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COVER LETTER

September 28, 2004

Cindy Hurtado  
San Juan Refining  
#50 CR 4990  
Bloomfield, NM 87413  
TEL: (505) 632-4161  
FAX (505) 632-3911

RE: Inj Well 3rd Qtr 2004

Order No.: 0409046

Dear Cindy Hurtado:

Hall Environmental Analysis Laboratory received 2 samples on 9/3/2004 for the analyses presented in the following report.

These were analyzed according to EPA procedures or equivalent.

Reporting limits are determined by EPA methodology. No determination of compounds below these (denoted by the ND or < sign) has been made.

Please don't hesitate to contact HEAL for any additional information or clarifications.

Sincerely,



Andy Freeman, Business Manager  
Nancy McDuffie, Laboratory Manager



**Hall Environmental Analysis Laboratory**

Date: 28-Sep-04

CLIENT: San Juan Refining  
Project: Inj Well 3rd Qtr 2004  
Lab Order: 0409046

**CASE NARRATIVE**

METHOD: 8270C

HOLDING TIMES: All holding times for Preparation and Analysis were met.

METHOD: Preparations: 3510  
Analysis: 8270C

PREPARATION: Sample preparation proceeded normally.

**ANALYSIS:**

1. Calibration: All acceptance criteria were met.
2. Blanks: The initial blank was spiked along with the first 3 samples. Results for these are reported from the re-extraction.
3. Internal Standards: Due to an incorrect internal standard batch surrogates failed in the initial extraction batch 6442
4. Surrogates: Surrogates were reported from QC batch 6493.
5. Spikes: Initial LCS/LCSD failed due to bad internal standards.
6. Samples: Samples were re-extracted out of holdtime to confirm internal standards. All target analytes are non detect and results from both extraction batches are used.

Andy with Hall called -  
our 8270 sample was inadvertently spiked - the lab tech  
thought it was a blank.  
8270 was re-extracted but out of hold time

# Hall Environmental Analysis Laboratory

Date: 28-Sep-04

CLIENT: San Juan Refining  
 Lab Order: 0409046  
 Project: Inj Well 3rd Qtr 2004  
 Lab ID: 0409046-01

Client Sample ID: Injection Well 3rd  
 Collection Date: 9/2/2004 9:20:00 AM  
 Matrix: AQUEOUS

| Analyses                            | Result | PQL | Qual | Units      | DF | Date Analyzed        |
|-------------------------------------|--------|-----|------|------------|----|----------------------|
| <b>EPA METHOD 300.0: ANIONS</b>     |        |     |      |            |    | Analyst: MAP         |
| Chloride                            | 1200   | 5.0 |      | mg/L       | 50 | 9/15/2004 2:49:36 PM |
| Sulfate                             | 750    | 25  |      | mg/L       | 50 | 9/15/2004 2:49:36 PM |
| <b>EPA METHOD 310.1: ALKALINITY</b> |        |     |      |            |    | Analyst: CMC         |
| Alkalinity, Total (As CaCO3)        | 520    | 4.0 |      | mg/L CaCO3 | 2  | 9/10/2004            |
| Carbonate                           | 16     | 4.0 |      | mg/L CaCO3 | 2  | 9/10/2004            |
| Bicarbonate                         | 510    | 4.0 |      | mg/L CaCO3 | 2  | 9/10/2004            |
| <b>EPA METHOD 8260B: VOLATILES</b>  |        |     |      |            |    | Analyst: BDH         |
| Benzene                             | ND     | 10  |      | µg/L       | 10 | 9/8/2004             |
| Toluene                             | ND     | 10  |      | µg/L       | 10 | 9/8/2004             |
| Ethylbenzene                        | ND     | 10  |      | µg/L       | 10 | 9/8/2004             |
| Methyl tert-butyl ether (MTBE)      | 1300   | 10  |      | µg/L       | 10 | 9/8/2004             |
| 1,2,4-Trimethylbenzene              | ND     | 10  |      | µg/L       | 10 | 9/8/2004             |
| 1,3,5-Trimethylbenzene              | 17     | 10  |      | µg/L       | 10 | 9/8/2004             |
| 1,2-Dichloroethane (EDC)            | ND     | 10  |      | µg/L       | 10 | 9/8/2004             |
| 1,2-Dibromoethane (EDB)             | ND     | 10  |      | µg/L       | 10 | 9/8/2004             |
| Naphthalene                         | ND     | 20  |      | µg/L       | 10 | 9/8/2004             |
| 1-Methylnaphthalene                 | ND     | 40  |      | µg/L       | 10 | 9/8/2004             |
| 2-Methylnaphthalene                 | ND     | 40  |      | µg/L       | 10 | 9/8/2004             |
| Acetone                             | ND     | 100 |      | µg/L       | 10 | 9/8/2004             |
| Bromobenzene                        | ND     | 10  |      | µg/L       | 10 | 9/8/2004             |
| Bromochloromethane                  | ND     | 10  |      | µg/L       | 10 | 9/8/2004             |
| Bromodichloromethane                | ND     | 10  |      | µg/L       | 10 | 9/8/2004             |
| Bromoform                           | ND     | 10  |      | µg/L       | 10 | 9/8/2004             |
| Bromomethane                        | ND     | 20  |      | µg/L       | 10 | 9/8/2004             |
| 2-Butanone                          | ND     | 100 |      | µg/L       | 10 | 9/8/2004             |
| Carbon disulfide                    | ND     | 100 |      | µg/L       | 10 | 9/8/2004             |
| Carbon Tetrachloride                | ND     | 10  |      | µg/L       | 10 | 9/8/2004             |
| Chlorobenzene                       | ND     | 10  |      | µg/L       | 10 | 9/8/2004             |
| Chloroethane                        | ND     | 20  |      | µg/L       | 10 | 9/8/2004             |
| Chloroform                          | ND     | 10  |      | µg/L       | 10 | 9/8/2004             |
| Chloromethane                       | ND     | 10  |      | µg/L       | 10 | 9/8/2004             |
| 2-Chlorotoluene                     | ND     | 10  |      | µg/L       | 10 | 9/8/2004             |
| 4-Chlorotoluene                     | ND     | 10  |      | µg/L       | 10 | 9/8/2004             |
| cis-1,2-DCE                         | ND     | 10  |      | µg/L       | 10 | 9/8/2004             |
| cis-1,3-Dichloropropene             | ND     | 10  |      | µg/L       | 10 | 9/8/2004             |
| 1,2-Dibromo-3-chloropropane         | ND     | 20  |      | µg/L       | 10 | 9/8/2004             |
| Dibromochloromethane                | ND     | 10  |      | µg/L       | 10 | 9/8/2004             |
| Dibromomethane                      | ND     | 20  |      | µg/L       | 10 | 9/8/2004             |
| 1,2-Dichlorobenzene                 | ND     | 10  |      | µg/L       | 10 | 9/8/2004             |
| 1,3-Dichlorobenzene                 | ND     | 10  |      | µg/L       | 10 | 9/8/2004             |

Qualifiers: ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits  
 J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits  
 B - Analyte detected in the associated Method Blank      E - Value above quantitation range  
 \* - Value exceeds Maximum Contaminant Level

# Hall Environmental Analysis Laboratory

Date: 28-Sep-04

CLIENT: San Juan Refining  
 Lab Order: 0409046  
 Project: Inj Well 3rd Qtr 2004  
 Lab ID: 0409046-01

Client Sample ID: Injection Well 3rd  
 Collection Date: 9/2/2004 9:20:00 AM

Matrix: AQUEOUS

| Analyses                    | Result | PQL      | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|----------|------|-------|----|---------------|
| 1,4-Dichlorobenzene         | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| Dichlorodifluoromethane     | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| 1,1-Dichloroethane          | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| 1,1-Dichloroethene          | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| 1,2-Dichloropropane         | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| 1,3-Dichloropropane         | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| 2,2-Dichloropropane         | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| 1,1-Dichloropropene         | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| Hexachlorobutadiene         | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| 2-Hexanone                  | ND     | 100      |      | µg/L  | 10 | 9/8/2004      |
| Isopropylbenzene            | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| 4-Isopropyltoluene          | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| 4-Methyl-2-pentanone        | ND     | 100      |      | µg/L  | 10 | 9/8/2004      |
| Methylene Chloride          | ND     | 30       |      | µg/L  | 10 | 9/8/2004      |
| n-Butylbenzene              | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| n-Propylbenzene             | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| sec-Butylbenzene            | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| Styrene                     | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| tert-Butylbenzene           | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| 1,1,1,2-Tetrachloroethane   | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| 1,1,1,2,2-Tetrachloroethane | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| Tetrachloroethene (PCE)     | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| trans-1,2-DCE               | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| trans-1,3-Dichloropropene   | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| 1,2,3-Trichlorobenzene      | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| 1,2,4-Trichlorobenzene      | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| 1,1,1-Trichloroethane       | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| 1,1,2-Trichloroethane       | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| Trichloroethene (TCE)       | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| Trichlorofluoromethane      | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| 1,2,3-Trichloropropane      | ND     | 20       |      | µg/L  | 10 | 9/8/2004      |
| Vinyl chloride              | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| Xylenes, Total              | ND     | 10       |      | µg/L  | 10 | 9/8/2004      |
| Surr: 1,2-Dichloroethane-d4 | 93.1   | 70.6-124 |      | %REC  | 10 | 9/8/2004      |
| Surr: 4-Bromofluorobenzene  | 102    | 76.4-130 |      | %REC  | 10 | 9/8/2004      |
| Surr: Dibromofluoromethane  | 91.6   | 67.2-131 |      | %REC  | 10 | 9/8/2004      |
| Surr: Toluene-d8            | 99.2   | 82.1-123 |      | %REC  | 10 | 9/8/2004      |

## EPA METHOD 8270C: SEMIVOLATILES

Analyst: GAB

|                |    |    |  |      |   |           |
|----------------|----|----|--|------|---|-----------|
| Acenaphthene   | ND | 50 |  | µg/L | 1 | 9/18/2004 |
| Acenaphthylene | ND | 50 |  | µg/L | 1 | 9/18/2004 |
| Aniline        | ND | 50 |  | µg/L | 1 | 9/18/2004 |
| Anthracene     | ND | 50 |  | µg/L | 1 | 9/23/2004 |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits

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# Hall Environmental Analysis Laboratory

Date: 28-Sep-04

CLIENT: San Juan Refining  
 Lab Order: 0409046  
 Project: Inj Well 3rd Qtr 2004  
 Lab ID: 0409046-01

