

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP021524\  
 Data File : BP019732.D  
 Acq On : 15 Feb 2024 20:18  
 Operator : MA/JU  
 Sample : P1395-04DL 10X  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 B-2(6-8)DL

Quant Time: Feb 16 00:55:33 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP013124.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Feb 01 00:51:48 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.905  | 152  | 75695    | 20.000 | ng    | -0.01    |
| 21) Naphthalene-d8            | 10.704 | 136  | 291828   | 20.000 | ng    | -0.01    |
| 39) Acenaphthene-d10          | 14.540 | 164  | 212826   | 20.000 | ng    | -0.01    |
| 64) Phenanthrene-d10          | 17.298 | 188  | 456363   | 20.000 | ng    | 0.00     |
| 76) Chrysene-d12              | 21.386 | 240  | 414544   | 20.000 | ng    | 0.00     |
| 86) Perylene-d12              | 23.810 | 264  | 479144   | 20.000 | ng    | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 5) 2-Fluorophenol             | 5.481  | 112  | 22316    | 4.752  | ng    | 0.00     |
| 7) Phenol-d6                  | 7.075  | 99   | 29886    | 4.720  | ng    | -0.01    |
| 23) Nitrobenzene-d5           | 9.069  | 82   | 21861    | 3.246  | ng    | -0.02    |
| 42) 2,4,6-Tribromophenol      | 16.034 | 330  | 18971    | 6.105  | ng    | 0.00     |
| 45) 2-Fluorobiphenyl          | 13.163 | 172  | 53610    | 3.114  | ng    | -0.02    |
| 79) Terphenyl-d14             | 19.857 | 244  | 96386    | 3.824  | ng    | -0.01    |
| Target Compounds              |        |      |          |        |       |          |
|                               |        |      |          |        |       | Qvalue   |
| 10) Phenol                    | 7.105  | 94   | 338434   | 53.628 | ng    | 99       |
| 13) 1,4-Dichlorobenzene       | 7.940  | 146  | 30679    | 5.133  | ng    | 97       |
| 17) 2-Methylphenol            | 8.352  | 107  | 113320   | 26.672 | ng    | 99       |
| 20) 3+4-Methylphenols         | 8.675  | 107  | 167423   | 28.874 | ng    | 95       |
| 27) 2,4-Dimethylphenol        | 9.887  | 122  | 107644   | 32.559 | ng    | 96       |
| 31) Naphthalene               | 10.751 | 128  | 35629    | 2.226  | ng    | 97       |
| 46) 1,1'-Biphenyl             | 13.375 | 154  | 234570   | 13.809 | ng    | 98       |
| 74) Di-n-butylphthalate       | 18.263 | 149  | 443046   | 16.835 | ng    | 99       |
| 84) Bis(2-ethylhexyl)phtha... | 21.316 | 149  | 1490619  | 89.595 | ng    | 98       |
| -----                         |        |      |          |        |       |          |

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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