

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF102824\  
 Data File : BF140081.D  
 Acq On : 28 Oct 2024 12:27  
 Operator : RC/JU  
 Sample : P4547-08  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 BP-F-20

Quant Time: Oct 28 13:01:19 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF101824.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Oct 18 15:07:50 2024  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.887  | 152  | 105972   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.169  | 136  | 413518   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.922  | 164  | 227492   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.410 | 188  | 389316   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.045 | 240  | 199247   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.527 | 264  | 233774   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.516  | 112  | 934938   | 138.115 | ng    | 0.02     |
| 7) Phenol-d6                | 6.510  | 99   | 1149935  | 131.162 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.445  | 82   | 794686   | 106.511 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.710 | 330  | 348850   | 163.942 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.239  | 172  | 1429396  | 103.844 | ng    | -0.01    |
| 79) Terphenyl-d14           | 12.998 | 244  | 1312236  | 107.317 | ng    | 0.00     |

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

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

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