Client Sample ID: Injection Well 3rd  
 Collection Date: 9/2/2004 9:20:00 AM  
 Matrix: AQUEOUS

| Analyses                    | Result | PQL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|-----|------|-------|----|---------------|
| Azobenzene                  | ND     | 50  |      | µg/L  | 1  | 9/18/2004     |
| Benz(a)anthracene           | ND     | 75  |      | µg/L  | 1  | 9/23/2004     |
| Benzo(a)pyrene              | ND     | 50  |      | µg/L  | 1  | 9/23/2004     |
| Benzo(b)fluoranthene        | ND     | 50  |      | µg/L  | 1  | 9/23/2004     |
| Benzo(g,h,i)perylene        | ND     | 50  |      | µg/L  | 1  | 9/18/2004     |
| Benzo(k)fluoranthene        | ND     | 50  |      | µg/L  | 1  | 9/23/2004     |
| Benzoic acid                | ND     | 250 |      | µg/L  | 1  | 9/18/2004     |
| Benzyl alcohol              | ND     | 100 |      | µg/L  | 1  | 9/18/2004     |
| Bis(2-chloroethoxy)methane  | ND     | 50  |      | µg/L  | 1  | 9/18/2004     |
| Bis(2-chloroethyl)ether     | ND     | 75  |      | µg/L  | 1  | 9/18/2004     |
| Bis(2-chloroisopropyl)ether | ND     | 75  |      | µg/L  | 1  | 9/18/2004     |
| Bis(2-ethylhexyl)phthalate  | ND     | 75  |      | µg/L  | 1  | 9/23/2004     |
| 4-Bromophenyl phenyl ether  | ND     | 50  |      | µg/L  | 1  | 9/18/2004     |
| Butyl benzyl phthalate      | ND     | 75  |      | µg/L  | 1  | 9/23/2004     |
| Carbazole                   | ND     | 50  |      | µg/L  | 1  | 9/18/2004     |
| 4-Chloro-3-methylphenol     | ND     | 100 |      | µg/L  | 1  | 9/18/2004     |
| 4-Chloroaniline             | ND     | 100 |      | µg/L  | 1  | 9/18/2004     |
| 2-Chloronaphthalene         | ND     | 50  |      | µg/L  | 1  | 9/18/2004     |
| 2-Chlorophenol              | ND     | 50  |      | µg/L  | 1  | 9/18/2004     |
| 4-Chlorophenyl phenyl ether | ND     | 75  |      | µg/L  | 1  | 9/18/2004     |
| Chrysene                    | ND     | 75  |      | µg/L  | 1  | 9/23/2004     |
| Di-n-butyl phthalate        | ND     | 50  |      | µg/L  | 1  | 9/23/2004     |
| Di-n-octyl phthalate        | ND     | 75  |      | µg/L  | 1  | 9/23/2004     |
| Dibenz(a,h)anthracene       | ND     | 50  |      | µg/L  | 1  | 9/23/2004     |
| Dibenzofuran                | ND     | 50  |      | µg/L  | 1  | 9/23/2004     |
| 1,2-Dichlorobenzene         | ND     | 50  |      | µg/L  | 1  | 9/23/2004     |
| 1,3-Dichlorobenzene         | ND     | 50  |      | µg/L  | 1  | 9/23/2004     |
| 1,4-Dichlorobenzene         | ND     | 50  |      | µg/L  | 1  | 9/23/2004     |
| 3,3'-Dichlorobenzidine      | ND     | 75  |      | µg/L  | 1  | 9/18/2004     |
| Diethyl phthalate           | ND     | 50  |      | µg/L  | 1  | 9/23/2004     |
| Dimethyl phthalate          | ND     | 50  |      | µg/L  | 1  | 9/23/2004     |
| 2,4-Dichlorophenol          | ND     | 50  |      | µg/L  | 1  | 9/18/2004     |
| 2,4-Dimethylphenol          | ND     | 50  |      | µg/L  | 1  | 9/18/2004     |
| 4,6-Dinitro-2-methylphenol  | ND     | 250 |      | µg/L  | 1  | 9/18/2004     |
| 2,4-Dinitrophenol           | ND     | 250 |      | µg/L  | 1  | 9/18/2004     |
| 2,4-Dinitrotoluene          | ND     | 50  |      | µg/L  | 1  | 9/18/2004     |
| 2,6-Dinitrotoluene          | ND     | 50  |      | µg/L  | 1  | 9/18/2004     |
| Fluoranthene                | ND     | 50  |      | µg/L  | 1  | 9/23/2004     |
| Fluorene                    | ND     | 50  |      | µg/L  | 1  | 9/18/2004     |
| Hexachlorobenzene           | ND     | 50  |      | µg/L  | 1  | 9/18/2004     |
| Hexachlorobutadiene         | ND     | 50  |      | µg/L  | 1  | 9/18/2004     |
| Hexachlorocyclopentadiene   | ND     | 50  |      | µg/L  | 1  | 9/18/2004     |

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 Lab Order: 0409046  
 Project: Inj Well 3rd Qtr 2004  
 Lab ID: 0409046-01

Client Sample ID: Injection Well 3rd  
 Collection Date: 9/2/2004 9:20:00 AM

Matrix: AQUEOUS

| Analyses                   | Result | PQL       | Qual | Units | DF | Date Analyzed |
|----------------------------|--------|-----------|------|-------|----|---------------|
| Hexachloroethane           | ND     | 50        |      | µg/L  | 1  | 9/18/2004     |
| Indeno(1,2,3-cd)pyrene     | ND     | 50        |      | µg/L  | 1  | 9/23/2004     |
| Isophorone                 | ND     | 50        |      | µg/L  | 1  | 9/18/2004     |
| 2-Methylnaphthalene        | ND     | 50        |      | µg/L  | 1  | 9/18/2004     |
| 2-Methylphenol             | ND     | 75        |      | µg/L  | 1  | 9/18/2004     |
| 3+4-Methylphenol           | ND     | 50        |      | µg/L  | 1  | 9/18/2004     |
| N-Nitrosodi-n-propylamine  | ND     | 50        |      | µg/L  | 1  | 9/18/2004     |
| N-Nitrosodimethylamine     | ND     | 50        |      | µg/L  | 1  | 9/18/2004     |
| N-Nitrosodiphenylamine     | ND     | 50        |      | µg/L  | 1  | 9/18/2004     |
| Naphthalene                | ND     | 50        |      | µg/L  | 1  | 9/23/2004     |
| 2-Nitroaniline             | ND     | 250       |      | µg/L  | 1  | 9/18/2004     |
| 3-Nitroaniline             | ND     | 250       |      | µg/L  | 1  | 9/18/2004     |
| 4-Nitroaniline             | ND     | 100       |      | µg/L  | 1  | 9/18/2004     |
| Nitrobenzene               | ND     | 50        |      | µg/L  | 1  | 9/18/2004     |
| 2-Nitrophenol              | ND     | 75        |      | µg/L  | 1  | 9/18/2004     |
| 4-Nitrophenol              | ND     | 250       |      | µg/L  | 1  | 9/18/2004     |
| Pentachlorophenol          | ND     | 250       |      | µg/L  | 1  | 9/18/2004     |
| Phenanthrene               | ND     | 50        |      | µg/L  | 1  | 9/23/2004     |
| Phenol                     | ND     | 50        |      | µg/L  | 1  | 9/18/2004     |
| Pyrene                     | ND     | 75        |      | µg/L  | 1  | 9/23/2004     |
| Pyridine                   | ND     | 150       |      | µg/L  | 1  | 9/23/2004     |
| 1,2,4-Trichlorobenzene     | ND     | 50        |      | µg/L  | 1  | 9/18/2004     |
| 2,4,5-Trichlorophenol      | ND     | 50        |      | µg/L  | 1  | 9/18/2004     |
| 2,4,6-Trichlorophenol      | ND     | 75        |      | µg/L  | 1  | 9/18/2004     |
| Surr: 2,4,6-Tribromophenol | 80.0   | 16.6-115  |      | %REC  | 1  | 9/23/2004     |
| Surr: 2,4,6-Tribromophenol | 37.3   | 16.6-115  |      | %REC  | 1  | 9/18/2004     |
| Surr: 2-Fluorobiphenyl     | 64.0   | 37-95.7   |      | %REC  | 1  | 9/23/2004     |
| Surr: 2-Fluorophenol       | 25.7   | 9.54-89.8 |      | %REC  | 1  | 9/18/2004     |
| Surr: 2-Fluorophenol       | 49.0   | 9.54-89.8 |      | %REC  | 1  | 9/23/2004     |
| Surr: 4-Terphenyl-d14      | 71.4   | 47.9-115  |      | %REC  | 1  | 9/23/2004     |
| Surr: Nitrobenzene-d5      | 35.9   | 38-106    | S    | %REC  | 1  | 9/18/2004     |
| Surr: Nitrobenzene-d5      | 58.1   | 38-106    |      | %REC  | 1  | 9/23/2004     |
| Surr: Phenol-d6            | 22.2   | 10.7-63.4 |      | %REC  | 1  | 9/18/2004     |
| Surr: Phenol-d6            | 37.2   | 10.7-63.4 |      | %REC  | 1  | 9/23/2004     |

**EPA 120.1: SPECIFIC CONDUCTANCE**

Analyst: CMC

Specific Conductance 5200 0.010 µmhos/cm 1 9/8/2004

**EPA METHOD 7470: MERCURY**

Analyst: CMC

Mercury 0.017 0.0010 mg/L 5 9/7/2004

**EPA 6010C: TOTAL RECOVERABLE METALS**

Analyst: NMO

Arsenic 0.045 0.020 mg/L 1 9/17/2004 3:02:59 PM

Barium 0.14 0.020 mg/L 1 9/17/2004 3:02:59 PM

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 Project: Inj Well 3rd Qtr 2004  
 Lab ID: 0409046-01

Client Sample ID: Injection Well 3rd  
 Collection Date: 9/2/2004 9:20:00 AM  
 Matrix: AQUEOUS

| Analyses                    | Result | PQL    | Qual | Units    | DF  | Date Analyzed        |
|-----------------------------|--------|--------|------|----------|-----|----------------------|
| Cadmium                     | ND     | 0.0020 |      | mg/L     | 1   | 9/17/2004 3:02:59 PM |
| Calcium                     | 35     | 1.0    |      | mg/L     | 1   | 9/17/2004 3:02:59 PM |
| Chromium                    | 0.0080 | 0.0060 |      | mg/L     | 1   | 9/17/2004 3:02:59 PM |
| Lead                        | ND     | 0.0050 |      | mg/L     | 1   | 9/17/2004 3:02:59 PM |
| Magnesium                   | 14     | 1.0    |      | mg/L     | 1   | 9/17/2004 3:02:59 PM |
| Potassium                   | 13     | 1.0    |      | mg/L     | 1   | 9/17/2004 3:02:59 PM |
| Selenium                    | ND     | 0.050  |      | mg/L     | 1   | 9/17/2004 3:02:59 PM |
| Silver                      | ND     | 0.0050 |      | mg/L     | 1   | 9/20/2004 8:12:42 AM |
| Sodium                      | 1000   | 100    |      | mg/L     | 100 | 9/17/2004 3:42:18 PM |
| <b>EPA METHOD 150.1: PH</b> |        |        |      |          |     | Analyst: CMC         |
| pH                          | 8.36   | 0.010  |      | pH units | 1   | 9/10/2004            |

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# Hall Environmental Analysis Laboratory

Date: 28-Sep-04

CLIENT: San Juan Refining  
 Lab Order: 0409046  
 Project: Inj Well 3rd Qtr 2004  
 Lab ID: 0409046-02

