# Chemical Analysis Report

**WQAP-2022-08-25-03**

Florida DEP Laboratory
2600 Blair Stone Road
Tallahassee, FL 32399-2400
DOH Accreditation E31780

Event Description: **Algal Bloom Response -SOROC**
Request ID: **RQ-2022-02-28-44**
Customer: **WQAP**
Project ID: **BLOOM-RESP**

Send Reports to:
FL Dept. of Environmental Protection
2295 Victoria AVE
Suite 364
Fort Myers, FL 33901
Attn: Kirby Wolfe

For additional information please contact
Colin Wright, Ph.D.
Liang-Tsair Lin, Ph.D.
Kerry Tate, Ph.D.
Dr. rer. nat. Bettina Steinbock
Thekkekalathil Chandrasekhar, Ph.D, QA Officer
Phone (850) 245-8085

Certified by: Colin Wright, Program Administrator

Date Certified: 13-SEP-2022 14:39

## Case Narrative

Unless otherwise noted, all samples included in this report were received in accordance with protocols referenced in Chapter 62-160, Florida Administrative Code (F.A.C.). Results published in this report pertain only to the samples as submitted to, and received by the laboratory. All times in this report are adjusted to the applicable Eastern Time Zone (EST or EDT).

Results for the following analytical groups are included in this report: Nutrients, Pesticides and Priority Organic Pollutants.

Scientific notation may be used in reporting very large or small values. Values reported using scientific notation will take the form of the following example: 1.3E+03, which is equivalent to $1.3 \times 10^3$ or 1300.

Unless otherwise noted, analytical values for soil and sediment samples are reported on a dry weight basis, and analytical values for waste and tissue samples are reported on a wet weight basis.

Results for TNI accredited tests met requirements established by The NELAC Institute. A double asterisk (**) is used to indicate an analyte/matrix/method for which the laboratory is not TNI accredited by the Florida Department of Health Environmental Laboratory Certification Program or where accreditation for that field of testing is not applicable.

Any significant anomalies or deviations from established protocols are documented in Non-Conformance Reports, which, where appropriate, are included within this analytical report. Additional comments related to specific analytical tests may be included as remarks following the analytical results for each sample. Such comments and remarks are for informational purposes only and are not intended to convey judgement about the usability of the reported data.

A quality control report on the performance of the test method for the submitted samples is included. Uncertainty associated with the analytical results contained in this report can be estimated from the reported quality assurance results and from published quality control acceptance limits for each analytical test. Matrix quality control results (matrix spike recoveries and matrix sample precision) pertain only to the matrix sample tested and do not necessarily reflect test method performance for other samples.

Typical matrix quality control (QC) measurements may include matrix spike recovery, matrix spike duplicate recovery, matrix spike precision and matrix sample precision. Not all matrix QC results may be available or reportable; where they are not an explanation is provided. Typical reasons for unavailable QC results include, but are not limited to, a) insufficient matrix sample to perform some or all QC measurements; b) analyte concentration in the sample replicated was too low for a meaningful measurement of precision and c) analyte concentration in the matrix sample spiked was too high (relative to the amount of analyte spiked) for a meaningful measurement of recovery. Where matrix QC results are unavailable, other method performance metrics (e.g., LCS recovery, LCS precision, surrogate recovery) may be used to assess performance of the method. Comments explaining any missing QC measurements are not intended to convey any adverse conclusions about the quality of the reported data.

Precision is reported as relative percent difference unless otherwise noted.

Quality Control codes as defined below may be used in this report to indicate results that are associated with one or more quality control elements which did not fall within established test method criteria. Such results may be qualified as estimates using a J qualifier as required by 62-160 F.A.C. Explanations are included in the report for any results that were reported as estimates for other reasons.

QC Codes used in this report may include:

- LCS – Recovery for the batch Laboratory Control Sample (LCS) was outside existing control limits;
- MS – Recovery for the batch matrix spike (MS) was outside existing control limits;
- CCV – Recovery for a continuing calibration verification (CCV) standard was outside existing control limits;
- SUR – Recovery of a surrogate (SUR) for associated analytes was outside existing control limits;
- RPD – The precision, measured as relative percent difference (RPD), of batch replicate measurements was outside existing control limits;
- RSD – The precision, measured as relative standard deviation (RSD), of batch replicate measurements was outside existing control limits;
- SMP – Sample - used precision derived from replicate analyses of a sample;

The following data qualifiers are used, where applicable, in this report as specified in 62-160 F.A.C.

- A - Value reported is the mean of two or more determinations.
- B - Results based on colony counts outside the acceptable range.
- I - The reported value is between the laboratory method detection limit and the laboratory practical quantitation limit.
- J - Estimated value and/or the analysis did not meet established quality control criteria.
- K - Actual value is known to be less than value given.
- L - Actual value is known to be greater than value given.
- N - Presumptive evidence of presence of material.
- O - Sampled, but analysis lost or not performed.
- Q - Sample held beyond normal holding time.
- T - Value reported is less than the criterion of detection.
- U - Material was analyzed for but not detected. The reported value is the method detection limit for the sample analyzed.
- V - Analyte was detected in both sample and method blank.
- X - Too few individuals to calculate SCI value.
- Y - The laboratory analysis was from an unpreserved or improperly preserved sample. The data may not be accurate.
- Z - Colonies were too numerous to count (TNTC).

Quality control information from overflow laboratories may not be included in this report. Please refer to the associated report from the overflow laboratory for additional information.

**Sample Location: Lake Clay** **Collection Date/Time: 08/24/2022 14:58**

**Field ID: HAB-SD-082422-1505** Matrix: W-SURF-FRH

| Sample ID | Ref. Method | Component | Result | Code | Units | Batch ID | QC Codes |
|---|---|---|---|---|---|---|---|
| 2352292 | EPA 8270E | Acenaphthene | 0.026 | U | ug/L | P418679 | |
| | | Acenaphthylene | 0.026 | U | ug/L | P418679 | |
| | | Acetophenone | 0.21 | U | ug/L | P418679 | |
| | | 2-Acetylaminofluorene | 1.0 | U | ug/L | P418679 | |
| | | 4-Aminobiphenyl | 4.2 | U | ug/L | P418679 | |
| | | Aniline | 1.0 | U | ug/L | P418679 | |
| | | Anthracene | 0.052 | U | ug/L | P418679 | |
| | | Azobenzene/1,2-Diphenylhydrazine** | 0.052 | U | ug/L | P418679 | |
| | | Benzidine | 10 | U | ug/L | P418679 | |
| | | Benzo(a)anthracene | 0.026 | U | ug/L | P418679 | |
| | | Benzo(a)pyrene | 0.026 | U | ug/L | P418679 | |
| | | Benzo(b)fluoranthene | 0.026 | U | ug/L | P418679 | |
| | | Benzo(k)fluoranthene | 0.026 | U | ug/L | P418679 | |
| | | Benzo(g,h,i)perylene | 0.026 | U | ug/L | P418679 | |
| | | Benzyl alcohol | 0.21 | U | ug/L | P418679 | |
| | | Bis(2-chloroethoxy)methane | 0.052 | U | ug/L | P418679 | |
| | | Bis(2-chloroethyl)ether | 0.052 | U | ug/L | P418679 | |
| | | Bis(2-chloroisopropyl)ether | 0.052 | U | ug/L | P418679 | |
| | | Bis(2-ethylhexyl)phthalate | 5.2 | U | ug/L | P418679 | |
| | | Butyl benzyl phthalate | 1.0 | U | ug/L | P418679 | |
| | | 4-Bromophenyl phenyl ether | 0.052 | U | ug/L | P418679 | |
| | | 2-Chloronaphthalene | 0.052 | U | ug/L | P418679 | |
| | | 4-Chlorophenyl phenyl ether | 0.052 | U | ug/L | P418679 | |
| | | Carbazole | 0.052 | U | ug/L | P418679 | |
| | | Chrysene | 0.026 | U | ug/L | P418679 | |
| | | m,p-Cresols | 0.052 | U | ug/L | P418679 | |
| | | o-Cresol | 0.052 | U | ug/L | P418679 | |
| | | Di-n-butyl phthalate | 2.1 | U | ug/L | P418679 | |
| | | Di-n-octyl phthalate | 0.052 | U | ug/L | P418679 | |
| | | Dibenzo(a,h)anthracene | 0.026 | U | ug/L | P418679 | |
| | | Dibenzofuran | 0.052 | U | ug/L | P418679 | |
| | | 3,3'-Dichlorobenzidine | 10 | U | ug/L | P418679 | |
| | | Diethyl phthalate | 2.1 | U | ug/L | P418679 | |
| | | Dimethyl phthalate | 0.21 | U | ug/L | P418679 | |
| | | Dimethylaminoazobenzene | 0.052 | U | ug/L | P418679 | |
| | | 7,12-Dimethylbenz(a)anthracene | 0.10 | U | ug/L | P418679 | |
| | | 1,3-Dinitrobenzene | 0.21 | U | ug/L | P418679 | |
| | | 2,4-Dinitrotoluene | 0.052 | U | ug/L | P418679 | |
| | | 2,6-Dinitrotoluene | 0.052 | U | ug/L | P418679 | |
| | | Dinoseb** | 4.2 | U | ug/L | P418679 | |
| | | Ethyl methanesulfonate | 1.0 | U | ug/L | P418679 | |
| | | Fluoranthene | 0.052 | U | ug/L | P418679 | |
| | | Fluorene | 0.026 | U | ug/L | P418679 | |
| | | Hexachlorobenzene | 0.052 | U | ug/L | P418679 | |

**Field ID: HAB-SD-082422-1505** Matrix: W-SURF-FRH

| Sample ID | Ref. Method | Component | Result | Code | Units | Batch ID | QC Codes |
|---|---|---|---|---|---|---|---|
| 2352292 | EPA 8270E | Hexachlorobutadiene | 0.052 | U | ug/L | P418679 | |
| | | Hexachlorocyclopentadiene | 0.052 | U | ug/L | P418679 | |
| | | Hexachloroethane | 0.052 | U | ug/L | P418679 | |
| | | Hexachloropropene | 0.052 | U | ug/L | P418679 | |
| | | Indeno(1,2,3-cd)pyrene | 0.026 | U | ug/L | P418679 | |
| | | Isophorone | 0.052 | U | ug/L | P418679 | |
| | | Isosafrole | 0.052 | U | ug/L | P418679 | |
| | | 3-Methylcholanthrene | 0.10 | U | ug/L | P418679 | |
| | | 2-Methylnaphthalene | 0.10 | U | ug/L | P418679 | |
| | | Naphthalene | 0.10 | U | ug/L | P418679 | |
| | | 1-Naphthylamine | 10 | U | ug/L | P418679 | |
| | | 2-Naphthylamine | 10 | U | ug/L | P418679 | |
| | | 2-Nitroaniline | 0.052 | U | ug/L | P418679 | |
| | | Nitrobenzene | 0.052 | U | ug/L | P418679 | |
| | | 5-Nitro-o-toluidine | 0.10 | U | ug/L | P418679 | |
| | | N-Nitrosodi-n-butylamine | 0.052 | U | ug/L | P418679 | |
| | | N-Nitrosodiethylamine | 1.0 | U | ug/L | P418679 | |
| | | N-Nitrosodimethylamine | 2.1 | U | ug/L | P418679 | |
| | | N-Nitrosodi-n-propylamine | 0.052 | U | ug/L | P418679 | |
| | | N-Nitrosomethylethylamine | 2.1 | U | ug/L | P418679 | |
| | | N-Nitrosomorpholine | 0.052 | U | ug/L | P418679 | |
| | | N-Nitrosopiperidine | 0.052 | U | ug/L | P418679 | |
| | | N-Nitrosopyrrolidine | 0.10 | U | ug/L | P418679 | |
| | | Pentachlorobenzene | 0.052 | U | ug/L | P418679 | |
| | | Pentachloroethane | 0.052 | U | ug/L | P418679 | |
| | | Pentachloronitrobenzene | 0.052 | U | ug/L | P418679 | |
| | | Phenacetin | 1.0 | U | ug/L | P418679 | |
| | | Phenanthrene | 0.10 | U | ug/L | P418679 | |
| | | 2-Picoline | 1.0 | U | ug/L | P418679 | |
| | | Pyrene | 0.10 | U | ug/L | P418679 | |
| | | Pyridine | 4.2 | U | ug/L | P418679 | |
| | | Safrole | 0.052 | U | ug/L | P418679 | |
| | | 1,2,4,5-Tetrachlorobenzene | 0.052 | U | ug/L | P418679 | |
| | | o-Toluidine | 0.10 | U | ug/L | P418679 | |
| | | 1,2,4-Trichlorobenzene | 0.052 | U | ug/L | P418679 | |
| | | 1,3,5-Trinitrobenzene | 0.21 | U | ug/L | P418679 | |
| | | 4-Chloro-3-methylphenol | 0.052 | U | ug/L | P418679 | |
| | | 2-Chlorophenol | 0.052 | U | ug/L | P418679 | |
| | | 2,4-Dichlorophenol | 0.052 | U | ug/L | P418679 | |
| | | 2,6-Dichlorophenol | 0.052 | U | ug/L | P418679 | |
| | | 2,4-Dimethylphenol | 0.052 | U | ug/L | P418679 | |
| | | 2,4-Dinitrophenol | 10 | U | ug/L | P418679 | |
| | | 2-Methyl-4,6-dinitrophenol | 3.1 | U | ug/L | P418679 | |
| | | 2-Nitrophenol | 0.052 | U | ug/L | P418679 | |
| | | 4-Nitrophenol | 10 | U | ug/L | P418679 | |

