW-2 - Locher Rd. SW 2 umber: 16099					_	•	ole Date: (cted By:			
CAS ID#	Parameter	Result	PQL	MDL	Units	ÐF	Method	Analyzed	Analy	st Batch	Comment
E-10139	HYDROGEN ION (pH)	7.25			pH Units	1	SM4500-H+ B	5/29/09	MAK	PH_090529	
E-10617	TURBIDITY	0.96	0.05	0.02	NTU	ı	180.1	5/29/09	MAK	TURB_090529	
14797-55-8	NITRATE-N	1.38	0.100	0.015	mg/L	1	300.0	5/30/09	BJ	1090529A	
16887-00-6	CHLORIDE	6.2	0.1	0.012	mg/L	1	300.0	5/29/09	BJ	1090529A	
E-10173	TOTAL DISSOLVED SOLIDS	198	10		mg/L	1	SM2540 C	6/3/09	CCN	TD\$_090603	
14265-44-2	ORTHO-PHOSPHATE	0.66	0.01	0.002	mg/L	1	SM4500-P F	5/29/09	so	OPHO\$-090529B	
E-10184	ELECTRICAL CONDUCTIVITY	310	10		uS/cm	1	SM2510 B	6/3/09	CCN	EC_090603	
15541-45-4	BROMATE	ND	0.005	0.0016	m g/ L	1	300.1	6/1/09	MVP	D090601A	
E-11 778	HARDNESS	126	3.30	0.055	mg CaCO3/L	1	200.7	6/1/09	BJ	200.7-090601B	
E-10117	CHEMICAL OXYGEN DEMAND	ND	8.0	2.47	mg/L	1	SM5220 D	6/2/09	MAK	COD_090602	

E-11778

E-10117

142

ND

3.30

8.0

ND = Not detected above the listed practical quantitation limit (PQL) or not above the Method Detection Limit (MDL), if requested.

PQL = Practical Quantitation Limit is the lowest level that can be acheived within specified limits of precision and accuracy during routine laboratory operating conditions.

D.F. - Dilution Factor



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Reference Number: 09-07745 Report Date:6/11/09

Data Report

Sample Description: L-INT - Locher Rd. Intake

Sample Date: 5/28/09

Lab N	umber: 16100						Colle	cted By: U	Jnknov	₩n	
CAS ID#	Parameter	Result	PQL	MDL	Units	DF	Method	Analyzed	Analys	st Batch	Comment
E-10139	HYDROGEN ION (pH)	6.89			pH Units	1	SM4500-H+ B	5/29/09	MAK	PH_090529	
E-10617	TURBIDITY	11.9	0.05	0.02	NTU	1	180.1	5/29/09	MAK	TURB_090529	
14797-55-8	NITRATE-N	0.25	0.100	0.015	mg/L	1	300.0	5/30/09	ВJ	1090529A	
16887-00-6	CHLORIDE	0.74	0,1	0.012	mg/L	1	300.0	5/29/09	BJ	1090529A	
E-10173	TOTAL DISSOLVED SOLIDS	61	10		mg/L	1	SM2540 C	6/3/09	CCN	TDS_090603	
14265-44-2	ORTHO-PHOSPHATE	0.14	0.01	0.002	mg/L	1	SM4500-P F	5/29/09	\$0	OPHOS-090529B	
E-10184	ELECTRICAL CONDUCTIVITY	66. 4	10		uS/cm	1	SM2510 B	6/3/09	CCN	EC_090603	
15541-45-4	BROMATE	ND	0.005	0.0016	mg/L	1	300.1	6/9/09	MVP	D090609A	
E-11778	HARDNESS	25.2	3.30	0.055	mg CaCO3/L	1	200.7	6/1/09	BJ	200.7-090601B	
E-10117	CHEMICAL OXYGEN DEMAND	ND	8.0	2.47	mg/L	1	SM5220 D	6/2/09	MAK	COD_090602	