Client Sample ID: Trip Blank  
 Collection Date:  
 Matrix: TRIP BLANK

| Analyses                           | Result | PQL | Qual | Units | DF | Date Analyzed |
|------------------------------------|--------|-----|------|-------|----|---------------|
| <b>EPA METHOD 8260B: VOLATILES</b> |        |     |      |       |    | Analyst: BDH  |
| Benzene                            | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| Toluene                            | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| Ethylbenzene                       | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| Methyl tert-butyl ether (MTBE)     | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| 1,2,4-Trimethylbenzene             | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| 1,3,5-Trimethylbenzene             | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| 1,2-Dichloroethane (EDC)           | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| 1,2-Dibromoethane (EDB)            | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| Naphthalene                        | ND     | 2.0 |      | µg/L  | 1  | 9/8/2004      |
| 1-Methylnaphthalene                | ND     | 4.0 |      | µg/L  | 1  | 9/8/2004      |
| 2-Methylnaphthalene                | ND     | 4.0 |      | µg/L  | 1  | 9/8/2004      |
| Acetone                            | ND     | 10  |      | µg/L  | 1  | 9/8/2004      |
| Bromobenzene                       | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| Bromochloromethane                 | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| Bromodichloromethane               | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| Bromoform                          | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| Bromomethane                       | ND     | 2.0 |      | µg/L  | 1  | 9/8/2004      |
| 2-Butanone                         | ND     | 10  |      | µg/L  | 1  | 9/8/2004      |
| Carbon disulfide                   | ND     | 10  |      | µg/L  | 1  | 9/8/2004      |
| Carbon Tetrachloride               | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| Chlorobenzene                      | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| Chloroethane                       | ND     | 2.0 |      | µg/L  | 1  | 9/8/2004      |
| Chloroform                         | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| Chloromethane                      | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| 2-Chlorotoluene                    | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| 4-Chlorotoluene                    | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| cis-1,2-DCE                        | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| cis-1,3-Dichloropropene            | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| 1,2-Dibromo-3-chloropropane        | ND     | 2.0 |      | µg/L  | 1  | 9/8/2004      |
| Dibromochloromethane               | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| Dibromomethane                     | ND     | 2.0 |      | µg/L  | 1  | 9/8/2004      |
| 1,2-Dichlorobenzene                | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| 1,3-Dichlorobenzene                | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| 1,4-Dichlorobenzene                | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| Dichlorodifluoromethane            | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| 1,1-Dichloroethane                 | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| 1,1-Dichloroethene                 | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| 1,2-Dichloropropane                | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| 1,3-Dichloropropane                | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| 2,2-Dichloropropane                | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |
| 1,1-Dichloropropene                | ND     | 1.0 |      | µg/L  | 1  | 9/8/2004      |

Qualifiers: ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits  
 J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits  
 B - Analyte detected in the associated Method Blank      E - Value above quantitation range  
 \* - Value exceeds Maximum Contaminant Level



# Hall Environmental Analysis Laboratory

Date: 28-Sep-04

CLIENT: San Juan Refining  
 Lab Order: 0409046  
 Project: Inj Well 3rd Qtr 2004  
 Lab ID: 0409046-02

Client Sample ID: Trip Blank  
 Collection Date:  
 Matrix: TRIP BLANK

| Analyses                    | Result | PQL      | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|----------|------|-------|----|---------------|
| Hexachlorobutadiene         | ND     | 1.0      |      | µg/L  | 1  | 9/8/2004      |
| 2-Hexanone                  | ND     | 10       |      | µg/L  | 1  | 9/8/2004      |
| Isopropylbenzene            | ND     | 1.0      |      | µg/L  | 1  | 9/8/2004      |
| 4-Isopropyltoluene          | ND     | 1.0      |      | µg/L  | 1  | 9/8/2004      |
| 4-Methyl-2-pentanone        | ND     | 10       |      | µg/L  | 1  | 9/8/2004      |
| Methylene Chloride          | ND     | 3.0      |      | µg/L  | 1  | 9/8/2004      |
| n-Butylbenzene              | ND     | 1.0      |      | µg/L  | 1  | 9/8/2004      |
| n-Propylbenzene             | ND     | 1.0      |      | µg/L  | 1  | 9/8/2004      |
| sec-Butylbenzene            | ND     | 1.0      |      | µg/L  | 1  | 9/8/2004      |
| Styrene                     | ND     | 1.0      |      | µg/L  | 1  | 9/8/2004      |
| tert-Butylbenzene           | ND     | 1.0      |      | µg/L  | 1  | 9/8/2004      |
| 1,1,1,2-Tetrachloroethane   | ND     | 1.0      |      | µg/L  | 1  | 9/8/2004      |
| 1,1,2,2-Tetrachloroethane   | ND     | 1.0      |      | µg/L  | 1  | 9/8/2004      |
| Tetrachloroethene (PCE)     | ND     | 1.0      |      | µg/L  | 1  | 9/8/2004      |
| trans-1,2-DCE               | ND     | 1.0      |      | µg/L  | 1  | 9/8/2004      |
| trans-1,3-Dichloropropene   | ND     | 1.0      |      | µg/L  | 1  | 9/8/2004      |
| 1,2,3-Trichlorobenzene      | ND     | 1.0      |      | µg/L  | 1  | 9/8/2004      |
| 1,2,4-Trichlorobenzene      | ND     | 1.0      |      | µg/L  | 1  | 9/8/2004      |
| 1,1,1-Trichloroethane       | ND     | 1.0      |      | µg/L  | 1  | 9/8/2004      |
| 1,1,2-Trichloroethane       | ND     | 1.0      |      | µg/L  | 1  | 9/8/2004      |
| Trichloroethene (TCE)       | ND     | 1.0      |      | µg/L  | 1  | 9/8/2004      |
| Trichlorofluoromethane      | ND     | 1.0      |      | µg/L  | 1  | 9/8/2004      |
| 1,2,3-Trichloropropane      | ND     | 2.0      |      | µg/L  | 1  | 9/8/2004      |
| Vinyl chloride              | ND     | 1.0      |      | µg/L  | 1  | 9/8/2004      |
| Xylenes, Total              | ND     | 1.0      |      | µg/L  | 1  | 9/8/2004      |
| Surr: 1,2-Dichloroethane-d4 | 93.5   | 70.6-124 |      | %REC  | 1  | 9/8/2004      |
| Surr: 4-Bromofluorobenzene  | 103    | 76.4-130 |      | %REC  | 1  | 9/8/2004      |
| Surr: Dibromofluoromethane  | 88.3   | 67.2-131 |      | %REC  | 1  | 9/8/2004      |
| Surr: Toluene-d8            | 106    | 82.1-123 |      | %REC  | 1  | 9/8/2004      |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits  
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits  
 B - Analyte detected in the associated Method Blank E - Value above quantitation range  
 \* - Value exceeds Maximum Contaminant Level



LABORATORY ANALYTICAL REPORT

Client: Hall Environmental-Albuquerque  
 Project: Inj Well  
 Lab ID: B04090453-001  
 Client Sample ID: Injection Well 3rd, 0409046-01E

Report Date: 09/17/04  
 Collection Date: 09/02/04 09:20  
 Date Received: 09/08/04  
 Matrix: Aqueous

| Analyses                   | Result | Units | Qual | MCL/ |     | Method     | Analysis Date / By   |
|----------------------------|--------|-------|------|------|-----|------------|----------------------|
|                            |        |       |      | RL   | QCL |            |                      |
| <b>INORGANICS</b>          |        |       |      |      |     |            |                      |
| Sulfide                    | ND     | mg/L  |      | 1    |     | E376.1     | 09/08/04 13:00 / pwc |
| <b>IGNITABILITY</b>        |        |       |      |      |     |            |                      |
| Flash Point (Ignitability) | >200   | °F    |      | 30   |     | SW1010     | 09/08/04 13:30 / mgs |
| <b>CORROSIVITY</b>         |        |       |      |      |     |            |                      |
| pH of Liquid Waste         | 8.21   | s.u.  |      | 0.10 |     | SW9040B    | 09/08/04 13:00 / mgs |
| <b>REACTIVITY</b>          |        |       |      |      |     |            |                      |
| Cyanide, Reactive          | ND     | mg/kg |      | 0.05 | 250 | SW846 Ch 7 | 09/15/04 10:05 / kjp |
| Sulfide, Reactive          | ND     | mg/kg |      | 20   | 500 | SW846 Ch 7 | 09/14/04 08:00 / pwc |

Report Definitions: RL - Analyte reporting limit.  
 QCL - Quality control limit.

MCL - Maximum contaminant level.  
 ND - Not detected at the reporting limit.



## QA/QC Summary Report

Client: Hall Environmental-Albuquerque

Report Date: 09/09/04

Project: Inj Well

Work Order: B04090453

| Analyte                      | Result                        | Units | RL  | %REC | Low Limit | High Limit | RPD | RPDLimit | Qual                         |
|------------------------------|-------------------------------|-------|-----|------|-----------|------------|-----|----------|------------------------------|
| Method: E376.1               |                               |       |     |      |           |            |     |          | Batch: 040908A-SULFIDE-TTR-W |
| Sample ID: B04090456-002BDUP | Sample Duplicate              |       |     |      |           |            |     |          | 09/08/04 13:00               |
| Sulfide                      | 1320                          | mg/L  | 1.0 |      |           |            | 1.5 | 20       |                              |
| Sample ID: B04090456-002BMS  | Sample Matrix Spike           |       |     |      |           |            |     |          | 09/08/04 13:00               |
| Sulfide                      | 2400                          | mg/L  | 1.0 | 98.1 | 80        | 120        |     |          |                              |
| Sample ID: B04090456-002BMSD | Sample Matrix Spike Duplicate |       |     |      |           |            |     |          | 09/08/04 13:00               |
| Sulfide                      | 2400                          | mg/L  | 1.0 | 98.1 | 80        | 120        | 0   | 20       |                              |

**Qualifiers:**

RL - Analyte reporting limit.

ND - Not detected at the reporting limit.



### QA/QC Summary Report

Client: Hall Environmental-Albuquerque

Report Date: 09/15/04

Project: Inj Well

Work Order: B04090453

| Analyte                      | Result                   | Units | RL   | %REC | Low Limit | High Limit | RPD | RPDLimit      | Qual           |
|------------------------------|--------------------------|-------|------|------|-----------|------------|-----|---------------|----------------|
| <b>Method: SW846 Ch 7</b>    |                          |       |      |      |           |            |     | Batch: 12267  |                |
| <b>Sample ID: MB-12267</b>   | Method Blank             |       |      |      |           |            |     |               | 09/15/04 10:07 |
| Cyanide, Reactive            | ND                       | mg/kg | 0.05 |      |           |            |     |               |                |
| <b>Method: SW846 Ch 7</b>    |                          |       |      |      |           |            |     | Batch: R47626 |                |
| <b>Sample ID: MB-R47626</b>  | Method Blank             |       |      |      |           |            |     |               | 09/14/04 08:00 |
| Sulfide, Reactive            | ND                       | mg/kg | 10   |      |           |            |     |               |                |
| <b>Sample ID: LCS-R47626</b> | Laboratory Control Spike |       |      |      |           |            |     |               | 09/14/04 08:00 |
| Sulfide, Reactive            | 22                       | mg/kg | 20   | 71   | 50        | 150        |     |               |                |

**Qualifiers:**

RL - Analyte reporting limit.

ND - Not detected at the reporting limit.

