

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP021624\  
 Data File : BP019750.D  
 Acq On : 16 Feb 2024 18:06  
 Operator : MA/JU  
 Sample : P1414-07  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SB-10(10-11)

Quant Time: Feb 19 00:27:18 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   | 7.899  | 152  | 65620    | 20.000  | ng    | -0.02    |
| 21) Naphthalene-d8          | 10.698 | 136  | 238644   | 20.000  | ng    | -0.02    |
| 39) Acenaphthene-d10        | 14.534 | 164  | 161484   | 20.000  | ng    | -0.02    |
| 64) Phenanthrene-d10        | 17.292 | 188  | 409603   | 20.000  | ng    | -0.01    |
| 76) Chrysene-d12            | 21.386 | 240  | 440178   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 23.810 | 264  | 491630   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.475  | 112  | 550041   | 135.115 | ng    | -0.01    |
| 7) Phenol-d6                | 7.075  | 99   | 696605   | 126.907 | ng    | -0.01    |
| 23) Nitrobenzene-d5         | 9.069  | 82   | 482755   | 87.656  | ng    | -0.02    |
| 42) 2,4,6-Tribromophenol    | 16.033 | 330  | 450459   | 191.052 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.163 | 172  | 1134247  | 86.831  | ng    | -0.02    |
| 79) Terphenyl-d14           | 19.863 | 244  | 2489844  | 93.023  | ng    | 0.00     |

|                  |        |
|------------------|--------|
| Target Compounds | Qvalue |
| -----            |        |

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

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