**Field ID: HAB-SD-082422-1505** Matrix: W-SURF-FRH

| Sample ID | Ref. Method | Component | Result | Code | Units | Batch ID | QC Codes |
|---|---|---|---|---|---|---|---|
| 2352292 | EPA 8270E | Pentachlorophenol | 0.52 | U | ug/L | P418679 | |
| | | Phenol | 0.052 | U | ug/L | P418679 | |
| | | 2,3,4,6-Tetrachlorophenol | 0.10 | U | ug/L | P418679 | |
| | | 2,4,5-Trichlorophenol | 0.052 | U | ug/L | P418679 | |
| | | 2,4,6-Trichlorophenol | 0.052 | U | ug/L | P418679 | |
| | | 1-Methylnaphthalene | 0.10 | U | ug/L | P418679 | |
| 2352293 | FL-PRO 2018 | TRPH | 0.67 | I J | mg/L | P418756 | MS, RPD |
| 2352294 | EPA 8321B | Anatoxin-a | 0.25 | U | ug/L | P418570 | |
| | | Cylindrospermopsin | 0.10 | U | ug/L | P418569 | |
| | | Desmethyl microcystin LR | 0.25 | U | ug/L | P418569 | |
| | | Microcystin HilR | 0.25 | U | ug/L | P418569 | |
| | | Microcystin HtyR | 0.25 | U | ug/L | P418569 | |
| | | Microcystin LA | 0.10 | U | ug/L | P418569 | |
| | | Microcystin LF | 0.10 | U | ug/L | P418569 | |
| | | Microcystin LR | 0.25 | U | ug/L | P418569 | |
| | | Microcystin LW | 0.25 | U | ug/L | P418569 | |
| | | Microcystin LY | 0.10 | U | ug/L | P418569 | |
| | | Microcystin RR | 0.10 | U | ug/L | P418569 | |
| | | Microcystin WR | 0.50 | U | ug/L | P418569 | |
| | | Microcystin YR | 0.25 | U | ug/L | P418569 | |
| | | Nodularin-R | 0.10 | U | ug/L | P418569 | |
| 2352296 | EPA 350.1 Rev. 2.0 | Ammonia-N | 0.004 | I | mg N/L | P418886 | |
| | EPA 351.2 Rev. 2.0 | Kjeldahl Nitrogen | 0.99 | | mg N/L | P418823 | |
| | EPA 353.2 Rev. 2.0 | NO2NO3-N | 0.004 | U | mg N/L | P419050 | |
| | EPA 365.1 Rev. 2.0 | Total-P | 0.018 | | mg P/L | P419131 | |
| 2352297 | EPA 365.1 Rev. 2.0 dissolved | O-Phosphate-P | 0.004 | U | mg P/L | P418697 | |

Ref. Method and Comment:
EPA 8270E: Precision for benzidine is not available due to low recoveries in the duplicate matrix spikes.
FL-PRO 2018: Refer to the Lab Analysis Report for an explanation of QC Codes.

# Quality Assurance Report
# Method Blank Results

**Reference Method: EPA 350.1 Rev. 2.0**
**Batch ID: P418886**

| Component | Result | Code | Units |
|---|---|---|---|
| Ammonia-N | 0.002 | U | mg N/L |

**Reference Method: EPA 351.2 Rev. 2.0**
**Batch ID: P418823**

| Component | Result | Code | Units |
|---|---|---|---|
| Kjeldahl Nitrogen | 0.080 | U | mg N/L |

**Reference Method: EPA 353.2 Rev. 2.0**
**Batch ID: P419050**

| Component | Result | Code | Units |
|---|---|---|---|
| NO2NO3-N | 0.004 | U | mg N/L |

**Reference Method: EPA 365.1 Rev. 2.0**
**Batch ID: P419131**

| Component | Result | Code | Units |
|---|---|---|---|
| Total-P | 0.002 | U | mg P/L |

**Reference Method: EPA 365.1 Rev. 2.0 dissolved**
**Batch ID: P418697**

| Component | Result | Code | Units |
|---|---|---|---|
| O-Phosphate-P | 0.004 | U | mg P/L |

**Reference Method: EPA 8270E**
**Batch ID: P418679**

| Component | Result | Code | Units |
|---|---|---|---|
| 1-Methylnaphthalene | 0.10 | U | ug/L |
| 1-Naphthylamine | 10 | U | ug/L |
| 1,2,4-Trichlorobenzene | 0.050 | U | ug/L |
| 1,2,4,5-Tetrachlorobenzene | 0.050 | U | ug/L |
| 1,3-Dinitrobenzene | 0.20 | U | ug/L |
| 1,3,5-Trinitrobenzene | 0.20 | U | ug/L |
| 2-Acetylaminofluorene | 1.0 | U | ug/L |
| 2-Chloronaphthalene | 0.050 | U | ug/L |
| 2-Chlorophenol | 0.050 | U | ug/L |
| 2-Methyl-4,6-dinitrophenol | 3.0 | U | ug/L |
| 2-Methylnaphthalene | 0.10 | U | ug/L |
| 2-Naphthylamine | 10 | U | ug/L |
| 2-Nitroaniline | 0.050 | U | ug/L |
| 2-Nitrophenol | 0.050 | U | ug/L |
| 2-Picoline | 1.0 | U | ug/L |
| 2,3,4,6-Tetrachlorophenol | 0.10 | U | ug/L |
| 2,4-Dichlorophenol | 0.050 | U | ug/L |
| 2,4-Dimethylphenol | 0.050 | U | ug/L |
| 2,4-Dinitrophenol | 10 | U | ug/L |
| 2,4-Dinitrotoluene | 0.050 | U | ug/L |
| 2,4,5-Trichlorophenol | 0.050 | U | ug/L |
| 2,4,6-Trichlorophenol | 0.050 | U | ug/L |

# Quality Assurance Report
# Method Blank Results

**Reference Method: EPA 8270E**
**Batch ID: P418679**

| Component | Result | Code | Units |
|---|---|---|---|
| 2,6-Dichlorophenol | 0.050 | U | ug/L |
| 2,6-Dinitrotoluene | 0.050 | U | ug/L |
| 3-Methylcholanthrene | 0.10 | U | ug/L |
| 3,3'-Dichlorobenzidine | 10 | U | ug/L |
| 4-Aminobiphenyl | 4.0 | U | ug/L |
| 4-Bromophenyl phenyl ether | 0.050 | U | ug/L |
| 4-Chloro-3-methylphenol | 0.050 | U | ug/L |
| 4-Chlorophenyl phenyl ether | 0.050 | U | ug/L |
| 4-Nitrophenol | 10 | U | ug/L |
| 5-Nitro-o-toluidine | 0.10 | U | ug/L |
| 7,12-Dimethylbenz(a)anthracene | 0.10 | U | ug/L |
| Acenaphthene | 0.025 | U | ug/L |
| Acenaphthylene | 0.025 | U | ug/L |
| Acetophenone | 0.20 | U | ug/L |
| Aniline | 1.0 | U | ug/L |
| Anthracene | 0.050 | U | ug/L |
| Azobenzene/1,2-Diphenylhydrazine | 0.050 | U | ug/L |
| Benzidine | 10 | U | ug/L |
| Benzo(a)anthracene | 0.025 | U | ug/L |
| Benzo(a)pyrene | 0.025 | U | ug/L |
| Benzo(b)fluoranthene | 0.025 | U | ug/L |
| Benzo(g,h,i)perylene | 0.025 | U | ug/L |
| Benzo(k)fluoranthene | 0.025 | U | ug/L |
| Benzyl alcohol | 0.20 | U | ug/L |
| Bis(2-chloroethoxy)methane | 0.050 | U | ug/L |
| Bis(2-chloroethyl)ether | 0.050 | U | ug/L |
| Bis(2-chloroisopropyl)ether | 0.050 | U | ug/L |
| Bis(2-ethylhexyl)phthalate | 5.0 | U | ug/L |
| Butyl benzyl phthalate | 1.0 | U | ug/L |
| Carbazole | 0.050 | U | ug/L |
| Chrysene | 0.025 | U | ug/L |
| Di-n-butyl phthalate | 2.0 | U | ug/L |
| Di-n-octyl phthalate | 0.050 | U | ug/L |
| Dibenzo(a,h)anthracene | 0.025 | U | ug/L |
| Dibenzofuran | 0.050 | U | ug/L |
| Diethyl phthalate | 2.0 | U | ug/L |
| Dimethyl phthalate | 0.20 | U | ug/L |
| Dimethylaminoazobenzene | 0.050 | U | ug/L |
| Dinoseb | 4.0 | U | ug/L |
| Ethyl methanesulfonate | 1.0 | U | ug/L |
| Fluoranthene | 0.050 | U | ug/L |
| Fluorene | 0.025 | U | ug/L |
| Hexachlorobenzene | 0.050 | U | ug/L |
| Hexachlorobutadiene | 0.050 | U | ug/L |
| Hexachlorocyclopentadiene | 0.050 | U | ug/L |
| Hexachloroethane | 0.050 | U | ug/L |
| Hexachloropropene | 0.050 | U | ug/L |
| Indeno(1,2,3-cd)pyrene | 0.025 | U | ug/L |
| Isophorone | 0.050 | U | ug/L |
| Isosafrole | 0.050 | U | ug/L |
| m,p-Cresols | 0.050 | U | ug/L |
| N-Nitrosodi-n-butylamine | 0.050 | U | ug/L |