Hall Environmental Analysis Laboratory

Date: 28-Sep-04

CLIENT: San Juan Refining  
 Work Order: 0409046  
 Project: Inj Well 3rd Qtr 2004

QC SUMMARY REPORT  
 Method Blank

| Sample ID  | Batch ID | Test Code          | Units     | Analysis Date         | Prep Date |          |           |             |      |          |      |
|------------|----------|--------------------|-----------|-----------------------|-----------|----------|-----------|-------------|------|----------|------|
| MBLK       | R13106   | E300               | mg/L      | 9/13/2004 12:12:52 PM |           |          |           |             |      |          |      |
| Client ID: |          | Run ID: LC_040913A |           | SeqNo: 304865         |           |          |           |             |      |          |      |
| Analyte    | Result   | PQL                | SPK value | SPK Ref Val           | %REC      | LowLimit | HighLimit | RPD Ref Val | %RPD | RPDLimit | Qual |
| Chloride   | ND       | 0.1                |           |                       |           |          |           |             |      |          |      |
| Sulfate    | ND       | 0.5                |           |                       |           |          |           |             |      |          |      |

| Sample ID  | Batch ID | Test Code          | Units     | Analysis Date        | Prep Date |          |           |             |      |          |      |
|------------|----------|--------------------|-----------|----------------------|-----------|----------|-----------|-------------|------|----------|------|
| MBLK       | R13133   | E300               | mg/L      | 9/15/2004 2:15:58 PM |           |          |           |             |      |          |      |
| Client ID: |          | Run ID: LC_040915A |           | SeqNo: 305674        |           |          |           |             |      |          |      |
| Analyte    | Result   | PQL                | SPK value | SPK Ref Val          | %REC      | LowLimit | HighLimit | RPD Ref Val | %RPD | RPDLimit | Qual |
| Chloride   | ND       | 0.1                |           |                      |           |          |           |             |      |          |      |
| Sulfate    | ND       | 0.5                |           |                      |           |          |           |             |      |          |      |

| Sample ID  | Batch ID | Test Code          | Units     | Analysis Date         | Prep Date |          |           |             |      |          |      |
|------------|----------|--------------------|-----------|-----------------------|-----------|----------|-----------|-------------|------|----------|------|
| MBLK       | R13133   | E300               | mg/L      | 9/16/2004 12:37:40 AM |           |          |           |             |      |          |      |
| Client ID: |          | Run ID: LC_040915A |           | SeqNo: 305695         |           |          |           |             |      |          |      |
| Analyte    | Result   | PQL                | SPK value | SPK Ref Val           | %REC      | LowLimit | HighLimit | RPD Ref Val | %RPD | RPDLimit | Qual |
| Chloride   | 0.1383   | 0.1                |           |                       |           |          |           |             |      |          |      |
| Sulfate    | ND       | 0.5                |           |                       |           |          |           |             |      |          |      |

12/28

Qualifiers: ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits      B - Analyte detected in the associated Method Blank  
 J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits

**CLIENT:** San Juan Refining  
**Work Order:** 0409046  
**Project:** Inj Well 3rd Qtr 2004

**QC SUMMARY REPORT**  
 Method Blank

Sample ID **mb-6493**      Batch ID: **6493**      Test Code: **SW8270A**      Units: **µg/L**      Analysis Date **9/23/2004**      Prep Date **9/16/2004**  
 Client ID:      Run ID: **ELMO\_040922A**      SeqNo: **308219**

| Analyte                     | Result | PQL | SPK value | SPK Ref Val | %REC | LowLimit | HighLimit | RPD Ref Val | %RPD | RPDLimit | Qual |
|-----------------------------|--------|-----|-----------|-------------|------|----------|-----------|-------------|------|----------|------|
| Acenaphthene                | ND     | 10  |           |             |      |          |           |             |      |          |      |
| Acenaphthylene              | ND     | 10  |           |             |      |          |           |             |      |          |      |
| Aniline                     | ND     | 10  |           |             |      |          |           |             |      |          |      |
| Anthracene                  | ND     | 10  |           |             |      |          |           |             |      |          |      |
| Azobenzene                  | ND     | 10  |           |             |      |          |           |             |      |          |      |
| Benzo(a)anthracene          | ND     | 15  |           |             |      |          |           |             |      |          |      |
| Benzo(a)pyrene              | ND     | 10  |           |             |      |          |           |             |      |          |      |
| Benzo(b)fluoranthene        | ND     | 10  |           |             |      |          |           |             |      |          |      |
| Benzo(g,h,i)perylene        | ND     | 10  |           |             |      |          |           |             |      |          |      |
| Benzo(k)fluoranthene        | ND     | 10  |           |             |      |          |           |             |      |          |      |
| Benzoic acid                | ND     | 50  |           |             |      |          |           |             |      |          |      |
| Benzyl alcohol              | ND     | 20  |           |             |      |          |           |             |      |          |      |
| Bis(2-chloroethoxy)methane  | ND     | 10  |           |             |      |          |           |             |      |          |      |
| Bis(2-chloroethyl)ether     | ND     | 15  |           |             |      |          |           |             |      |          |      |
| Bis(2-chloroisopropyl)ether | ND     | 15  |           |             |      |          |           |             |      |          |      |
| Bis(2-ethylhexyl)phthalate  | ND     | 15  |           |             |      |          |           |             |      |          |      |
| 4-Bromophenyl phenyl ether  | ND     | 10  |           |             |      |          |           |             |      |          |      |
| Butyl benzyl phthalate      | ND     | 15  |           |             |      |          |           |             |      |          |      |
| Carbazole                   | ND     | 10  |           |             |      |          |           |             |      |          |      |
| 4-Chloro-3-methylphenol     | ND     | 20  |           |             |      |          |           |             |      |          |      |
| 4-Chloroaniline             | ND     | 20  |           |             |      |          |           |             |      |          |      |
| 2-Chloronaphthalene         | ND     | 10  |           |             |      |          |           |             |      |          |      |
| 2-Chlorophenol              | ND     | 10  |           |             |      |          |           |             |      |          |      |
| 4-Chlorophenyl phenyl ether | ND     | 15  |           |             |      |          |           |             |      |          |      |
| Chrysene                    | ND     | 15  |           |             |      |          |           |             |      |          |      |
| Di-n-butyl phthalate        | ND     | 10  |           |             |      |          |           |             |      |          |      |
| Di-n-octyl phthalate        | ND     | 15  |           |             |      |          |           |             |      |          |      |
| Dibenz(a,h)anthracene       | ND     | 10  |           |             |      |          |           |             |      |          |      |

13/28

**Qualifiers:**      ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits      B - Analyte detected in the associated Method Blank  
                          J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits

**CLIENT:** San Juan Refining  
**Work Order:** 0409046  
**Project:** Inj Well 3rd Qtr 2004

## QC SUMMARY REPORT

Method Blank

|                            |    |    |
|----------------------------|----|----|
| Dibenzofuran               | ND | 10 |
| 1,2-Dichlorobenzene        | ND | 10 |
| 1,3-Dichlorobenzene        | ND | 10 |
| 1,4-Dichlorobenzene        | ND | 10 |
| 3,3'-Dichlorobenzidine     | ND | 15 |
| Diethyl phthalate          | ND | 10 |
| Dimethyl phthalate         | ND | 10 |
| 2,4-Dichlorophenol         | ND | 10 |
| 2,4-Dimethylphenol         | ND | 10 |
| 4,6-Dinitro-2-methylphenol | ND | 50 |
| 2,4-Dinitrophenol          | ND | 50 |
| 2,4-Dinitrotoluene         | ND | 10 |
| 2,6-Dinitrotoluene         | ND | 10 |
| Fluoranthene               | ND | 10 |
| Fluorene                   | ND | 10 |
| Hexachlorobenzene          | ND | 10 |
| Hexachlorobutadiene        | ND | 10 |
| Hexachlorocyclopentadiene  | ND | 10 |
| Hexachloroethane           | ND | 10 |
| Indeno(1,2,3-cd)pyrene     | ND | 10 |
| Isophorone                 | ND | 10 |
| 2-Methylnaphthalene        | ND | 10 |
| 2-Methylphenol             | ND | 15 |
| 3+4-Methylphenol           | ND | 10 |
| N-Nitrosodi-n-propylamine  | ND | 10 |
| N-Nitrosodimethylamine     | ND | 10 |
| N-Nitrosodiphenylamine     | ND | 10 |
| Naphthalene                | ND | 10 |
| 2-Nitroaniline             | ND | 50 |
| 3-Nitroaniline             | ND | 50 |
| 4-Nitroaniline             | ND | 20 |
| Nitrobenzene               | ND | 10 |
| 2-Nitrophenol              | ND | 15 |

14/28

**Qualifiers:** ND - Not Detected at the Reporting Limit  
 J - Analyte detected below quantitation limits

S - Spike Recovery outside accepted recovery limits  
 R - RPD outside accepted recovery limits

B - Analyte detected in the associated Method Blank

CLIENT: San Juan Refining  
 Work Order: 0409046  
 Project: Inj Well 3rd Qtr 2004

**QC SUMMARY REPORT**

Method Blank

|                            |       |    |     |   |      |      |      |   |  |
|----------------------------|-------|----|-----|---|------|------|------|---|--|
| 4-Nitrophenol              | ND    | 50 |     |   |      |      |      |   |  |
| Pentachlorophenol          | ND    | 50 |     |   |      |      |      |   |  |
| Phenanthrene               | ND    | 10 |     |   |      |      |      |   |  |
| Phenol                     | ND    | 10 |     |   |      |      |      |   |  |
| Pyrene                     | ND    | 15 |     |   |      |      |      |   |  |
| Pyridine                   | ND    | 30 |     |   |      |      |      |   |  |
| 1,2,4-Trichlorobenzene     | ND    | 10 |     |   |      |      |      |   |  |
| 2,4,5-Trichlorophenol      | ND    | 10 |     |   |      |      |      |   |  |
| 2,4,6-Trichlorophenol      | ND    | 15 |     |   |      |      |      |   |  |
| Surr: 2,4,6-Tribromophenol | 71.32 | 0  | 200 | 0 | 35.7 | 16.6 | 115  | 0 |  |
| Surr: 2-Fluorobiphenyl     | 58.44 | 0  | 100 | 0 | 58.4 | 37   | 95.7 | 0 |  |
| Surr: 2-Fluorophenol       | 106.8 | 0  | 200 | 0 | 53.4 | 9.54 | 89.8 | 0 |  |
| Surr: 4-Terphenyl-d14      | 96.36 | 0  | 100 | 0 | 96.4 | 51.2 | 125  | 0 |  |
| Surr: Nitrobenzene-d5      | 63.66 | 0  | 100 | 0 | 63.7 | 38   | 106  | 0 |  |
| Surr: Phenol-d6            | 76.32 | 0  | 200 | 0 | 38.2 | 10.7 | 63.4 | 0 |  |

|            |           |           |                  |             |        |          |           |               |          |           |          |
|------------|-----------|-----------|------------------|-------------|--------|----------|-----------|---------------|----------|-----------|----------|
| Sample ID  | MB-6436   | Batch ID: | 6436             | Test Code:  | SW7470 | Units:   | mg/L      | Analysis Date | 9/7/2004 | Prep Date | 9/7/2004 |
| Client ID: |           | Run ID:   | MI-LA254_040907A | SeqNo:      | 302914 |          |           |               |          |           |          |
| Analyte    | Result    | PQL       | SPK value        | SPK Ref Val | %REC   | LowLimit | HighLimit | RPD Ref Val   | %RPD     | RPDLimit  | Qual     |
| Mercury    | 0.0001224 | 0.0002    |                  |             |        |          |           |               |          |           | J        |

Qualifiers: ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits      B - Analyte detected in the associated Method Blank  
 J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits

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**CLIENT:** San Juan Refining  
**Work Order:** 0409046  
**Project:** Inj Well 3rd Qtr 2004