# Quality Assurance Report
# Method Blank Results

**Reference Method: EPA 8270E**
**Batch ID: P418679**

| Component | Result | Code | Units |
|---|---|---|---|
| N-Nitrosodi-n-propylamine | 0.050 | U | ug/L |
| N-Nitrosodiethylamine | 1.0 | U | ug/L |
| N-Nitrosodimethylamine | 2.0 | U | ug/L |
| N-Nitrosomethylethylamine | 2.0 | U | ug/L |
| N-Nitrosomorpholine | 0.050 | U | ug/L |
| N-Nitrosopiperidine | 0.050 | U | ug/L |
| N-Nitrosopyrrolidine | 0.10 | U | ug/L |
| Naphthalene | 0.10 | U | ug/L |
| Nitrobenzene | 0.050 | U | ug/L |
| o-Cresol | 0.050 | U | ug/L |
| o-Toluidine | 0.10 | U | ug/L |
| Pentachlorobenzene | 0.050 | U | ug/L |
| Pentachloroethane | 0.050 | U | ug/L |
| Pentachloronitrobenzene | 0.050 | U | ug/L |
| Pentachlorophenol | 0.50 | U | ug/L |
| Phenacetin | 1.0 | U | ug/L |
| Phenanthrene | 0.10 | U | ug/L |
| Phenol | 0.050 | U | ug/L |
| Pyrene | 0.10 | U | ug/L |
| Pyridine | 4.0 | U | ug/L |
| Safrole | 0.050 | U | ug/L |

**Reference Method: EPA 8321B**
**Batch ID: P418569**

| Component | Result | Code | Units |
|---|---|---|---|
| Cylindrospermopsin | 0.10 | U | ug/L |
| Desmethyl microcystin LR | 0.25 | U | ug/L |
| Microcystin HilR | 0.25 | U | ug/L |
| Microcystin HtyR | 0.25 | U | ug/L |
| Microcystin LA | 0.10 | U | ug/L |
| Microcystin LF | 0.10 | U | ug/L |
| Microcystin LR | 0.25 | U | ug/L |
| Microcystin LW | 0.25 | U | ug/L |
| Microcystin LY | 0.10 | U | ug/L |
| Microcystin RR | 0.10 | U | ug/L |
| Microcystin WR | 0.50 | U | ug/L |
| Microcystin YR | 0.25 | U | ug/L |
| Nodularin-R | 0.10 | U | ug/L |

**Reference Method: EPA 8321B**
**Batch ID: P418570**

| Component | Result | Code | Units |
|---|---|---|---|
| Anatoxin-a | 0.25 | U | ug/L |

**Reference Method: FL-PRO 2018**
**Batch ID: P418756**

| Component | Result | Code | Units |
|---|---|---|---|
| TRPH | 0.45 | U | mg/L |

# Quality Assurance Report
# Laboratory Control Sample Accuracy

**Reference Method: EPA 350.1 Rev. 2.0**
**Batch ID: P418886**

| Component | % Rec.1 | % Rec.2 | Pass/Fail | Control Limits |
|---|---|---|---|---|
| Ammonia-N | 96.8 | | P | 90 - 110 |

**Reference Method: EPA 351.2 Rev. 2.0**
**Batch ID: P418823**

| Component | % Rec.1 | % Rec.2 | Pass/Fail | Control Limits |
|---|---|---|---|---|
| Kjeldahl Nitrogen | 104 | | P | 90 - 110 |

**Reference Method: EPA 353.2 Rev. 2.0**
**Batch ID: P419050**

| Component | % Rec.1 | % Rec.2 | Pass/Fail | Control Limits |
|---|---|---|---|---|
| NO2NO3-N | 102 | | P | 90 - 110 |

**Reference Method: EPA 365.1 Rev. 2.0**
**Batch ID: P419131**

| Component | % Rec.1 | % Rec.2 | Pass/Fail | Control Limits |
|---|---|---|---|---|
| Total-P | 104 | | P | 90 - 110 |

**Reference Method: EPA 365.1 Rev. 2.0 dissolved**
**Batch ID: P418697**

| Component | % Rec.1 | % Rec.2 | Pass/Fail | Control Limits |
|---|---|---|---|---|
| O-Phosphate-P | 102 | | P | 90 - 110 |

**Reference Method: EPA 8270E**
**Batch ID: P418679**

| Component | % Rec.1 | % Rec.2 | Pass/Fail | Control Limits |
|---|---|---|---|---|
| 1-Methylnaphthalene | 92.7 | | P | 50 - 130 |
| 1-Naphthylamine | 65.2 | | P | 20 - 130 |
| 1,2,4-Trichlorobenzene | 84.1 | | P | 50 - 130 |
| 1,2,4,5-Tetrachlorobenzene | 96.0 | | P | 50 - 130 |
| 1,3-Dinitrobenzene | 103 | | P | 50 - 130 |
| 1,3,5-Trinitrobenzene | 124 | | P | 50 - 150 |
| 2-Acetylaminofluorene | 117 | | P | 50 - 130 |
| 2-Chloronaphthalene | 91.6 | | P | 50 - 130 |
| 2-Chlorophenol | 87.2 | | P | 50 - 130 |
| 2-Methyl-4,6-dinitrophenol | 99.3 | | P | 50 - 150 |
| 2-Methylnaphthalene | 92.3 | | P | 50 - 130 |
| 2-Naphthylamine | 66.1 | | P | 20 - 130 |
| 2-Nitroaniline | 104 | | P | 50 - 130 |
| 2-Nitrophenol | 92.5 | | P | 50 - 130 |
| 2-Picoline | 80.5 | | P | 40 - 130 |
| 2,3,4,6-Tetrachlorophenol | 107 | | P | 50 - 130 |
| 2,4-Dichlorophenol | 92.1 | | P | 50 - 130 |
| 2,4-Dimethylphenol | 88.3 | | P | 50 - 130 |
| 2,4-Dinitrophenol | 41.8 | | P | 30 - 160 |
| 2,4-Dinitrotoluene | 99.1 | | P | 50 - 130 |
| 2,4,5-Trichlorophenol | 87.8 | | P | 50 - 130 |
| 2,4,6-Trichlorophenol | 90.3 | | P | 50 - 130 |

# Quality Assurance Report
# Laboratory Control Sample Accuracy

**Reference Method: EPA 8270E**
**Batch ID: P418679**

| Component | % Rec.1 | % Rec.2 | Pass/Fail | Control Limits |
|---|---|---|---|---|
| 2,6-Dichlorophenol | 109 | | P | 50 - 130 |
| 2,6-Dinitrotoluene | 99.4 | | P | 50 - 130 |
| 3-Methylcholanthrene | 82.1 | | P | 50 - 130 |
| 3,3'-Dichlorobenzidine | 195 | | P | 0.0 - 200 |
| 4-Aminobiphenyl | 98.7 | | P | 30 - 130 |
| 4-Bromophenyl phenyl ether | 90.3 | | P | 50 - 130 |
| 4-Chloro-3-methylphenol | 91.1 | | P | 50 - 130 |
| 4-Chlorophenyl phenyl ether | 98.7 | | P | 50 - 130 |
| 4-Nitrophenol | 58.8 | | P | 15 - 110 |
| 5-Nitro-o-toluidine | 120 | | P | 50 - 130 |
| 7,12-Dimethylbenz(a)anthracene | 106 | | P | 50 - 130 |
| Acenaphthene | 96.8 | | P | 50 - 130 |
| Acenaphthylene | 95.8 | | P | 50 - 130 |
| Acetophenone | 114 | | P | 50 - 130 |
| Aniline | 90.5 | | P | 30 - 130 |
| Anthracene | 94.3 | | P | 50 - 130 |
| Azobenzene/1,2-Diphenylhydrazine | 99.1 | | P | 50 - 130 |
| Benzidine | 65.0 | | P | 0.0 - 240 |
| Benzo(a)anthracene | 97.8 | | P | 50 - 130 |
| Benzo(a)pyrene | 92.0 | | P | 50 - 130 |
| Benzo(b)fluoranthene | 97.3 | | P | 50 - 130 |
| Benzo(g,h,i)perylene | 74.9 | | P | 50 - 130 |
| Benzo(k)fluoranthene | 91.7 | | P | 50 - 130 |
| Benzyl alcohol | 91.6 | | P | 50 - 130 |
| Bis(2-chloroethoxy)methane | 91.9 | | P | 50 - 130 |
| Bis(2-chloroethyl)ether | 91.6 | | P | 50 - 160 |
| Bis(2-chloroisopropyl)ether | 90.1 | | P | 50 - 130 |
| Bis(2-ethylhexyl)phthalate | 116 | | P | 50 - 160 |
| Butyl benzyl phthalate | 109 | | P | 50 - 160 |
| Carbazole | 118 | | P | 50 - 130 |
| Chrysene | 98.2 | | P | 50 - 130 |
| Di-n-butyl phthalate | 102 | | P | 50 - 160 |
| Di-n-octyl phthalate | 107 | | P | 50 - 130 |
| Dibenzo(a,h)anthracene | 80.9 | | P | 50 - 130 |
| Dibenzofuran | 96.8 | | P | 50 - 130 |
| Diethyl phthalate | 110 | | P | 50 - 130 |
| Dimethyl phthalate | 104 | | P | 50 - 130 |
| Dimethylaminoazobenzene | 113 | | P | 50 - 130 |
| Dinoseb | 126 | | P | 50 - 150 |
| Ethyl methanesulfonate | 96.2 | | P | 50 - 130 |
| Fluoranthene | 94.5 | | P | 50 - 130 |
| Fluorene | 100 | | P | 50 - 130 |
| Hexachlorobenzene | 90.9 | | P | 50 - 130 |
| Hexachlorobutadiene | 84.5 | | P | 20 - 130 |
| Hexachlorocyclopentadiene | 51.9 | | P | 20 - 130 |
| Hexachloroethane | 84.0 | | P | 40 - 130 |
| Hexachloropropene | 97.0 | | P | 50 - 130 |
| Indeno(1,2,3-cd)pyrene | 78.8 | | P | 50 - 130 |
| Isophorone | 96.2 | | P | 50 - 130 |
| Isosafrole | 108 | | P | 50 - 130 |
| m,p-Cresols | 82.5 | | P | 50 - 130 |
| N-Nitrosodi-n-butylamine | 107 | | P | 50 - 130 |

# Quality Assurance Report
# Laboratory Control Sample Accuracy

**Reference Method: EPA 8270E**
**Batch ID: P418679**

| Component | % Rec.1 | % Rec.2 | Pass/Fail | Control Limits |
|---|---|---|---|---|
| N-Nitrosodi-n-propylamine | 89.1 | | P | 50 - 130 |
| N-Nitrosodiethylamine | 104 | | P | 50 - 130 |
| N-Nitrosodimethylamine | 92.4 | | P | 30 - 130 |
| N-Nitrosomethylethylamine | 104 | | P | 50 - 130 |
| N-Nitrosomorpholine | 127 | | P | 50 - 150 |
| N-Nitrosopiperidine | 109 | | P | 50 - 130 |
| N-Nitrosopyrrolidine | 96.1 | | P | 50 - 130 |
| Naphthalene | 90.7 | | P | 50 - 130 |
| Nitrobenzene | 95.9 | | P | 50 - 130 |
| o-Cresol | 86.7 | | P | 50 - 130 |
| o-Toluidine | 111 | | P | 50 - 130 |
| Pentachlorobenzene | 99.7 | | P | 50 - 130 |
| Pentachloroethane | 95.8 | | P | 50 - 130 |
| Pentachloronitrobenzene | 130 | | P | 50 - 130 |
| Pentachlorophenol | 91.5 | | P | 50 - 130 |
| Phenacetin | 130 | | P | 50 - 130 |
| Phenanthrene | 95.2 | | P | 50 - 130 |
| Phenol | 54.8 | | P | 15 - 110 |
| Pyrene | 98.3 | | P | 50 - 130 |
| Pyridine | 70.4 | | P | 20 - 130 |
| Safrole | 106 | | P | 50 - 130 |

**Reference Method: EPA 8321B**
**Batch ID: P418569**

| Component | % Rec.1 | % Rec.2 | Pass/Fail | Control Limits |
|---|---|---|---|---|
| Cylindrospermopsin | 79.8 | | P | 40 - 150 |
| Desmethyl microcystin LR | 73.2 | | P | 40 - 150 |
| Microcystin HilR | 77.2 | | P | 40 - 150 |
| Microcystin HtyR | 78.1 | | P | 40 - 150 |
| Microcystin LA | 80.6 | | P | 40 - 150 |
| Microcystin LF | 66.4 | | P | 40 - 150 |
| Microcystin LR | 76.4 | | P | 40 - 150 |
| Microcystin LW | 64.4 | | P | 40 - 150 |
| Microcystin LY | 66.1 | | P | 40 - 150 |
| Microcystin RR | 77.1 | | P | 40 - 150 |
| Microcystin WR | 73.7 | | P | 40 - 150 |
| Microcystin YR | 80.0 | | P | 40 - 150 |
| Nodularin-R | 80.8 | | P | 40 - 150 |

**Reference Method: EPA 8321B**
**Batch ID: P418570**

| Component | % Rec.1 | % Rec.2 | Pass/Fail | Control Limits |
|---|---|---|---|---|
| Anatoxin-a | 75.5 | | P | 40 - 150 |

**Reference Method: FL-PRO 2018**
**Batch ID: P418756**

| Component | % Rec.1 | % Rec.2 | Pass/Fail | Control Limits |
|---|---|---|---|---|
| TRPH | 90.9 | | P | 66 - 116 |

# Quality Assurance Report
# Matrix Spike Accuracy

**Reference Method: EPA 350.1 Rev. 2.0**
**Batch ID: P418886**

| Spiked Sample | Component | % Rec.1 | % Rec.2 | Pass/Fail | Control Limits |
|---|---|---|---|---|---|
| 2352296 | Ammonia-N | 95.2 | 95.4 | P/P | 90 - 110 |

**Reference Method: EPA 351.2 Rev. 2.0**
**Batch ID: P418823**

| Spiked Sample | Component | % Rec.1 | % Rec.2 | Pass/Fail | Control Limits |
|---|---|---|---|---|---|
| 2352296 | Kjeldahl Nitrogen | 100 | 96.9 | P/P | 90 - 110 |

**Reference Method: EPA 353.2 Rev. 2.0**
**Batch ID: P419050**

| Spiked Sample | Component | % Rec.1 | % Rec.2 | Pass/Fail | Control Limits |
|---|---|---|---|---|---|
| 2352331 | NO2NO3-N | 99.5 | 102 | P/P | 90 - 110 |

**Reference Method: EPA 365.1 Rev. 2.0**
**Batch ID: P419131**

| Spiked Sample | Component | % Rec.1 | % Rec.2 | Pass/Fail | Control Limits |
|---|---|---|---|---|---|
| 2352281 | Total-P | 104 | 106 | P/P | 90 - 110 |

**Reference Method: EPA 365.1 Rev. 2.0 dissolved**
**Batch ID: P418697**

| Spiked Sample | Component | % Rec.1 | % Rec.2 | Pass/Fail | Control Limits |
|---|---|---|---|---|---|
| 2352277 | O-Phosphate-P | 101 | 100 | P/P | 90 - 110 |

**Reference Method: EPA 8270E**
**Batch ID: P418679**

| Spiked Sample | Component | % Rec.1 | % Rec.2 | Pass/Fail | Control Limits |
|---|---|---|---|---|---|
| 2352292 | 1-Methylnaphthalene | 93.7 | 89.6 | P/P | 50 - 130 |
| 2352292 | 1-Naphthylamine | 106 | 87.5 | P/P | 20 - 130 |
| 2352292 | 1,2,4-Trichlorobenzene | 84.9 | 78.9 | P/P | 50 - 130 |
| 2352292 | 1,2,4,5-Tetrachlorobenzene | 96.4 | 89.1 | P/P | 50 - 130 |
| 2352292 | 1,3-Dinitrobenzene | 97.2 | 101 | P/P | 50 - 130 |
| 2352292 | 1,3,5-Trinitrobenzene | 112 | 117 | P/P | 50 - 150 |
| 2352292 | 2-Acetylaminofluorene | 115 | 121 | P/P | 50 - 130 |
| 2352292 | 2-Chloronaphthalene | 91.8 | 87.4 | P/P | 50 - 130 |
| 2352292 | 2-Chlorophenol | 86.9 | 81.1 | P/P | 50 - 130 |
| 2352292 | 2-Methyl-4,6-dinitrophenol | 115 | 116 | P/P | 50 - 150 |
| 2352292 | 2-Methylnaphthalene | 92.3 | 87.5 | P/P | 50 - 130 |
| 2352292 | 2-Naphthylamine | 75.7 | 71.5 | P/P | 20 - 130 |
| 2352292 | 2-Nitroaniline | 103 | 98.0 | P/P | 50 - 130 |
| 2352292 | 2-Nitrophenol | 90.8 | 87.7 | P/P | 50 - 130 |
| 2352292 | 2-Picoline | 99.4 | 92.6 | P/P | 40 - 130 |
| 2352292 | 2,3,4,6-Tetrachlorophenol | 104 | 107 | P/P | 50 - 130 |
| 2352292 | 2,4-Dichlorophenol | 92.8 | 88.7 | P/P | 50 - 130 |
| 2352292 | 2,4-Dimethylphenol | 91.5 | 88.0 | P/P | 50 - 130 |
| 2352292 | 2,4-Dinitrophenol | 74.4 | 72.6 | P/P | 30 - 160 |
| 2352292 | 2,4-Dinitrotoluene | 99.8 | 96.2 | P/P | 50 - 130 |
| 2352292 | 2,4,5-Trichlorophenol | 90.0 | 88.9 | P/P | 50 - 130 |
| 2352292 | 2,4,6-Trichlorophenol | 90.6 | 88.3 | P/P | 50 - 130 |

# Quality Assurance Report
# Matrix Spike Accuracy

**Reference Method: EPA 8270E**
**Batch ID: P418679**

| Spiked Sample | Component | % Rec.1 | % Rec.2 | Pass/Fail | Control Limits |
|---|---|---|---|---|---|
| 2352292 | 2,6-Dichlorophenol | 109 | 103 | P/P | 50 - 130 |
| 2352292 | 2,6-Dinitrotoluene | 100 | 95.8 | P/P | 50 - 130 |
| 2352292 | 3-Methylcholanthrene | 92.4 | 97.2 | P/P | 50 - 130 |
| 2352292 | 3,3'-Dichlorobenzidine | 180 | 169 | P/P | 0.0 - 200 |
| 2352292 | 4-Aminobiphenyl | 103 | 98.1 | P/P | 30 - 130 |
| 2352292 | 4-Bromophenyl phenyl ether | 88.2 | 88.1 | P/P | 50 - 130 |
| 2352292 | 4-Chloro-3-methylphenol | 92.1 | 89.4 | P/P | 50 - 130 |
| 2352292 | 4-Chlorophenyl phenyl ether | 100 | 94.6 | P/P | 50 - 130 |
| 2352292 | 4-Nitrophenol | 53.7 | 53.0 | P/P | 15 - 110 |
| 2352292 | 5-Nitro-o-toluidine | 114 | 121 | P/P | 50 - 130 |
| 2352292 | 7,12-Dimethylbenz(a)anthracene | 85.6 | 97.4 | P/P | 50 - 130 |
| 2352292 | Acenaphthene | 97.8 | 93.4 | P/P | 50 - 130 |
| 2352292 | Acenaphthylene | 96.1 | 90.8 | P/P | 50 - 130 |
| 2352292 | Acetophenone | 114 | 104 | P/P | 50 - 130 |
| 2352292 | Aniline | 89.7 | 85.1 | P/P | 30 - 130 |
| 2352292 | Anthracene | 91.5 | 91.3 | P/P | 50 - 130 |
| 2352292 | Azobenzene/1,2-Diphenylhydrazine | 98.7 | 100 | P/P | 50 - 130 |
| 2352292 | Benzidine | 23.9 | 15.6 | P/P | 0.0 - 240 |
| 2352292 | Benzo(a)anthracene | 93.4 | 92.7 | P/P | 50 - 130 |
| 2352292 | Benzo(a)pyrene | 96.0 | 88.0 | P/P | 50 - 130 |
| 2352292 | Benzo(b)fluoranthene | 99.2 | 90.7 | P/P | 50 - 130 |
| 2352292 | Benzo(g,h,i)perylene | 81.8 | 79.0 | P/P | 50 - 130 |
| 2352292 | Benzo(k)fluoranthene | 94.5 | 87.7 | P/P | 50 - 130 |
| 2352292 | Benzyl alcohol | 95.1 | 83.2 | P/P | 50 - 130 |
| 2352292 | Bis(2-chloroethoxy)methane | 91.7 | 88.9 | P/P | 50 - 130 |
| 2352292 | Bis(2-chloroethyl)ether | 89.7 | 82.8 | P/P | 50 - 160 |
| 2352292 | Bis(2-chloroisopropyl)ether | 89.3 | 83.8 | P/P | 50 - 130 |
| 2352292 | Bis(2-ethylhexyl)phthalate | 115 | 114 | P/P | 50 - 160 |
| 2352292 | Butyl benzyl phthalate | 104 | 105 | P/P | 50 - 160 |
| 2352292 | Carbazole | 126 | 128 | P/P | 50 - 130 |
| 2352292 | Chrysene | 91.6 | 92.1 | P/P | 50 - 130 |
| 2352292 | Di-n-butyl phthalate | 97.5 | 98.0 | P/P | 50 - 160 |
| 2352292 | Di-n-octyl phthalate | 111 | 106 | P/P | 50 - 130 |
| 2352292 | Dibenzo(a,h)anthracene | 87.0 | 84.4 | P/P | 50 - 130 |
| 2352292 | Dibenzofuran | 97.6 | 92.8 | P/P | 50 - 130 |
| 2352292 | Diethyl phthalate | 107 | 104 | P/P | 50 - 130 |
| 2352292 | Dimethyl phthalate | 102 | 98.7 | P/P | 50 - 130 |
| 2352292 | Dimethylaminoazobenzene | 93.0 | 103 | P/P | 50 - 130 |
| 2352292 | Dinoseb | 128 | 128 | P/P | 50 - 150 |
| 2352292 | Ethyl methanesulfonate | 97.0 | 89.9 | P/P | 50 - 130 |
| 2352292 | Fluoranthene | 90.6 | 91.1 | P/P | 50 - 130 |
| 2352292 | Fluorene | 99.9 | 96.0 | P/P | 50 - 130 |
| 2352292 | Hexachlorobenzene | 87.9 | 89.1 | P/P | 50 - 130 |
| 2352292 | Hexachlorobutadiene | 85.1 | 80.6 | P/P | 20 - 130 |
| 2352292 | Hexachlorocyclopentadiene | 49.8 | 43.7 | P/P | 20 - 130 |
| 2352292 | Hexachloroethane | 80.7 | 72.8 | P/P | 40 - 130 |
| 2352292 | Hexachloropropene | 89.5 | 76.3 | P/P | 50 - 130 |
| 2352292 | Indeno(1,2,3-cd)pyrene | 85.3 | 81.7 | P/P | 50 - 130 |
| 2352292 | Isophorone | 96.1 | 92.3 | P/P | 50 - 130 |
| 2352292 | Isosafrole | 101 | 108 | P/P | 50 - 130 |
| 2352292 | m,p-Cresols | 81.4 | 78.5 | P/P | 50 - 130 |
| 2352292 | N-Nitrosodi-n-butylamine | 109 | 101 | P/P | 50 - 130 |