## QC SUMMARY REPORT

Method Blank

| Sample ID  | Batch ID | Test Code | Units       | Analysis Date        | Prep Date |          |           |             |      |          |      |
|------------|----------|-----------|-------------|----------------------|-----------|----------|-----------|-------------|------|----------|------|
| MB-6492    | 6492     | SW6010A   | mg/L        | 9/17/2004 2:52:50 PM | 9/16/2004 |          |           |             |      |          |      |
| Client ID: |          | Run ID:   | ICP_040917A | SeqNo:               | 306647    |          |           |             |      |          |      |
| Analyte    | Result   | PQL       | SPK value   | SPK Ref Val          | %REC      | LowLimit | HighLimit | RPD Ref Val | %RPD | RPDLimit | Qual |
| Arsenic    | ND       | 0.02      |             |                      |           |          |           |             |      |          |      |
| Barium     | ND       | 0.02      |             |                      |           |          |           |             |      |          |      |
| Cadmium    | 0.001101 | 0.002     |             |                      |           |          |           |             |      |          | J    |
| Calcium    | ND       | 1         |             |                      |           |          |           |             |      |          |      |
| Chromium   | ND       | 0.006     |             |                      |           |          |           |             |      |          |      |
| Lead       | 0.003158 | 0.005     |             |                      |           |          |           |             |      |          | J    |
| Magnesium  | ND       | 1         |             |                      |           |          |           |             |      |          |      |
| Potassium  | ND       | 1         |             |                      |           |          |           |             |      |          |      |
| Selenium   | ND       | 0.05      |             |                      |           |          |           |             |      |          |      |
| Sodium     | ND       | 1         |             |                      |           |          |           |             |      |          |      |

| Sample ID  | Batch ID | Test Code | Units       | Analysis Date         | Prep Date |          |           |             |      |          |      |
|------------|----------|-----------|-------------|-----------------------|-----------|----------|-----------|-------------|------|----------|------|
| MB-6492    | 6492     | SW6010A   | mg/L        | 9/20/2004 11:27:57 AM | 9/16/2004 |          |           |             |      |          |      |
| Client ID: |          | Run ID:   | ICP_040920A | SeqNo:                | 306678    |          |           |             |      |          |      |
| Analyte    | Result   | PQL       | SPK value   | SPK Ref Val           | %REC      | LowLimit | HighLimit | RPD Ref Val | %RPD | RPDLimit | Qual |
| Arsenic    | ND       | 0.02      |             |                       |           |          |           |             |      |          |      |
| Barium     | ND       | 0.02      |             |                       |           |          |           |             |      |          |      |
| Cadmium    | ND       | 0.002     |             |                       |           |          |           |             |      |          |      |
| Calcium    | ND       | 1         |             |                       |           |          |           |             |      |          |      |
| Chromium   | ND       | 0.006     |             |                       |           |          |           |             |      |          |      |
| Lead       | ND       | 0.005     |             |                       |           |          |           |             |      |          |      |
| Magnesium  | ND       | 1         |             |                       |           |          |           |             |      |          |      |
| Potassium  | ND       | 1         |             |                       |           |          |           |             |      |          |      |
| Selenium   | ND       | 0.05      |             |                       |           |          |           |             |      |          |      |
| Sodium     | ND       | 1         |             |                       |           |          |           |             |      |          |      |

**Qualifiers:** ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits      B - Analyte detected in the associated Method Blank  
 J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits

16/28

**CLIENT:** San Juan Refining  
**Work Order:** 0409046  
**Project:** Inj Well 3rd Qtr 2004

## QC SUMMARY REPORT

Method Blank

|            |         |           |             |            |           |             |      |               |                       |           |             |      |          |      |
|------------|---------|-----------|-------------|------------|-----------|-------------|------|---------------|-----------------------|-----------|-------------|------|----------|------|
| Sample ID  | MB-6492 | Batch ID: | 6492        | Test Code: | SW6010A   | Units:      | mg/L | Analysis Date | 9/20/2004 12:31:59 PM | Prep Date | 9/16/2004   |      |          |      |
| Client ID: |         | Run ID:   | ICP_040920A | SeqNo:     | 306682    |             |      |               |                       |           |             |      |          |      |
| Analyte    |         | Result    |             | PQL        | SPK value | SPK Ref Val |      | %REC          | LowLimit              | HighLimit | RPD Ref Val | %RPD | RPDLimit | Qual |
| Silver     |         | ND        |             | 0.005      |           |             |      |               |                       |           |             |      |          |      |

|            |         |           |             |            |           |             |      |               |                      |           |             |      |          |      |
|------------|---------|-----------|-------------|------------|-----------|-------------|------|---------------|----------------------|-----------|-------------|------|----------|------|
| Sample ID  | MB-6492 | Batch ID: | 6492        | Test Code: | SW6010A   | Units:      | mg/L | Analysis Date | 9/20/2004 8:05:04 AM | Prep Date | 9/16/2004   |      |          |      |
| Client ID: |         | Run ID:   | ICP_040920E | SeqNo:     | 307231    |             |      |               |                      |           |             |      |          |      |
| Analyte    |         | Result    |             | PQL        | SPK value | SPK Ref Val |      | %REC          | LowLimit             | HighLimit | RPD Ref Val | %RPD | RPDLimit | Qual |
| Silver     |         | ND        |             | 0.005      |           |             |      |               |                      |           |             |      |          |      |

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**Qualifiers:** ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits      B - Analyte detected in the associated Method Blank  
 J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits

Hall Environmental Analysis Laboratory

Date: 28-Sep-04

CLIENT: San Juan Refining  
 Work Order: 0409046  
 Project: Inj Well 3rd Qtr 2004

QC SUMMARY REPORT

Method Blank

| Sample ID                      | Batch ID | Test Code | Units        | Analysis Date | Prep Date |          |           |             |      |          |      |
|--------------------------------|----------|-----------|--------------|---------------|-----------|----------|-----------|-------------|------|----------|------|
| 5ml rb                         | R13035   | SW8260B   | µg/L         | 9/7/2004      |           |          |           |             |      |          |      |
| Client ID:                     |          | Run ID:   | THOR_040907A | SeqNo:        | 303134    |          |           |             |      |          |      |
| Analyte                        | Result   | PQL       | SPK value    | SPK Ref Val   | %REC      | LowLimit | HighLimit | RPD Ref Val | %RPD | RPDLimit | Qual |
| Benzene                        | ND       | 1         |              |               |           |          |           |             |      |          |      |
| Toluene                        | ND       | 1         |              |               |           |          |           |             |      |          |      |
| Ethylbenzene                   | ND       | 1         |              |               |           |          |           |             |      |          |      |
| Methyl tert-butyl ether (MTBE) | ND       | 1         |              |               |           |          |           |             |      |          |      |
| 1,2,4-Trimethylbenzene         | ND       | 1         |              |               |           |          |           |             |      |          |      |
| 1,3,5-Trimethylbenzene         | ND       | 1         |              |               |           |          |           |             |      |          |      |
| 1,2-Dichloroethane (EDC)       | ND       | 1         |              |               |           |          |           |             |      |          |      |
| 1,2-Dibromoethane (EDB)        | ND       | 1         |              |               |           |          |           |             |      |          |      |
| Naphthalene                    | ND       | 2         |              |               |           |          |           |             |      |          |      |
| 1-Methylnaphthalene            | ND       | 4         |              |               |           |          |           |             |      |          |      |
| 2-Methylnaphthalene            | ND       | 4         |              |               |           |          |           |             |      |          |      |
| Acetone                        | ND       | 10        |              |               |           |          |           |             |      |          |      |
| Bromobenzene                   | ND       | 1         |              |               |           |          |           |             |      |          |      |
| Bromochloromethane             | ND       | 1         |              |               |           |          |           |             |      |          |      |
| Bromodichloromethane           | ND       | 1         |              |               |           |          |           |             |      |          |      |
| Bromoform                      | ND       | 1         |              |               |           |          |           |             |      |          |      |
| Bromomethane                   | 0.39     | 2         |              |               |           |          |           |             |      |          | J    |
| 2-Butanone                     | ND       | 10        |              |               |           |          |           |             |      |          |      |
| Carbon disulfide               | ND       | 10        |              |               |           |          |           |             |      |          |      |
| Carbon Tetrachloride           | ND       | 1         |              |               |           |          |           |             |      |          |      |
| Chlorobenzene                  | ND       | 1         |              |               |           |          |           |             |      |          |      |
| Chloroethane                   | ND       | 2         |              |               |           |          |           |             |      |          |      |
| Chloroform                     | ND       | 1         |              |               |           |          |           |             |      |          |      |
| Chloromethane                  | ND       | 1         |              |               |           |          |           |             |      |          |      |
| 2-Chlorotoluene                | ND       | 1         |              |               |           |          |           |             |      |          |      |
| 4-Chlorotoluene                | ND       | 1         |              |               |           |          |           |             |      |          |      |
| cis-1,2-DCE                    | ND       | 1         |              |               |           |          |           |             |      |          |      |

18/28

Qualifiers: ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits      B - Analyte detected in the associated Method Blank  
 J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits

**CLIENT:** San Juan Refining  
**Work Order:** 0409046  
**Project:** Inj Well 3rd Qtr 2004

**QC SUMMARY REPORT**

Method Blank

|                             |    |    |
|-----------------------------|----|----|
| cis-1,3-Dichloropropene     | ND | 1  |
| 1,2-Dibromo-3-chloropropane | ND | 2  |
| Dibromochloromethane        | ND | 1  |
| Dibromomethane              | ND | 2  |
| 1,2-Dichlorobenzene         | ND | 1  |
| 1,3-Dichlorobenzene         | ND | 1  |
| 1,4-Dichlorobenzene         | ND | 1  |
| Dichlorodifluoromethane     | ND | 1  |
| 1,1-Dichloroethane          | ND | 1  |
| 1,1-Dichloroethene          | ND | 1  |
| 1,2-Dichloropropane         | ND | 1  |
| 1,3-Dichloropropane         | ND | 1  |
| 2,2-Dichloropropane         | ND | 1  |
| 1,1-Dichloropropene         | ND | 1  |
| Hexachlorobutadiene         | ND | 1  |
| 2-Hexanone                  | ND | 10 |
| Isopropylbenzene            | ND | 1  |
| 4-Isopropyltoluene          | ND | 1  |
| 4-Methyl-2-pentanone        | ND | 10 |
| Methylene Chloride          | ND | 3  |
| n-Butylbenzene              | ND | 1  |
| n-Propylbenzene             | ND | 1  |
| sec-Butylbenzene            | ND | 1  |
| Styrene                     | ND | 1  |
| tert-Butylbenzene           | ND | 1  |
| 1,1,1,2-Tetrachloroethane   | ND | 1  |
| 1,1,2,2-Tetrachloroethane   | ND | 1  |
| Tetrachloroethene (PCE)     | ND | 1  |
| trans-1,2-DCE               | ND | 1  |
| trans-1,3-Dichloropropene   | ND | 1  |
| 1,2,3-Trichlorobenzene      | ND | 1  |
| 1,2,4-Trichlorobenzene      | ND | 1  |
| 1,1,1-Trichloroethane       | ND | 1  |

19/28

**Qualifiers:** ND - Not Detected at the Reporting Limit  
 J - Analyte detected below quantitation limits

S - Spike Recovery outside accepted recovery limits  
 R - RPD outside accepted recovery limits

B - Analyte detected in the associated Method Blank

**CLIENT:** San Juan Refining  
**Work Order:** 0409046  
**Project:** Inj Well 3rd Qtr 2004

**QC SUMMARY REPORT**

Method Blank

|                             |       |   |    |   |      |      |     |   |  |
|-----------------------------|-------|---|----|---|------|------|-----|---|--|
| 1,1,2-Trichloroethane       | ND    | 1 |    |   |      |      |     |   |  |
| Trichloroethene (TCE)       | ND    | 1 |    |   |      |      |     |   |  |
| Trichlorofluoromethane      | ND    | 1 |    |   |      |      |     |   |  |
| 1,2,3-Trichloropropane      | ND    | 2 |    |   |      |      |     |   |  |
| Vinyl chloride              | ND    | 1 |    |   |      |      |     |   |  |
| Xylenes, Total              | ND    | 1 |    |   |      |      |     |   |  |
| Surr: 1,2-Dichloroethane-d4 | 9.374 | 0 | 10 | 0 | 93.7 | 68.4 | 127 | 0 |  |
| Surr: 4-Bromofluorobenzene  | 10.18 | 0 | 10 | 0 | 102  | 70.4 | 126 | 0 |  |
| Surr: Dibromofluoromethane  | 9.394 | 0 | 10 | 0 | 93.9 | 70.2 | 126 | 0 |  |
| Surr: Toluene-d8            | 10.45 | 0 | 10 | 0 | 104  | 73.5 | 129 | 0 |  |