# Quality Assurance Report
# Matrix Spike Accuracy

**Reference Method: EPA 8270E**
**Batch ID: P418679**

| Spiked Sample | Component | % Rec.1 | % Rec.2 | Pass/Fail | Control Limits |
|---|---|---|---|---|---|
| 2352292 | N-Nitrosodi-n-propylamine | 90.8 | 85.1 | P/P | 50 - 130 |
| 2352292 | N-Nitrosodiethylamine | 103 | 95.4 | P/P | 50 - 130 |
| 2352292 | N-Nitrosodimethylamine | 104 | 97.7 | P/P | 30 - 130 |
| 2352292 | N-Nitrosomethylethylamine | 103 | 95.0 | P/P | 50 - 130 |
| 2352292 | N-Nitrosomorpholine | 126 | 117 | P/P | 50 - 150 |
| 2352292 | N-Nitrosopiperidine | 109 | 101 | P/P | 50 - 130 |
| 2352292 | N-Nitrosopyrrolidine | 94.8 | 88.9 | P/P | 50 - 130 |
| 2352292 | Naphthalene | 91.0 | 85.6 | P/P | 50 - 130 |
| 2352292 | Nitrobenzene | 95.1 | 91.7 | P/P | 50 - 130 |
| 2352292 | o-Cresol | 85.7 | 82.2 | P/P | 50 - 130 |
| 2352292 | o-Toluidine | 110 | 102 | P/P | 50 - 130 |
| 2352292 | Pentachlorobenzene | 92.3 | 102 | P/P | 50 - 130 |
| 2352292 | Pentachloroethane | 93.9 | 85.1 | P/P | 50 - 130 |
| 2352292 | Pentachloronitrobenzene | 124 | 127 | P/P | 50 - 130 |
| 2352292 | Pentachlorophenol | 96.7 | 98.4 | P/P | 50 - 130 |
| 2352292 | Phenacetin | 122 | 125 | P/P | 50 - 130 |
| 2352292 | Phenanthrene | 92.5 | 94.1 | P/P | 50 - 130 |
| 2352292 | Phenol | 54.8 | 51.6 | P/P | 15 - 110 |
| 2352292 | Pyrene | 94.2 | 93.8 | P/P | 50 - 130 |
| 2352292 | Pyridine | 77.5 | 73.5 | P/P | 20 - 130 |
| 2352292 | Safrole | 108 | 100 | P/P | 50 - 130 |

**Reference Method: EPA 8321B**
**Batch ID: P418569**

| Spiked Sample | Component | % Rec.1 | % Rec.2 | Pass/Fail | Control Limits |
|---|---|---|---|---|---|
| 2351625 | Cylindrospermopsin | 91.4 | 88.2 | P/P | 40 - 150 |
| 2351625 | Desmethyl microcystin LR | 106 | 112 | P/P | 40 - 150 |
| 2351625 | Microcystin HilR | 129 | 136 | P/P | 40 - 150 |
| 2351625 | Microcystin HtyR | 108 | 115 | P/P | 40 - 150 |
| 2351625 | Microcystin LA | 98.8 | 94.2 | P/P | 40 - 150 |
| 2351625 | Microcystin LF | 77.2 | 82.0 | P/P | 40 - 150 |
| 2351625 | Microcystin LR | 110 | 115 | P/P | 40 - 150 |
| 2351625 | Microcystin LW | 94.1 | 98.5 | P/P | 40 - 150 |
| 2351625 | Microcystin LY | 94.5 | 84.8 | P/P | 40 - 150 |
| 2351625 | Microcystin RR | 102 | 107 | P/P | 40 - 150 |
| 2351625 | Microcystin WR | 107 | 115 | P/P | 40 - 150 |
| 2351625 | Microcystin YR | 110 | 116 | P/P | 40 - 150 |
| 2351625 | Nodularin-R | 115 | 118 | P/P | 40 - 150 |

**Reference Method: EPA 8321B**
**Batch ID: P418570**

| Spiked Sample | Component | % Rec.1 | % Rec.2 | Pass/Fail | Control Limits |
|---|---|---|---|---|---|
| 2351625 | Anatoxin-a | 82.8 | 81.5 | P/P | 40 - 150 |

**Reference Method: FL-PRO 2018**
**Batch ID: P418756**

| Spiked Sample | Component | % Rec.1 | % Rec.2 | Pass/Fail | Control Limits |
|---|---|---|---|---|---|
| 2352293 | TRPH | 64.3 | 83.6 | F/P | 65 - 123 |

# Quality Assurance Report
# Precision

**Reference Method: EPA 350.1 Rev. 2.0**
**Batch ID: P418886**

| Replicated Lab Sample | Component | % RSD/RPD | Sample/Spike/LCS* | Pass/Fail | Control Limits |
|---|---|---|---|---|---|
| 2352296 | Ammonia-N | 0.195 | Spike | P | 0 - 20 |

**Reference Method: EPA 351.2 Rev. 2.0**
**Batch ID: P418823**

| Replicated Lab Sample | Component | % RSD/RPD | Sample/Spike/LCS* | Pass/Fail | Control Limits |
|---|---|---|---|---|---|
| 2352296 | Kjeldahl Nitrogen | 1.70 | Spike | P | 0 - 20 |

**Reference Method: EPA 353.2 Rev. 2.0**
**Batch ID: P419050**

| Replicated Lab Sample | Component | % RSD/RPD | Sample/Spike/LCS* | Pass/Fail | Control Limits |
|---|---|---|---|---|---|
| 2352331 | NO2NO3-N | 1.19 | Spike | P | 0 - 20 |

**Reference Method: EPA 365.1 Rev. 2.0**
**Batch ID: P419131**

| Replicated Lab Sample | Component | % RSD/RPD | Sample/Spike/LCS* | Pass/Fail | Control Limits |
|---|---|---|---|---|---|
| 2352281 | Total-P | 1.83 | Spike | P | 0 - 20 |

**Reference Method: EPA 365.1 Rev. 2.0 dissolved**
**Batch ID: P418697**

| Replicated Lab Sample | Component | % RSD/RPD | Sample/Spike/LCS* | Pass/Fail | Control Limits |
|---|---|---|---|---|---|
| 2352277 | O-Phosphate-P | 0.661 | Spike | P | 0 - 20 |

**Reference Method: EPA 8270E**
**Batch ID: P418679**

| Replicated Lab Sample | Component | % RSD/RPD | Sample/Spike/LCS* | Pass/Fail | Control Limits |
|---|---|---|---|---|---|
| 2352292 | 1-Methylnaphthalene | 4.47 | Spike | P | 0 - 40 |
| 2352292 | 1-Naphthylamine | 19.4 | Spike | P | 0 - 40 |
| 2352292 | 1,2,4-Trichlorobenzene | 7.33 | Spike | P | 0 - 40 |
| 2352292 | 1,2,4,5-Tetrachlorobenzene | 7.87 | Spike | P | 0 - 40 |
| 2352292 | 1,3-Dinitrobenzene | 3.54 | Spike | P | 0 - 40 |
| 2352292 | 1,3,5-Trinitrobenzene | 4.20 | Spike | P | 0 - 40 |
| 2352292 | 2-Acetylaminofluorene | 4.67 | Spike | P | 0 - 40 |
| 2352292 | 2-Chloronaphthalene | 4.91 | Spike | P | 0 - 40 |
| 2352292 | 2-Chlorophenol | 6.90 | Spike | P | 0 - 40 |
| 2352292 | 2-Methyl-4,6-dinitrophenol | 0.694 | Spike | P | 0 - 40 |
| 2352292 | 2-Methylnaphthalene | 5.34 | Spike | P | 0 - 40 |
| 2352292 | 2-Naphthylamine | 5.71 | Spike | P | 0 - 40 |
| 2352292 | 2-Nitroaniline | 4.68 | Spike | P | 0 - 40 |
| 2352292 | 2-Nitrophenol | 3.47 | Spike | P | 0 - 40 |
| 2352292 | 2-Picoline | 7.08 | Spike | P | 0 - 40 |
| 2352292 | 2,3,4,6-Tetrachlorophenol | 3.37 | Spike | P | 0 - 40 |

# Quality Assurance Report
# Precision

**Reference Method: EPA 8270E**
**Batch ID: P418679**

| Replicated Lab Sample | Component | % RSD/RPD | Sample/Spike/LCS* | Pass/Fail | Control Limits |
|---|---|---|---|---|---|
| 2352292 | 2,4-Dichlorophenol | 4.52 | Spike | P | 0 - 40 |
| 2352292 | 2,4-Dimethylphenol | 3.90 | Spike | P | 0 - 40 |
| 2352292 | 2,4-Dinitrophenol | 2.45 | Spike | P | 0 - 40 |
| 2352292 | 2,4-Dinitrotoluene | 3.67 | Spike | P | 0 - 40 |
| 2352292 | 2,4,5-Trichlorophenol | 1.23 | Spike | P | 0 - 40 |
| 2352292 | 2,4,6-Trichlorophenol | 2.57 | Spike | P | 0 - 40 |
| 2352292 | 2,6-Dichlorophenol | 5.66 | Spike | P | 0 - 40 |
| 2352292 | 2,6-Dinitrotoluene | 4.59 | Spike | P | 0 - 40 |
| 2352292 | 3-Methylcholanthrene | 5.06 | Spike | P | 0 - 40 |
| 2352292 | 3,3'-Dichlorobenzidine | 5.85 | Spike | P | 0 - 40 |
| 2352292 | 4-Aminobiphenyl | 4.97 | Spike | P | 0 - 40 |
| 2352292 | 4-Bromophenyl phenyl ether | 0.113 | Spike | P | 0 - 40 |
| 2352292 | 4-Chloro-3-methylphenol | 2.98 | Spike | P | 0 - 40 |
| 2352292 | 4-Chlorophenyl phenyl ether | 5.75 | Spike | P | 0 - 40 |
| 2352292 | 4-Nitrophenol | 1.31 | Spike | P | 0 - 40 |
| 2352292 | 5-Nitro-o-toluidine | 6.12 | Spike | P | 0 - 40 |
| 2352292 | 7,12-Dimethylbenz(a)anthracene | 12.9 | Spike | P | 0 - 40 |
| 2352292 | Acenaphthene | 4.60 | Spike | P | 0 - 40 |
| 2352292 | Acenaphthylene | 5.67 | Spike | P | 0 - 40 |
| 2352292 | Acetophenone | 9.21 | Spike | P | 0 - 40 |
| 2352292 | Aniline | 5.26 | Spike | P | 0 - 40 |
| 2352292 | Anthracene | 0.219 | Spike | P | 0 - 40 |
| 2352292 | Azobenzene/1,2-Diphenylhydrazine | 1.51 | Spike | P | 0 - 40 |
| 2352292 | Benzo(a)anthracene | 0.752 | Spike | P | 0 - 40 |
| 2352292 | Benzo(a)pyrene | 8.70 | Spike | P | 0 - 40 |
| 2352292 | Benzo(b)fluoranthene | 8.95 | Spike | P | 0 - 40 |
| 2352292 | Benzo(g,h,i)perylene | 3.48 | Spike | P | 0 - 40 |
| 2352292 | Benzo(k)fluoranthene | 7.46 | Spike | P | 0 - 40 |
| 2352292 | Benzyl alcohol | 13.3 | Spike | P | 0 - 40 |
| 2352292 | Bis(2-chloroethoxy)methane | 3.10 | Spike | P | 0 - 40 |
| 2352292 | Bis(2-chloroethyl)ether | 8.00 | Spike | P | 0 - 40 |
| 2352292 | Bis(2-chloroisopropyl)ether | 6.35 | Spike | P | 0 - 40 |
| 2352292 | Bis(2-ethylhexyl)phthalate | 1.05 | Spike | P | 0 - 40 |
| 2352292 | Butyl benzyl phthalate | 0.478 | Spike | P | 0 - 40 |
| 2352292 | Carbazole | 1.81 | Spike | P | 0 - 40 |
| 2352292 | Chrysene | 0.544 | Spike | P | 0 - 40 |
| 2352292 | Di-n-butyl phthalate | 0.512 | Spike | P | 0 - 40 |
| 2352292 | Di-n-octyl phthalate | 4.98 | Spike | P | 0 - 40 |
| 2352292 | Dibenzo(a,h)anthracene | 3.03 | Spike | P | 0 - 40 |
| 2352292 | Dibenzofuran | 5.04 | Spike | P | 0 - 40 |
| 2352292 | Diethyl phthalate | 2.75 | Spike | P | 0 - 40 |
| 2352292 | Dimethyl phthalate | 2.90 | Spike | P | 0 - 40 |
| 2352292 | Dimethylaminoazobenzene | 10.0 | Spike | P | 0 - 40 |
| 2352292 | Dinoseb | 0.391 | Spike | P | 0 - 40 |
| 2352292 | Ethyl methanesulfonate | 7.60 | Spike | P | 0 - 40 |
| 2352292 | Fluoranthene | 0.550 | Spike | P | 0 - 40 |
| 2352292 | Fluorene | 3.98 | Spike | P | 0 - 40 |
| 2352292 | Hexachlorobenzene | 1.36 | Spike | P | 0 - 40 |
| 2352292 | Hexachlorobutadiene | 5.43 | Spike | P | 0 - 40 |
| 2352292 | Hexachlorocyclopentadiene | 13.0 | Spike | P | 0 - 40 |
| 2352292 | Hexachloroethane | 10.3 | Spike | P | 0 - 40 |

# Quality Assurance Report
# Precision

**Reference Method: EPA 8270E**
**Batch ID: P418679**

| Replicated Lab Sample | Component | % RSD/RPD | Sample/Spike/LCS* | Pass/Fail | Control Limits |
|---|---|---|---|---|---|
| 2352292 | Hexachloropropene | 15.9 | Spike | P | 0 - 40 |
| 2352292 | Indeno(1,2,3-cd)pyrene | 4.31 | Spike | P | 0 - 40 |
| 2352292 | Isophorone | 4.03 | Spike | P | 0 - 40 |
| 2352292 | Isosafrole | 6.82 | Spike | P | 0 - 40 |
| 2352292 | m,p-Cresols | 3.63 | Spike | P | 0 - 40 |
| 2352292 | N-Nitrosodi-n-butylamine | 7.46 | Spike | P | 0 - 40 |
| 2352292 | N-Nitrosodi-n-propylamine | 6.48 | Spike | P | 0 - 40 |
| 2352292 | N-Nitrosodiethylamine | 7.37 | Spike | P | 0 - 40 |
| 2352292 | N-Nitrosodimethylamine | 6.05 | Spike | P | 0 - 40 |
| 2352292 | N-Nitrosomethylethylamine | 8.27 | Spike | P | 0 - 40 |
| 2352292 | N-Nitrosomorpholine | 7.50 | Spike | P | 0 - 40 |
| 2352292 | N-Nitrosopiperidine | 7.51 | Spike | P | 0 - 40 |
| 2352292 | N-Nitrosopyrrolidine | 6.42 | Spike | P | 0 - 40 |
| 2352292 | Naphthalene | 6.12 | Spike | P | 0 - 40 |
| 2352292 | Nitrobenzene | 3.64 | Spike | P | 0 - 40 |
| 2352292 | o-Cresol | 4.17 | Spike | P | 0 - 40 |
| 2352292 | o-Toluidine | 7.74 | Spike | P | 0 - 40 |
| 2352292 | Pentachlorobenzene | 9.69 | Spike | P | 0 - 40 |
| 2352292 | Pentachloroethane | 9.83 | Spike | P | 0 - 40 |
| 2352292 | Pentachloronitrobenzene | 2.15 | Spike | P | 0 - 40 |
| 2352292 | Pentachlorophenol | 1.74 | Spike | P | 0 - 40 |
| 2352292 | Phenacetin | 2.18 | Spike | P | 0 - 40 |
| 2352292 | Phenanthrene | 1.71 | Spike | P | 0 - 40 |
| 2352292 | Phenol | 6.02 | Spike | P | 0 - 40 |
| 2352292 | Pyrene | 0.426 | Spike | P | 0 - 40 |
| 2352292 | Pyridine | 5.30 | Spike | P | 0 - 40 |
| 2352292 | Safrole | 7.11 | Spike | P | 0 - 40 |

**Reference Method: EPA 8321B**
**Batch ID: P418569**

| Replicated Lab Sample | Component | % RSD/RPD | Sample/Spike/LCS* | Pass/Fail | Control Limits |
|---|---|---|---|---|---|
| 2351625 | Cylindrospermopsin | 3.49 | Spike | P | 0 - 30 |
| 2351625 | Desmethyl microcystin LR | 5.15 | Spike | P | 0 - 30 |
| 2351625 | Microcystin HilR | 5.47 | Spike | P | 0 - 30 |
| 2351625 | Microcystin HtyR | 5.96 | Spike | P | 0 - 30 |
| 2351625 | Microcystin LA | 4.76 | Spike | P | 0 - 30 |
| 2351625 | Microcystin LF | 6.09 | Spike | P | 0 - 30 |
| 2351625 | Microcystin LR | 3.94 | Spike | P | 0 - 30 |
| 2351625 | Microcystin LW | 4.58 | Spike | P | 0 - 30 |
| 2351625 | Microcystin LY | 10.8 | Spike | P | 0 - 30 |
| 2351625 | Microcystin RR | 4.42 | Spike | P | 0 - 30 |
| 2351625 | Microcystin WR | 7.17 | Spike | P | 0 - 30 |
| 2351625 | Microcystin YR | 4.96 | Spike | P | 0 - 30 |
| 2351625 | Nodularin-R | 3.05 | Spike | P | 0 - 30 |

# Quality Assurance Report
# Precision

**Reference Method: EPA 8321B**
**Batch ID: P418570**

| Replicated Lab Sample | Component | % RSD/RPD | Sample/Spike/LCS* | Pass/Fail | Control Limits |
|---|---|---|---|---|---|
| 2351625 | Anatoxin-a | 1.55 | Spike | P | 0 - 30 |

**Reference Method: FL-PRO 2018**
**Batch ID: P418756**

| Replicated Lab Sample | Component | % RSD/RPD | Sample/Spike/LCS* | Pass/Fail | Control Limits |
|---|---|---|---|---|---|
| 2352293 | TRPH | 22.1 | Spike | F | 0 - 20 |

* Sample, spike and/or laboratory control sample precision (LCS) is reported.

# Quality Assurance Report
# Surrogates

**Lab Sample ID: 2352292**
**Field Sample ID: HAB-SD-082422-1505**

| Reference Method | Surrogate | % Rec. | Pass/Fail | Control Limits |
|---|---|---|---|---|
| EPA 8270E | 2-Fluorobiphenyl | 107 | P | 30 - 150 |
| EPA 8270E | 2-Fluorophenol | 65.2 | P | 20 - 150 |
| EPA 8270E | 2,4,6-Tribromophenol | 95.1 | P | 30 - 150 |
| EPA 8270E | Nitrobenzene-d5 | 120 | P | 30 - 150 |
| EPA 8270E | Phenol-d5 | 58.6 | P | 20 - 150 |
| EPA 8270E | Terphenyl-d14 | 127 | P | 30 - 150 |

**Lab Sample ID: 2352293**
**Field Sample ID: HAB-SD-082422-1505**

| Reference Method | Surrogate | % Rec. | Pass/Fail | Control Limits |
|---|---|---|---|---|
| FL-PRO 2018 | Nonatriacontane(C39) | 70.3 | P | 40 - 120 |
| FL-PRO 2018 | o-Terphenyl | 93.5 | P | 66 - 137 |

**Lab Sample ID: 2352294**
**Field Sample ID: HAB-SD-082422-1505**

| Reference Method | Surrogate | % Rec. | Pass/Fail | Control Limits |
|---|---|---|---|---|
| EPA 8321B | Microcystin LR ethyl-d5 | 104 | P | 30 - 160 |
| EPA 8321B | Neosaxitoxin-15N7 | 75.5 | P | 30 - 160 |

# Quality Assurance Report
# Calibration Verification

**Reference Method: EPA 8321B**
**Run ID: A114331**
**Included Lab Sample IDs: 2352294**

| Component | % Rec.1 | % Rec.2 | Pass/Fail* | Control Limits |
|---|---|---|---|---|
| Anatoxin-a | 88.9 | 86.3 | P/P | 50 - 160 |

**Reference Method: EPA 365.1 Rev. 2.0 dissolved**
**Run ID: A114347**
**Included Lab Sample IDs: 2352297**

| Component | % Rec.1 | % Rec.2 | Pass/Fail* | Control Limits |
|---|---|---|---|---|
| O-Phosphate-P | 97.1 | 97.9 | P/P | 90 - 110 |

**Reference Method: EPA 8321B**
**Run ID: A114353**
**Included Lab Sample IDs: 2352294**

| Component | % Rec.1 | % Rec.2 | Pass/Fail* | Control Limits |
|---|---|---|---|---|
| Cylindrospermopsin | 87.8 | 79.2 | P/P | 50 - 160 |
| Desmethyl microcystin LR | 76.5 | 82.4 | P/P | 50 - 160 |
| Microcystin HilR | 84.2 | 84.1 | P/P | 50 - 160 |
| Microcystin HtyR | 83.3 | 86.6 | P/P | 50 - 160 |
| Microcystin LA | 79.6 | 87.7 | P/P | 50 - 160 |
| Microcystin LF | 77.3 | 73.4 | P/P | 50 - 160 |
| Microcystin LR | 85.2 | 83.5 | P/P | 50 - 160 |
| Microcystin LW | 73.4 | 74.3 | P/P | 50 - 160 |
| Microcystin LY | 79.7 | 70.0 | P/P | 50 - 160 |
| Microcystin RR | 79.0 | 80.2 | P/P | 50 - 160 |
| Microcystin WR | 78.4 | 87.1 | P/P | 50 - 160 |
| Microcystin YR | 77.5 | 79.2 | P/P | 50 - 160 |
| Nodularin-R | 94.0 | 92.9 | P/P | 50 - 160 |

**Reference Method: FL-PRO 2018**
**Run ID: A114416**
**Included Lab Sample IDs: 2352293**

| Component | % Rec.1 | % Rec.2 | Pass/Fail* | Control Limits |
|---|---|---|---|---|
| TRPH | 96.5 | 92.5 | P/P | 75 - 125 |

**Reference Method: EPA 8270E**
**Run ID: A114431**
**Included Lab Sample IDs: 2352292**

| Component | % Rec.1 | % Rec.2 | Pass/Fail* | Control Limits |
|---|---|---|---|---|
| 1-Methylnaphthalene | 83.6 | | P | 70 - 130 |
| 1,2,4-Trichlorobenzene | 99.5 | | P | 70 - 130 |
| 1,2,4,5-Tetrachlorobenzene | 91.9 | | P | 70 - 130 |
| 1,3-Dinitrobenzene | 75.8 | | P | 70 - 130 |
| 1,3,5-Trinitrobenzene | 74.9 | | P | 70 - 130 |
| 2-Chloronaphthalene | 104 | | P | 70 - 130 |
| 2-Chlorophenol | 92.0 | | P | 70 - 130 |
| 2-Methylnaphthalene | 114 | | P | 70 - 130 |
| 2-Nitroaniline | 90.1 | | P | 70 - 130 |
| 2-Nitrophenol | 104 | | P | 70 - 130 |
| 2,3,4,6-Tetrachlorophenol | 87.7 | | P | 70 - 130 |