20/28

**Qualifiers:** ND - Not Detected at the Reporting Limit  
 J - Analyte detected below quantitation limits

S - Spike Recovery outside accepted recovery limits  
 R - RPD outside accepted recovery limits

B - Analyte detected in the associated Method Blank

Hall Environmental Analysis Laboratory

Date: 28-Sep-04

**CLIENT:** San Juan Refining  
**Work Order:** 0409046  
**Project:** Inj Well 3rd Qtr 2004

**QC SUMMARY REPORT**  
 Laboratory Control Spike - generic

|            |     |           |            |            |           |             |      |               |                       |           |             |      |          |      |
|------------|-----|-----------|------------|------------|-----------|-------------|------|---------------|-----------------------|-----------|-------------|------|----------|------|
| Sample ID  | LCS | Batch ID: | R13106     | Test Code: | E300      | Units:      | mg/L | Analysis Date | 9/13/2004 12:29:41 PM | Prep Date |             |      |          |      |
| Client ID: |     | Run ID:   | LC_040913A | SeqNo:     | 304866    |             |      |               |                       |           |             |      |          |      |
| Analyte    |     | Result    |            | PQL        | SPK value | SPK Ref Val |      | %REC          | LowLimit              | HighLimit | RPD Ref Val | %RPD | RPDLimit | Qual |
| Chloride   |     | 4.506     |            | 0.1        | 5         | 0           |      | 90.1          | 90                    | 110       | 0           |      |          |      |
| Sulfate    |     | 9.159     |            | 0.5        | 10        | 0           |      | 91.6          | 90                    | 110       | 0           |      |          |      |

|            |     |           |            |            |           |             |      |               |                      |           |             |      |          |      |
|------------|-----|-----------|------------|------------|-----------|-------------|------|---------------|----------------------|-----------|-------------|------|----------|------|
| Sample ID  | LCS | Batch ID: | R13133     | Test Code: | E300      | Units:      | mg/L | Analysis Date | 9/15/2004 2:32:47 PM | Prep Date |             |      |          |      |
| Client ID: |     | Run ID:   | LC_040915A | SeqNo:     | 305675    |             |      |               |                      |           |             |      |          |      |
| Analyte    |     | Result    |            | PQL        | SPK value | SPK Ref Val |      | %REC          | LowLimit             | HighLimit | RPD Ref Val | %RPD | RPDLimit | Qual |
| Chloride   |     | 5.051     |            | 0.1        | 5         | 0.1383      |      | 98.3          | 90                   | 110       | 0           |      |          |      |
| Sulfate    |     | 9.823     |            | 0.5        | 10        | 0           |      | 98.2          | 90                   | 110       | 0           |      |          |      |

|            |     |           |            |            |           |             |      |               |                       |           |             |      |          |      |
|------------|-----|-----------|------------|------------|-----------|-------------|------|---------------|-----------------------|-----------|-------------|------|----------|------|
| Sample ID  | LCS | Batch ID: | R13133     | Test Code: | E300      | Units:      | mg/L | Analysis Date | 9/16/2004 12:54:28 AM | Prep Date |             |      |          |      |
| Client ID: |     | Run ID:   | LC_040915A | SeqNo:     | 305696    |             |      |               |                       |           |             |      |          |      |
| Analyte    |     | Result    |            | PQL        | SPK value | SPK Ref Val |      | %REC          | LowLimit              | HighLimit | RPD Ref Val | %RPD | RPDLimit | Qual |
| Chloride   |     | 4.74      |            | 0.1        | 5         | 0.1383      |      | 92.0          | 90                    | 110       | 0           |      |          |      |
| Sulfate    |     | 9.821     |            | 0.5        | 10        | 0           |      | 98.2          | 90                    | 110       | 0           |      |          |      |

|                       |           |           |              |            |           |             |      |               |          |           |             |      |          |      |
|-----------------------|-----------|-----------|--------------|------------|-----------|-------------|------|---------------|----------|-----------|-------------|------|----------|------|
| Sample ID             | 100ng lcs | Batch ID: | R13035       | Test Code: | SW8260B   | Units:      | µg/L | Analysis Date | 9/7/2004 | Prep Date |             |      |          |      |
| Client ID:            |           | Run ID:   | THOR_040907A | SeqNo:     | 303135    |             |      |               |          |           |             |      |          |      |
| Analyte               |           | Result    |              | PQL        | SPK value | SPK Ref Val |      | %REC          | LowLimit | HighLimit | RPD Ref Val | %RPD | RPDLimit | Qual |
| Benzene               |           | 20.98     |              | 1          | 20        | 0           |      | 105           | 75.3     | 128       | 0           |      |          |      |
| Toluene               |           | 19.56     |              | 1          | 20        | 0           |      | 97.8          | 77.8     | 122       | 0           |      |          |      |
| Chlorobenzene         |           | 19.87     |              | 1          | 20        | 0           |      | 99.3          | 76.2     | 130       | 0           |      |          |      |
| 1,1-Dichloroethene    |           | 18.38     |              | 1          | 20        | 0           |      | 91.9          | 70.2     | 119       | 0           |      |          |      |
| Trichloroethene (TCE) |           | 18.2      |              | 1          | 20        | 0           |      | 91.0          | 76.9     | 130       | 0           |      |          |      |

**Qualifiers:** ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits      B - Analyte detected in the associated Method Blank  
 J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits

21/28

**CLIENT:** San Juan Refining  
**Work Order:** 0409046  
**Project:** Inj Well 3rd Qtr 2004

**QC SUMMARY REPORT**  
 Laboratory Control Spike - generic

| Sample ID               | Batch ID | Test Code            | Units     | Analysis Date | Prep Date |          |           |             |      |          |      |
|-------------------------|----------|----------------------|-----------|---------------|-----------|----------|-----------|-------------|------|----------|------|
| Ics-6442                | 6442     | SW8270A              | µg/L      | 9/18/2004     | 9/7/2004  |          |           |             |      |          |      |
| Client ID:              |          | Run ID: ELMO_040918A |           | SeqNo: 306378 |           |          |           |             |      |          |      |
| Analyte                 | Result   | PQL                  | SPK value | SPK Ref Val   | %REC      | LowLimit | HighLimit | RPD Ref Val | %RPD | RPDLimit | Qual |
| Acenaphthene            | 60.94    | 10                   | 100       | 0             | 60.9      | 44.6     | 87.9      | 0           |      |          |      |
| 4-Chloro-3-methylphenol | 126      | 20                   | 200       | 0             | 63.0      | 46       | 92.2      | 0           |      |          |      |
| 2-Chlorophenol          | 126.8    | 10                   | 200       | 0             | 63.4      | 46.3     | 91.1      | 0           |      |          |      |
| 2,4-Dinitrotoluene      | 73       | 10                   | 100       | 0             | 73.0      | 53.2     | 94.9      | 0           |      |          |      |
| 4-Nitrophenol           | 69.28    | 50                   | 200       | 0             | 34.6      | 3.95     | 65.2      | 0           |      |          |      |
| Pentachlorophenol       | 116.6    | 50                   | 200       | 0             | 58.3      | 21.6     | 105       | 0           |      |          |      |
| Phenol                  | 80.96    | 10                   | 200       | 0             | 40.5      | 26.3     | 53.8      | 0           |      |          |      |
| 1,2,4-Trichlorobenzene  | 53.34    | 10                   | 100       | 0             | 53.3      | 37.1     | 90.7      | 0           |      |          |      |

| Sample ID               | Batch ID | Test Code            | Units     | Analysis Date | Prep Date |          |           |             |      |          |      |
|-------------------------|----------|----------------------|-----------|---------------|-----------|----------|-----------|-------------|------|----------|------|
| Icsd-6442               | 6442     | SW8270A              | µg/L      | 9/18/2004     | 9/7/2004  |          |           |             |      |          |      |
| Client ID:              |          | Run ID: ELMO_040918A |           | SeqNo: 306379 |           |          |           |             |      |          |      |
| Analyte                 | Result   | PQL                  | SPK value | SPK Ref Val   | %REC      | LowLimit | HighLimit | RPD Ref Val | %RPD | RPDLimit | Qual |
| Acenaphthene            | 66.3     | 10                   | 100       | 0             | 66.3      | 44.6     | 87.9      | 60.94       | 8.43 | 30.5     |      |
| 4-Chloro-3-methylphenol | 123.8    | 20                   | 200       | 0             | 61.9      | 46       | 92.2      | 126         | 1.70 | 28.6     |      |
| 2-Chlorophenol          | 122.4    | 10                   | 200       | 0             | 61.2      | 46.3     | 91.1      | 126.8       | 3.55 | 107      |      |
| 2,4-Dinitrotoluene      | 69.28    | 10                   | 100       | 0             | 69.3      | 53.2     | 94.9      | 73          | 5.23 | 14.7     |      |
| 4-Nitrophenol           | 72.64    | 50                   | 200       | 0             | 36.3      | 3.95     | 65.2      | 69.28       | 4.74 | 36.3     |      |
| Pentachlorophenol       | 122.6    | 50                   | 200       | 0             | 61.3      | 21.6     | 105       | 116.6       | 5.08 | 49       |      |
| Phenol                  | 75.16    | 10                   | 200       | 0             | 37.6      | 26.3     | 53.8      | 80.96       | 7.43 | 52.4     |      |
| 1,2,4-Trichlorobenzene  | 61.66    | 10                   | 100       | 0             | 61.7      | 37.1     | 90.7      | 53.34       | 14.5 | 36.4     |      |

**Qualifiers:** ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits      B - Analyte detected in the associated Method Blank  
 J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits

22/28

CLIENT: San Juan Refining  
 Work Order: 0409046  
 Project: Inj Well 3rd Qtr 2004