# Quality Assurance Report
# Calibration Verification

**Reference Method: EPA 8270E**
**Run ID: A114431**
**Included Lab Sample IDs: 2352292**

| Component | % Rec.1 | % Rec.2 | Pass/Fail* | Control Limits |
|---|---|---|---|---|
| 2,4-Dichlorophenol | 88.0 | | P | 70 - 130 |
| 2,4-Dimethylphenol | 110 | | P | 70 - 130 |
| 2,4-Dinitrotoluene | 99.2 | | P | 70 - 130 |
| 2,4,5-Trichlorophenol | 124 | | P | 70 - 130 |
| 2,4,6-Trichlorophenol | 113 | | P | 70 - 130 |
| 2,6-Dichlorophenol | 99.0 | | P | 70 - 130 |
| 2,6-Dinitrotoluene | 105 | | P | 70 - 130 |
| 3-Methylcholanthrene | 99.4 | | P | 70 - 130 |
| 4-Bromophenyl phenyl ether | 102 | | P | 70 - 130 |
| 4-Chloro-3-methylphenol | 97.8 | | P | 70 - 130 |
| 4-Chlorophenyl phenyl ether | 101 | | P | 70 - 130 |
| 5-Nitro-o-toluidine | 90.1 | | P | 70 - 130 |
| 7,12-Dimethylbenz(a)anthracene | 81.9 | | P | 70 - 130 |
| Acenaphthene | 101 | | P | 70 - 130 |
| Acenaphthylene | 103 | | P | 70 - 130 |
| Acetophenone | 94.9 | | P | 70 - 130 |
| Anthracene | 104 | | P | 70 - 130 |
| Azobenzene/1,2-Diphenylhydrazine | 111 | | P | 70 - 130 |
| Benzo(a)anthracene | 108 | | P | 70 - 130 |
| Benzo(a)pyrene | 106 | | P | 70 - 130 |
| Benzo(b)fluoranthene | 96.3 | | P | 70 - 130 |
| Benzo(g,h,i)perylene | 105 | | P | 70 - 130 |
| Benzo(k)fluoranthene | 99.6 | | P | 70 - 130 |
| Benzyl alcohol | 81.7 | | P | 70 - 130 |
| Bis(2-chloroethoxy)methane | 97.3 | | P | 70 - 130 |
| Bis(2-chloroethyl)ether | 92.7 | | P | 70 - 130 |
| Bis(2-chloroisopropyl)ether | 101 | | P | 70 - 130 |
| Carbazole | 98.3 | | P | 70 - 130 |
| Chrysene | 102 | | P | 70 - 130 |
| Di-n-octyl phthalate | 121 | | P | 70 - 130 |
| Dibenzo(a,h)anthracene | 117 | | P | 70 - 130 |
| Dibenzofuran | 101 | | P | 70 - 130 |
| Dimethyl phthalate | 100 | | P | 70 - 130 |
| Dimethylaminoazobenzene | 88.6 | | P | 70 - 130 |
| Fluoranthene | 103 | | P | 70 - 130 |
| Fluorene | 102 | | P | 70 - 130 |
| Hexachlorobenzene | 96.4 | | P | 70 - 130 |
| Hexachlorobutadiene | 96.6 | | P | 70 - 130 |
| Hexachlorocyclopentadiene | 101 | | P | 70 - 130 |
| Hexachloroethane | 96.0 | | P | 70 - 130 |
| Hexachloropropene | 93.4 | | P | 70 - 130 |
| Indeno(1,2,3-cd)pyrene | 114 | | P | 70 - 130 |
| Isophorone | 111 | | P | 70 - 130 |
| Isosafrole | 95.4 | | P | 70 - 130 |
| m,p-Cresols | 86.6 | | P | 70 - 130 |
| N-Nitrosodi-n-butylamine | 87.4 | | P | 70 - 130 |
| N-Nitrosodi-n-propylamine | 94.9 | | P | 70 - 130 |
| N-Nitrosomorpholine | 88.1 | | P | 70 - 130 |
| N-Nitrosopiperidine | 93.4 | | P | 70 - 130 |
| N-Nitrosopyrrolidine | 76.4 | | P | 70 - 130 |
| Naphthalene | 101 | | P | 70 - 130 |

# Quality Assurance Report
# Calibration Verification

**Reference Method: EPA 8270E**
**Run ID: A114431**
**Included Lab Sample IDs: 2352292**

| Component | % Rec.1 | % Rec.2 | Pass/Fail* | Control Limits |
|---|---|---|---|---|
| Nitrobenzene | 78.4 | | P | 70 - 130 |
| o-Cresol | 95.8 | | P | 70 - 130 |
| o-Toluidine | 105 | | P | 70 - 130 |
| Pentachlorobenzene | 91.7 | | P | 70 - 130 |
| Pentachloroethane | 88.9 | | P | 70 - 130 |
| Pentachloronitrobenzene | 91.2 | | P | 70 - 130 |
| Phenanthrene | 96.7 | | P | 70 - 130 |
| Phenol | 82.4 | | P | 70 - 130 |
| Pyrene | 104 | | P | 70 - 130 |
| Safrole | 89.9 | | P | 70 - 130 |

**Reference Method: EPA 8270E**
**Run ID: A114433**
**Included Lab Sample IDs: 2352292**

| Component | % Rec.1 | % Rec.2 | Pass/Fail* | Control Limits |
|---|---|---|---|---|
| 1-Naphthylamine | 124 | | P | 60 - 130 |
| 2-Acetylaminofluorene | 98.8 | | P | 70 - 150 |
| 2-Methyl-4,6-dinitrophenol | 107 | | P | 70 - 130 |
| 2-Naphthylamine | 88.7 | | P | 60 - 130 |
| 2-Picoline | 111 | | P | 70 - 130 |
| 2,4-Dinitrophenol | 74.6 | | P | 70 - 130 |
| 3,3'-Dichlorobenzidine | 128 | | P | 50 - 130 |
| 4-Aminobiphenyl | 91.9 | | P | 70 - 130 |
| 4-Nitrophenol | 98.3 | | P | 70 - 130 |
| Aniline | 106 | | P | 70 - 130 |
| Benzidine | 77.1 | | P | 50 - 130 |
| Bis(2-ethylhexyl)phthalate | 114 | | P | 70 - 130 |
| Butyl benzyl phthalate | 113 | | P | 70 - 130 |
| Di-n-butyl phthalate | 105 | | P | 70 - 130 |
| Diethyl phthalate | 108 | | P | 70 - 130 |
| Dinoseb | 91.8 | | P | 70 - 130 |
| Ethyl methanesulfonate | 109 | | P | 70 - 130 |
| N-Nitrosodiethylamine | 114 | | P | 70 - 130 |
| N-Nitrosodimethylamine | 128 | | P | 70 - 130 |
| N-Nitrosomethylethylamine | 115 | | P | 70 - 130 |
| Pentachlorophenol | 93.4 | | P | 70 - 130 |
| Phenacetin | 108 | | P | 70 - 130 |
| Pyridine | 113 | | P | 70 - 130 |

**Reference Method: EPA 350.1 Rev. 2.0**
**Run ID: A114437**
**Included Lab Sample IDs: 2352296**

| Component | % Rec.1 | % Rec.2 | Pass/Fail* | Control Limits |
|---|---|---|---|---|
| Ammonia-N | 98.5 | 96.1 | P/P | 90 - 110 |

# Quality Assurance Report
# Calibration Verification

**Reference Method: EPA 351.2 Rev. 2.0**
**Run ID: A114482**
**Included Lab Sample IDs: 2352296**

| Component | % Rec.1 | % Rec.2 | Pass/Fail* | Control Limits |
|---|---|---|---|---|
| Kjeldahl Nitrogen | 94.9 | 99.6 | P/P | 90 - 110 |

**Reference Method: EPA 353.2 Rev. 2.0**
**Run ID: A114510**
**Included Lab Sample IDs: 2352296**

| Component | % Rec.1 | % Rec.2 | Pass/Fail* | Control Limits |
|---|---|---|---|---|
| NO2NO3-N | 103 | 103 | P/P | 90 - 110 |

**Reference Method: EPA 365.1 Rev. 2.0**
**Run ID: A114650**
**Included Lab Sample IDs: 2352296**

| Component | % Rec.1 | % Rec.2 | Pass/Fail* | Control Limits |
|---|---|---|---|---|
| Total-P | 101 | 99.0 | P/P | 90 - 110 |

* Pass/Fail determinations are made for each bracketing calibration verification check.
Control limits for initial calibration checks may be different from those for continuing checks, depending on method requirements.
Where they are different, both control limits are provided.

# Quality Assurance Report
# Summary

| Ref. Method | Analyte | LCS % Recovery | MS % Recovery | | Precision | | |
|---|---|---|---|---|---|---|---|
| | | | | | LCS | SMP | MS |
| EPA 350.1 Rev. 2.0 | Ammonia-N | 96.8 | 95.2 | 95.4 | | | 0.195 |
| EPA 351.2 Rev. 2.0 | Kjeldahl Nitrogen | 104 | 100 | 96.9 | | | 1.70 |
| EPA 353.2 Rev. 2.0 | NO2NO3-N | 102 | 99.5 | 102 | | | 1.19 |
| EPA 365.1 Rev. 2.0 | Total-P | 104 | 104 | 106 | | | 1.83 |
| EPA 365.1 Rev. 2.0 dissolved | O-Phosphate-P | 102 | 101 | 100 | | | 0.661 |
| EPA 8270E | 1-Methylnaphthalene | 92.7 | 89.6 | 93.7 | | | 4.47 |
| | 1-Naphthylamine | 65.2 | 87.5 | 106 | | | 19.4 |
| | 1,2,4-Trichlorobenzene | 84.1 | 78.9 | 84.9 | | | 7.33 |
| | 1,2,4,5-Tetrachlorobenzene | 96.0 | 89.1 | 96.4 | | | 7.87 |
| | 1,3-Dinitrobenzene | 103 | 101 | 97.2 | | | 3.54 |
| | 1,3,5-Trinitrobenzene | 124 | 117 | 112 | | | 4.20 |
| | 2-Acetylaminofluorene | 117 | 121 | 115 | | | 4.67 |
| | 2-Chloronaphthalene | 91.6 | 87.4 | 91.8 | | | 4.91 |
| | 2-Chlorophenol | 87.2 | 81.1 | 86.9 | | | 6.90 |
| | 2-Methyl-4,6-dinitrophenol | 99.3 | 116 | 115 | | | 0.694 |
| | 2-Methylnaphthalene | 92.3 | 87.5 | 92.3 | | | 5.34 |
| | 2-Naphthylamine | 66.1 | 71.5 | 75.7 | | | 5.71 |
| | 2-Nitroaniline | 104 | 98.0 | 103 | | | 4.68 |
| | 2-Nitrophenol | 92.5 | 87.7 | 90.8 | | | 3.47 |
| | 2-Picoline | 80.5 | 92.6 | 99.4 | | | 7.08 |
| | 2,3,4,6-Tetrachlorophenol | 107 | 107 | 104 | | | 3.37 |
| | 2,4-Dichlorophenol | 92.1 | 88.7 | 92.8 | | | 4.52 |
| | 2,4-Dimethylphenol | 88.3 | 88.0 | 91.5 | | | 3.90 |
| | 2,4-Dinitrophenol | 41.8 | 72.6 | 74.4 | | | 2.45 |
| | 2,4-Dinitrotoluene | 99.1 | 96.2 | 99.8 | | | 3.67 |
| | 2,4,5-Trichlorophenol | 87.8 | 88.9 | 90.0 | | | 1.23 |
| | 2,4,6-Trichlorophenol | 90.3 | 88.3 | 90.6 | | | 2.57 |
| | 2,6-Dichlorophenol | 109 | 103 | 109 | | | 5.66 |
| | 2,6-Dinitrotoluene | 99.4 | 95.8 | 100 | | | 4.59 |
| | 3-Methylcholanthrene | 82.1 | 97.2 | 92.4 | | | 5.06 |
| | 3,3'-Dichlorobenzidine | 195 | 169 | 180 | | | 5.85 |
| | 4-Aminobiphenyl | 98.7 | 98.1 | 103 | | | 4.97 |
| | 4-Bromophenyl phenyl ether | 90.3 | 88.1 | 88.2 | | | 0.113 |
| | 4-Chloro-3-methylphenol | 91.1 | 89.4 | 92.1 | | | 2.98 |
| | 4-Chlorophenyl phenyl ether | 98.7 | 94.6 | 100 | | | 5.75 |
| | 4-Nitrophenol | 58.8 | 53.0 | 53.7 | | | 1.31 |
| | 5-Nitro-o-toluidine | 120 | 121 | 114 | | | 6.12 |
| | 7,12-Dimethylbenz(a)anthracene | 106 | 97.4 | 85.6 | | | 12.9 |
| | Acenaphthene | 96.8 | 93.4 | 97.8 | | | 4.60 |
| | Acenaphthylene | 95.8 | 90.8 | 96.1 | | | 5.67 |
| | Acetophenone | 114 | 104 | 114 | | | 9.21 |
| | Aniline | 90.5 | 85.1 | 89.7 | | | 5.26 |
| | Anthracene | 94.3 | 91.3 | 91.5 | | | 0.219 |
| | Azobenzene/1,2-Diphenylhydrazine | 99.1 | 100 | 98.7 | | | 1.51 |
| | Benzidine | 65.0 | 15.6 | 23.9 | | | |
| | Benzo(a)anthracene | 97.8 | 92.7 | 93.4 | | | 0.752 |
| | Benzo(a)pyrene | 92.0 | 88.0 | 96.0 | | | 8.70 |
| | Benzo(b)fluoranthene | 97.3 | 90.7 | 99.2 | | | 8.95 |
| | Benzo(g,h,i)perylene | 74.9 | 79.0 | 81.8 | | | 3.48 |
| | Benzo(k)fluoranthene | 91.7 | 87.7 | 94.5 | | | 7.46 |
| | Benzyl alcohol | 91.6 | 83.2 | 95.1 | | | 13.3 |
| | Bis(2-chloroethoxy)methane | 91.9 | 88.9 | 91.7 | | | 3.10 |
| | Bis(2-chloroethyl)ether | 91.6 | 82.8 | 89.7 | | | 8.00 |
| | Bis(2-chloroisopropyl)ether | 90.1 | 83.8 | 89.3 | | | 6.35 |

# Quality Assurance Report
# Summary

| Ref. Method | Analyte | LCS % Recovery | MS % Recovery | | Precision | | |
|---|---|---|---|---|---|---|---|
| | | | | | LCS | SMP | MS |
| EPA 8270E | Bis(2-ethylhexyl)phthalate | 116 | 114 | 115 | | | 1.05 |
| | Butyl benzyl phthalate | 109 | 105 | 104 | | | 0.478 |
| | Carbazole | 118 | 128 | 126 | | | 1.81 |
| | Chrysene | 98.2 | 92.1 | 91.6 | | | 0.544 |
| | Di-n-butyl phthalate | 102 | 98.0 | 97.5 | | | 0.512 |
| | Di-n-octyl phthalate | 107 | 106 | 111 | | | 4.98 |
| | Dibenzo(a,h)anthracene | 80.9 | 84.4 | 87.0 | | | 3.03 |
| | Dibenzofuran | 96.8 | 92.8 | 97.6 | | | 5.04 |
| | Diethyl phthalate | 110 | 104 | 107 | | | 2.75 |
| | Dimethyl phthalate | 104 | 98.7 | 102 | | | 2.90 |
| | Dimethylaminoazobenzene | 113 | 103 | 93.0 | | | 10.0 |
| | Dinoseb | 126 | 128 | 128 | | | 0.391 |
| | Ethyl methanesulfonate | 96.2 | 89.9 | 97.0 | | | 7.60 |
| | Fluoranthene | 94.5 | 91.1 | 90.6 | | | 0.550 |
| | Fluorene | 100 | 96.0 | 99.9 | | | 3.98 |
| | Hexachlorobenzene | 90.9 | 89.1 | 87.9 | | | 1.36 |
| | Hexachlorobutadiene | 84.5 | 80.6 | 85.1 | | | 5.43 |
| | Hexachlorocyclopentadiene | 51.9 | 43.7 | 49.8 | | | 13.0 |
| | Hexachloroethane | 84.0 | 72.8 | 80.7 | | | 10.3 |
| | Hexachloropropene | 97.0 | 76.3 | 89.5 | | | 15.9 |
| | Indeno(1,2,3-cd)pyrene | 78.8 | 81.7 | 85.3 | | | 4.31 |
| | Isophorone | 96.2 | 92.3 | 96.1 | | | 4.03 |
| | Isosafrole | 108 | 108 | 101 | | | 6.82 |
| | m,p-Cresols | 82.5 | 78.5 | 81.4 | | | 3.63 |
| | N-Nitrosodi-n-butylamine | 107 | 101 | 109 | | | 7.46 |
| | N-Nitrosodi-n-propylamine | 89.1 | 85.1 | 90.8 | | | 6.48 |
| | N-Nitrosodiethylamine | 104 | 95.4 | 103 | | | 7.37 |
| | N-Nitrosodimethylamine | 92.4 | 97.7 | 104 | | | 6.05 |
| | N-Nitrosomethylethylamine | 104 | 95.0 | 103 | | | 8.27 |
| | N-Nitrosomorpholine | 127 | 117 | 126 | | | 7.50 |
| | N-Nitrosopiperidine | 109 | 101 | 109 | | | 7.51 |
| | N-Nitrosopyrrolidine | 96.1 | 88.9 | 94.8 | | | 6.42 |
| | Naphthalene | 90.7 | 85.6 | 91.0 | | | 6.12 |
| | Nitrobenzene | 95.9 | 91.7 | 95.1 | | | 3.64 |
| | o-Cresol | 86.7 | 82.2 | 85.7 | | | 4.17 |
| | o-Toluidine | 111 | 102 | 110 | | | 7.74 |
| | Pentachlorobenzene | 99.7 | 102 | 92.3 | | | 9.69 |
| | Pentachloroethane | 95.8 | 85.1 | 93.9 | | | 9.83 |
| | Pentachloronitrobenzene | 130 | 127 | 124 | | | 2.15 |
| | Pentachlorophenol | 91.5 | 98.4 | 96.7 | | | 1.74 |
| | Phenacetin | 130 | 125 | 122 | | | 2.18 |
| | Phenanthrene | 95.2 | 94.1 | 92.5 | | | 1.71 |
| | Phenol | 54.8 | 51.6 | 54.8 | | | 6.02 |
| | Pyrene | 98.3 | 93.8 | 94.2 | | | 0.426 |
| | Pyridine | 70.4 | 73.5 | 77.5 | | | 5.30 |
| | Safrole | 106 | 100 | 108 | | | 7.11 |
| EPA 8321B | Anatoxin-a | 75.5 | 82.8 | 81.5 | | | 1.55 |
| | Cylindrospermopsin | 79.8 | 91.4 | 88.2 | | | 3.49 |
| | Desmethyl microcystin LR | 73.2 | 106 | 112 | | | 5.15 |
| | Microcystin HilR | 77.2 | 129 | 136 | | | 5.47 |
| | Microcystin HtyR | 78.1 | 108 | 115 | | | 5.96 |
| | Microcystin LA | 80.6 | 98.8 | 94.2 | | | 4.76 |
| | Microcystin LF | 66.4 | 77.2 | 82.0 | | | 6.09 |
| | Microcystin LR | 76.4 | 110 | 115 | | | 3.94 |

# Quality Assurance Report
# Summary

| Ref. Method | Analyte | LCS % Recovery | MS % Recovery | | Precision | | |
|---|---|---|---|---|---|---|---|
| | | | | | LCS | SMP | MS |
| EPA 8321B | Microcystin LW | 64.4 | 94.1 | 98.5 | | | 4.58 |
| | Microcystin LY | 66.1 | 94.5 | 84.8 | | | 10.8 |
| | Microcystin RR | 77.1 | 102 | 107 | | | 4.42 |
| | Microcystin WR | 73.7 | 107 | 115 | | | 7.17 |
| | Microcystin YR | 80.0 | 110 | 116 | | | 4.96 |
| | Nodularin-R | 80.8 | 115 | 118 | | | 3.05 |
| FL-PRO 2018 | TRPH | 90.9 | 64.3 | 83.6 | | | 22.1 |

# Reference Method Descriptions

| Method | Description | Associated Samples |
|---|---|---|
| EPA 350.1 Rev. 2.0 | Ammonia in aqueous matrices as mg N/L | 2352296 |
| EPA 351.2 Rev. 2.0 | Total Kjeldahl Nitrogen in aqueous matrices | 2352296 |
| EPA 353.2 Rev. 2.0 | Nitrite/Nitrate in aqueous matrices as mg N/L | 2352296 |
| EPA 365.1 Rev. 2.0 | Total Phosphorus in aqueous matrices as mg P/L | 2352296 |
| EPA 365.1 Rev. 2.0 dissolved | Ortho-phosphate, dissolved, in filtered, aqueous matrices as mg P/L | 2352297 |
| EPA 8270E | EPA Method 8270, Semi-volatile organic pollutants including PAHs, excluding PCBs and Toxaphene, in water matrices by GC/MS. | 2352292 |
| EPA 8321B | Microcystins in water matrices by HPLC/MS/MS | 2352294 |
| EPA 8321B | Saxitoxin in water matrices by HPLC/MS/MS | 2352294 |
| FL-PRO 2018 | Total recoverable petroleum hydrocarbons in water samples by GC-FID. | 2352293 |

# Preparation and Analysis Log

| Ref. Method | Received Date | Prep Date/Time | Prepared By | Analysis Date/Time | Analyzed By | Associated Samples |
|---|---|---|---|---|---|---|
| EPA 350.1 Rev. 2.0 | 08/25/2022 | | | 08/30/2022 14:42 | Judith K. Clark | 2352296 |
| EPA 351.2 Rev. 2.0 | 08/25/2022 | 08/30/2022 14:31 | Samantha L Bates | 09/01/2022 10:35 | Alexandra J Mattheus | 2352296 |
| EPA 353.2 Rev. 2.0 | 08/25/2022 | | | 09/02/2022 09:09 | Judith K. Clark | 2352296 |
| EPA 365.1 Rev. 2.0 | 08/25/2022 | 09/07/2022 16:10 | Simonne.V. Calhoun | 09/09/2022 13:04 | James L Waggerby | 2352296 |
| EPA 365.1 Rev. 2.0 dissolved | 08/25/2022 | | | 08/25/2022 14:41 | Elise M. Gillis | 2352297 |
| EPA 8270E | 08/25/2022 | 08/29/2022 09:00 | Hoor Shaik | 08/30/2022 11:51 | Mohammad Ghaffari | 2352292 |
| | 08/25/2022 | 08/29/2022 09:00 | Hoor Shaik | 08/30/2022 12:37 | Mohammad Ghaffari | 2352292 |
| EPA 8321B | 08/25/2022 | 08/25/2022 13:00 | Manjita Shrestha | 08/25/2022 20:10 | Manjita Shrestha | 2352294 |
| | 08/25/2022 | 08/25/2022 13:00 | Manjita Shrestha | 08/25/2022 23:34 | Manjita Shrestha | 2352294 |
| FL-PRO 2018 | 08/25/2022 | 08/26/2022 08:00 | Lipi Saha | 08/29/2022 13:42 | Lipi Saha | 2352293 |