**QC SUMMARY REPORT**  
 Laboratory Control Spike - generic

| Sample ID                 | Ics-6493             | Batch ID: 6493 | Test Code: SW8270A | Units: µg/L | Analysis Date | 9/23/2004 | Prep Date | 9/16/2004   |      |          |      |
|---------------------------|----------------------|----------------|--------------------|-------------|---------------|-----------|-----------|-------------|------|----------|------|
| Client ID:                | Run ID: ELMO_040922A |                |                    | SeqNo:      | 308221        |           |           |             |      |          |      |
| Analyte                   | Result               | PQL            | SPK value          | SPK Ref Val | %REC          | LowLimit  | HighLimit | RPD Ref Val | %RPD | RPDLimit | Qual |
| Acenaphthene              | 69.28                | 10             | 100                | 0           | 69.3          | 44.6      | 87.9      | 0           |      |          |      |
| 4-Chloro-3-methylphenol   | 141.5                | 20             | 200                | 0           | 70.8          | 46        | 92.2      | 0           |      |          |      |
| 2-Chlorophenol            | 168.6                | 10             | 200                | 0           | 84.3          | 46.3      | 91.1      | 0           |      |          |      |
| 1,4-Dichlorobenzene       | 71.36                | 10             | 100                | 0           | 71.4          | 37.8      | 91.3      | 0           |      |          |      |
| 2,4-Dinitrotoluene        | 68                   | 10             | 100                | 0           | 68.0          | 53.2      | 94.9      | 0           |      |          |      |
| N-Nitrosodi-n-propylamine | 75.1                 | 10             | 100                | 0           | 75.1          | 46.2      | 82.6      | 0           |      |          |      |
| Phenol                    | 93.6                 | 10             | 200                | 0           | 46.8          | 26.3      | 53.8      | 0           |      |          |      |
| Pyrene                    | 89.64                | 15             | 100                | 0           | 89.6          | 53.8      | 94.5      | 0           |      |          |      |
| 1,2,4-Trichlorobenzene    | 65.08                | 10             | 100                | 0           | 65.1          | 37.1      | 90.7      | 0           |      |          |      |

| Sample ID                 | Icsd-6493            | Batch ID: 6493 | Test Code: SW8270A | Units: µg/L | Analysis Date | 9/23/2004 | Prep Date | 9/16/2004   |       |          |      |
|---------------------------|----------------------|----------------|--------------------|-------------|---------------|-----------|-----------|-------------|-------|----------|------|
| Client ID:                | Run ID: ELMO_040922A |                |                    | SeqNo:      | 308233        |           |           |             |       |          |      |
| Analyte                   | Result               | PQL            | SPK value          | SPK Ref Val | %REC          | LowLimit  | HighLimit | RPD Ref Val | %RPD  | RPDLimit | Qual |
| Acenaphthene              | 68.66                | 10             | 100                | 0           | 68.7          | 44.6      | 87.9      | 69.28       | 0.899 | 30.5     |      |
| 4-Chloro-3-methylphenol   | 133.2                | 20             | 200                | 0           | 66.6          | 46        | 92.2      | 141.5       | 6.01  | 28.6     |      |
| 2-Chlorophenol            | 157.8                | 10             | 200                | 0           | 78.9          | 46.3      | 91.1      | 168.6       | 6.63  | 107      |      |
| 1,4-Dichlorobenzene       | 75.86                | 10             | 100                | 0           | 75.9          | 37.8      | 91.3      | 71.36       | 6.11  | 62.1     |      |
| 2,4-Dinitrotoluene        | 65.52                | 10             | 100                | 0           | 65.5          | 53.2      | 94.9      | 68          | 3.71  | 14.7     |      |
| N-Nitrosodi-n-propylamine | 69.14                | 10             | 100                | 0           | 69.1          | 46.2      | 82.6      | 75.1        | 8.26  | 30.3     |      |
| Phenol                    | 86.82                | 10             | 200                | 0           | 43.4          | 26.3      | 53.8      | 93.6        | 7.52  | 52.4     |      |
| Pyrene                    | 85.4                 | 15             | 100                | 0           | 85.4          | 53.8      | 94.5      | 89.64       | 4.84  | 16.3     |      |
| 1,2,4-Trichlorobenzene    | 62.56                | 10             | 100                | 0           | 62.6          | 37.1      | 90.7      | 65.08       | 3.95  | 36.4     |      |

Qualifiers: ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits      B - Analyte detected in the associated Method Blank  
 J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits

23/28



**CLIENT:** San Juan Refining  
**Work Order:** 0409046  
**Project:** Inj Well 3rd Qtr 2004

**QC SUMMARY REPORT**  
 Laboratory Control Spike - generic

|            |                 |           |                         |             |               |          |             |               |                 |           |                 |
|------------|-----------------|-----------|-------------------------|-------------|---------------|----------|-------------|---------------|-----------------|-----------|-----------------|
| Sample ID  | <b>LCS-6436</b> | Batch ID: | <b>6436</b>             | Test Code:  | <b>SW7470</b> | Units:   | <b>mg/L</b> | Analysis Date | <b>9/7/2004</b> | Prep Date | <b>9/7/2004</b> |
| Client ID: |                 | Run ID:   | <b>MI-LA254_040907A</b> | SeqNo:      | <b>302915</b> |          |             |               |                 |           |                 |
| Analyte    | Result          | PQL       | SPK value               | SPK Ref Val | %REC          | LowLimit | HighLimit   | RPD Ref Val   | %RPD            | RPDLimit  | Qual            |
| Mercury    | 0.005726        | 0.0002    | 0.005                   | 0.0001224   | 112           | 75.2     | 134         | 0             |                 |           |                 |

|            |                 |           |                         |             |               |          |             |               |                 |           |                 |
|------------|-----------------|-----------|-------------------------|-------------|---------------|----------|-------------|---------------|-----------------|-----------|-----------------|
| Sample ID  | <b>LCS-6439</b> | Batch ID: | <b>6436</b>             | Test Code:  | <b>SW7470</b> | Units:   | <b>mg/L</b> | Analysis Date | <b>9/7/2004</b> | Prep Date | <b>9/7/2004</b> |
| Client ID: |                 | Run ID:   | <b>MI-LA254_040907A</b> | SeqNo:      | <b>302927</b> |          |             |               |                 |           |                 |
| Analyte    | Result          | PQL       | SPK value               | SPK Ref Val | %REC          | LowLimit | HighLimit   | RPD Ref Val   | %RPD            | RPDLimit  | Qual            |
| Mercury    | 0.005047        | 0.0002    | 0.005                   | 0.0001224   | 98.5          | 75.2     | 134         | 0.005726      | 12.6            | 0         |                 |

|            |                 |           |                    |             |                |          |             |               |                             |           |                  |
|------------|-----------------|-----------|--------------------|-------------|----------------|----------|-------------|---------------|-----------------------------|-----------|------------------|
| Sample ID  | <b>LCS-6492</b> | Batch ID: | <b>6492</b>        | Test Code:  | <b>SW6010A</b> | Units:   | <b>mg/L</b> | Analysis Date | <b>9/17/2004 2:55:08 PM</b> | Prep Date | <b>9/16/2004</b> |
| Client ID: |                 | Run ID:   | <b>ICP_040917A</b> | SeqNo:      | <b>306648</b>  |          |             |               |                             |           |                  |
| Analyte    | Result          | PQL       | SPK value          | SPK Ref Val | %REC           | LowLimit | HighLimit   | RPD Ref Val   | %RPD                        | RPDLimit  | Qual             |
| Arsenic    | 0.5026          | 0.02      | 0.5                | 0           | 101            | 80       | 120         | 0             |                             |           |                  |
| Barium     | 0.4711          | 0.02      | 0.5                | 0           | 94.2           | 80       | 120         | 0             |                             |           |                  |
| Cadmium    | 0.4858          | 0.002     | 0.5                | 0.001101    | 96.9           | 80       | 120         | 0             |                             |           |                  |
| Calcium    | 44.51           | 1         | 50                 | 0           | 89.0           | 80       | 120         | 0             |                             |           |                  |
| Chromium   | 0.4789          | 0.006     | 0.5                | 0           | 95.8           | 80       | 120         | 0             |                             |           |                  |
| Lead       | 0.4833          | 0.005     | 0.5                | 0.003158    | 96.0           | 80       | 120         | 0             |                             |           |                  |
| Magnesium  | 46.73           | 1         | 50                 | 0           | 93.5           | 80       | 120         | 0             |                             |           |                  |
| Potassium  | 48.02           | 1         | 50                 | 0           | 96.0           | 80       | 120         | 0             |                             |           |                  |
| Selenium   | 0.4785          | 0.05      | 0.5                | 0           | 95.7           | 80       | 120         | 0             |                             |           |                  |
| Sodium     | 51.27           | 1         | 50                 | 0           | 103            | 80       | 120         | 0             |                             |           |                  |

24/28

**Qualifiers:** ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits      B - Analyte detected in the associated Method Blank  
 J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits

**CLIENT:** San Juan Refining  
**Work Order:** 0409046  
**Project:** Inj Well 3rd Qtr 2004

**QC SUMMARY REPORT**  
 Laboratory Control Spike Duplicate

| Sample ID  | Batch ID | Test Code | Units       | Analysis Date        | Prep Date |          |           |             |       |          |      |
|------------|----------|-----------|-------------|----------------------|-----------|----------|-----------|-------------|-------|----------|------|
| LCS-6492   | 6492     | SW6010A   | mg/L        | 9/17/2004 2:57:19 PM | 9/16/2004 |          |           |             |       |          |      |
| Client ID: |          | Run ID:   | ICP_040917A | SeqNo:               | 306649    |          |           |             |       |          |      |
| Analyte    | Result   | PQL       | SPK value   | SPK Ref Val          | %REC      | LowLimit | HighLimit | RPD Ref Val | %RPD  | RPDLimit | Qual |
| Arsenic    | 0.4921   | 0.02      | 0.5         | 0                    | 98.4      | 80       | 120       | 0.5026      | 2.13  | 20       |      |
| Barium     | 0.4702   | 0.02      | 0.5         | 0                    | 94.0      | 80       | 120       | 0.4711      | 0.206 | 20       |      |
| Cadmium    | 0.4836   | 0.002     | 0.5         | 0.001101             | 96.5      | 80       | 120       | 0.4858      | 0.465 | 20       |      |
| Calcium    | 45.37    | 1         | 50          | 0                    | 90.7      | 80       | 120       | 44.51       | 1.91  | 20       |      |
| Chromium   | 0.4754   | 0.006     | 0.5         | 0                    | 95.1      | 80       | 120       | 0.4789      | 0.732 | 20       |      |
| Lead       | 0.4817   | 0.005     | 0.5         | 0.003158             | 95.7      | 80       | 120       | 0.4833      | 0.323 | 20       |      |
| Magnesium  | 47.56    | 1         | 50          | 0                    | 95.1      | 80       | 120       | 46.73       | 1.76  | 20       |      |
| Potassium  | 48.69    | 1         | 50          | 0                    | 97.4      | 80       | 120       | 48.02       | 1.38  | 20       |      |
| Selenium   | 0.4981   | 0.05      | 0.5         | 0                    | 99.6      | 80       | 120       | 0.4785      | 4.01  | 20       |      |
| Sodium     | 51.74    | 1         | 50          | 0                    | 103       | 80       | 120       | 51.27       | 0.904 | 20       |      |

| Sample ID  | Batch ID | Test Code | Units       | Analysis Date         | Prep Date |          |           |             |      |          |      |
|------------|----------|-----------|-------------|-----------------------|-----------|----------|-----------|-------------|------|----------|------|
| LCS-6492   | 6492     | SW6010A   | mg/L        | 9/20/2004 10:46:14 AM | 9/16/2004 |          |           |             |      |          |      |
| Client ID: |          | Run ID:   | ICP_040920A | SeqNo:                | 306667    |          |           |             |      |          |      |
| Analyte    | Result   | PQL       | SPK value   | SPK Ref Val           | %REC      | LowLimit | HighLimit | RPD Ref Val | %RPD | RPDLimit | Qual |
| Arsenic    | 0.5658   | 0.02      | 0.5         | 0                     | 113       | 80       | 120       | 0           |      |          |      |
| Barium     | 0.5299   | 0.02      | 0.5         | 0                     | 106       | 80       | 120       | 0           |      |          |      |
| Cadmium    | 0.5455   | 0.002     | 0.5         | 0                     | 109       | 80       | 120       | 0           |      |          |      |
| Calcium    | 44.8     | 1         | 50          | 0                     | 89.6      | 80       | 120       | 0           |      |          |      |
| Chromium   | 0.533    | 0.006     | 0.5         | 0                     | 107       | 80       | 120       | 0           |      |          |      |
| Lead       | 0.5396   | 0.005     | 0.5         | 0                     | 108       | 80       | 120       | 0           |      |          |      |
| Magnesium  | 47.02    | 1         | 50          | 0                     | 94.0      | 80       | 120       | 0           |      |          |      |
| Potassium  | 47.87    | 1         | 50          | 0                     | 95.7      | 80       | 120       | 0           |      |          |      |
| Selenium   | 0.5398   | 0.05      | 0.5         | 0                     | 108       | 80       | 120       | 0           |      |          |      |
| Sodium     | 51.29    | 1         | 50          | 0                     | 103       | 80       | 120       | 0           |      |          |      |

**Qualifiers:** ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits      B - Analyte detected in the associated Method Blank  
 J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits

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CLIENT: San Juan Refining  
 Work Order: 0409046  
 Project: Inj Well 3rd Qtr 2004

**QC SUMMARY REPORT**  
 Laboratory Control Spike Duplicate

Sample ID **LCSD-6492** Batch ID: **6492** Test Code: **SW6010A** Units: **mg/L** Analysis Date **9/20/2004 10:49:14 AM** Prep Date **9/16/2004**  
 Client ID: Run ID: **ICP\_040920A** SeqNo: **306668**

| Analyte   | Result | PQL   | SPK value | SPK Ref Val | %REC | LowLimit | HighLimit | RPD Ref Val | %RPD  | RPDLimit | Qual |
|-----------|--------|-------|-----------|-------------|------|----------|-----------|-------------|-------|----------|------|
| Arsenic   | 0.568  | 0.02  | 0.5       | 0           | 114  | 80       | 120       | 0.5658      | 0.394 | 20       |      |
| Barium    | 0.5279 | 0.02  | 0.5       | 0           | 106  | 80       | 120       | 0.5299      | 0.391 | 20       |      |
| Cadmium   | 0.5408 | 0.002 | 0.5       | 0           | 108  | 80       | 120       | 0.5455      | 0.862 | 20       |      |
| Calcium   | 45.48  | 1     | 50        | 0           | 91.0 | 80       | 120       | 44.8        | 1.51  | 20       |      |
| Chromium  | 0.5294 | 0.006 | 0.5       | 0           | 106  | 80       | 120       | 0.533       | 0.673 | 20       |      |
| Lead      | 0.5329 | 0.005 | 0.5       | 0           | 107  | 80       | 120       | 0.5396      | 1.26  | 20       |      |
| Magnesium | 47.89  | 1     | 50        | 0           | 95.8 | 80       | 120       | 47.02       | 1.85  | 20       |      |
| Potassium | 48.94  | 1     | 50        | 0           | 97.9 | 80       | 120       | 47.87       | 2.21  | 20       |      |
| Selenium  | 0.5307 | 0.05  | 0.5       | 0           | 106  | 80       | 120       | 0.5398      | 1.69  | 20       |      |
| Sodium    | 52.2   | 1     | 50        | 0           | 104  | 80       | 120       | 51.29       | 1.76  | 20       |      |

Sample ID **LCS-6492** Batch ID: **6492** Test Code: **SW6010A** Units: **mg/L** Analysis Date **9/20/2004 12:34:15 PM** Prep Date **9/16/2004**  
 Client ID: Run ID: **ICP\_040920A** SeqNo: **306683**

| Analyte | Result | PQL   | SPK value | SPK Ref Val | %REC | LowLimit | HighLimit | RPD Ref Val | %RPD | RPDLimit | Qual |
|---------|--------|-------|-----------|-------------|------|----------|-----------|-------------|------|----------|------|
| Silver  | 0.5907 | 0.005 | 0.5       | 0           | 118  | 80       | 120       | 0           |      |          |      |

Sample ID **LCSD-6492** Batch ID: **6492** Test Code: **SW6010A** Units: **mg/L** Analysis Date **9/20/2004 12:36:33 PM** Prep Date **9/16/2004**  
 Client ID: Run ID: **ICP\_040920A** SeqNo: **306684**

| Analyte | Result | PQL   | SPK value | SPK Ref Val | %REC | LowLimit | HighLimit | RPD Ref Val | %RPD | RPDLimit | Qual |
|---------|--------|-------|-----------|-------------|------|----------|-----------|-------------|------|----------|------|
| Silver  | 0.5705 | 0.005 | 0.5       | 0           | 114  | 80       | 120       | 0.5907      | 3.48 | 20       |      |

Sample ID **LCS-6492** Batch ID: **6492** Test Code: **SW6010A** Units: **mg/L** Analysis Date **9/20/2004 8:07:23 AM** Prep Date **9/16/2004**  
 Client ID: Run ID: **ICP\_040920E** SeqNo: **307232**

| Analyte | Result | PQL   | SPK value | SPK Ref Val | %REC | LowLimit | HighLimit | RPD Ref Val | %RPD | RPDLimit | Qual |
|---------|--------|-------|-----------|-------------|------|----------|-----------|-------------|------|----------|------|
| Silver  | 0.498  | 0.005 | 0.5       | 0           | 99.6 | 80       | 120       | 0           |      |          |      |

**Qualifiers:** ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank  
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits

26/28

CLIENT: San Juan Refining  
Work Order: 0409046  
Project: Inj Well 3rd Qtr 2004

**QC SUMMARY REPORT**  
Laboratory Control Spike Duplicate

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|            |           |           |             |            |             |        |          |               |                      |           |           |      |
|------------|-----------|-----------|-------------|------------|-------------|--------|----------|---------------|----------------------|-----------|-----------|------|
| Sample ID  | LCSD-6492 | Batch ID: | 6492        | Test Code: | SW6010A     | Units: | mg/L     | Analysis Date | 9/20/2004 8:09:28 AM | Prep Date | 9/16/2004 |      |
| Client ID: |           | Run ID:   | ICP_040920E | SeqNo:     | 307233      |        |          |               |                      |           |           |      |
| Analyte    |           | Result    | PQL         | SPK value  | SPK Ref Val | %REC   | LowLimit | HighLimit     | RPD Ref Val          | %RPD      | RPDLimit  | Qual |
| Silver     |           | 0.4968    | 0.005       | 0.5        | 0           | 99.4   | 80       | 120           | 0.498                | 0.251     | 20        |      |

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Qualifiers: ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits      B - Analyte detected in the associated Method Blank  
                  J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits

# Hall Environmental Analysis Laboratory

## Sample Receipt Checklist

Client Name SJR

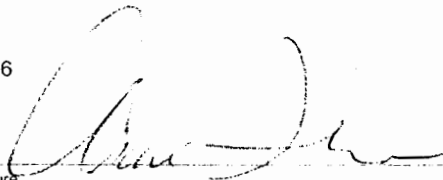
Date and Time Received:

9/3/2004

Work Order Number 0409046

Received by AT

Checklist completed by

  
Signature

9/3/04  
Date

Matrix

Carrier name UPS

- Shipping container/cooler in good condition? Yes  No  Not Present
- Custody seals intact on shipping container/cooler? Yes  No  Not Present  Not Shipped
- Custody seals intact on sample bottles? Yes  No  N/A
- Chain of custody present? Yes  No
- Chain of custody signed when relinquished and received? Yes  No
- Chain of custody agrees with sample labels? Yes  No
- Samples in proper container/bottle? Yes  No
- Sample containers intact? Yes  No
- Sufficient sample volume for indicated test? Yes  No
- All samples received within holding time? Yes  No
- Water - VOA vials have zero headspace? No VOA vials submitted  Yes  No
- Water - pH acceptable upon receipt? Yes  No  N/A

Container/Temp Blank temperature?

**16°**

*4° C ± 2 Acceptable*

If given sufficient time to cool.

COMMENTS:

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Client contacted \_\_\_\_\_ Date contacted: \_\_\_\_\_ Person contacted \_\_\_\_\_

Contacted by: \_\_\_\_\_ Regarding \_\_\_\_\_

Comments: \_\_\_\_\_

\_\_\_\_\_

\_\_\_\_\_

\_\_\_\_\_

\_\_\_\_\_

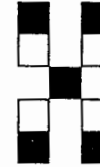
Corrective Action \_\_\_\_\_

\_\_\_\_\_

\_\_\_\_\_

# CHAIN-OF-CUSTODY RECORD

Accreditation Applied:  
 NELAC  USACE



**HALL ENVIRONMENTAL ANALYSIS LABORATORY**  
 4901 Hawkins NE, Suite D  
 Albuquerque, New Mexico 87109  
 Tel. 505.345.3975 Fax 505.345.4107  
 www.hallenvironmental.com

Client: SAN Juan Refining

Project Name: Injection Well 3<sup>rd</sup> Qtr 2004

Address: #50 Rd 4990  
Bloomfield, NM  
87413

Project #:

Project Manager:

Phone #: 505-632-4161

Sampler: Cindy Hurtado

Fax #: 505-632-3911

Sample Temperature: 16°

## ANALYSIS REQUEST

| Date    | Time  | Matrix | Sample I.D. No.                | Number/Volume | Preservative      |                  | HEAL No.   |
|---------|-------|--------|--------------------------------|---------------|-------------------|------------------|------------|
|         |       |        |                                |               | HgCl <sub>2</sub> | HNO <sub>3</sub> |            |
| 9-02-04 | 920AM | H2O    | Injection Well 3 <sup>rd</sup> | 3-VOA         | X                 |                  | 0409246.1  |
|         |       |        |                                | 1-liter       |                   |                  | Amber      |
|         |       |        |                                | 1-500         |                   | X                |            |
|         |       |        |                                | 1 liter       |                   |                  |            |
|         |       |        |                                | 1-250         |                   |                  | Na OH      |
|         |       |        |                                | 1-250         |                   |                  | Zn Acetate |
|         |       |        | Trip Blank                     |               |                   |                  | -2         |

| BTEX + MTBE + TMB's (8021) | BTEX + MTBE + TPH (Gasoline Only) | TPH Method 8015B (Gas/Diesel) | TPH (Method 418.1) | EDB (Method 504.1) | EDC (Method 8021) | 8310 (PNA or PAH) | PCRA 8 Metals, Na, K, Mg, Ca | Anions (F, Cl, NO <sub>3</sub> , NO <sub>2</sub> , PO <sub>4</sub> , SO <sub>4</sub> ) | 8081 Pesticides / PCB's (8082) | 8260B (VOA) | 8270 (Semi-VOA) | EC, PH, SO <sub>4</sub> , ALK, Cl | Ignitability, Corrosivity | Reactivity | Sulfide | Air Bubbles or Headspace (Y or N) |
|----------------------------|-----------------------------------|-------------------------------|--------------------|--------------------|-------------------|-------------------|------------------------------|--|--------------------------------|-------------|-----------------|-----------------------------------|---------------------------|------------|---------|-----------------------------------|
|                            |                                   |                               |                    |                    |                   |                   |                              |  |                                | X           |                 |                                   |                           |            |         |                                   |
|                            |                                   |                               |                    |                    |                   |                   |                              |  |                                |             | X               |                                   |                           |            |         |                                   |
|                            |                                   |                               |                    |                    |                   |                   | X                            |  |                                |             |                 | X                                 | X                         |            |         |                                   |
|                            |                                   |                               |                    |                    |                   |                   |                              |  |                                |             |                 |                                   | X                         |            |         |                                   |
|                            |                                   |                               |                    |                    |                   |                   |                              |  |                                |             |                 |                                   |                           | X          |         |                                   |

Date: 02-04 Time: 10AM Relinquished By: (Signature) Cindy Hurtado  
 Received By: (Signature) [Signature] 9/13/04  
 Relinquished By: (Signature) \_\_\_\_\_ Received By: (Signature) \_\_\_\_\_  
 1602

Remarks: