

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF081024\  
 Data File : BF138922.D  
 Acq On : 10 Aug 2024 21:25  
 Operator : RC/JU  
 Sample : P3467-01  
 Misc :  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 919-J-WS-080224

Quant Time: Aug 12 01:20:17 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF073024.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Jul 30 17:50:01 2024  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.840  | 152  | 37654    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.122  | 136  | 147641   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.869  | 164  | 79728    | 20.000  | ng    | -0.01    |
| 64) Phenanthrene-d10        | 11.357 | 188  | 124633   | 20.000  | ng    | -0.01    |
| 76) Chrysene-d12            | 13.998 | 240  | 68405    | 20.000  | ng    | # 0.00   |
| 86) Perylene-d12            | 15.462 | 264  | 79740    | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.475  | 112  | 181837   | 74.545  | ng    | 0.00     |
| 7) Phenol-d6                | 6.487  | 99   | 154533   | 47.186  | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.404  | 82   | 303638   | 100.550 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.663 | 330  | 96865    | 148.320 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.192  | 172  | 526889   | 99.294  | ng    | -0.01    |
| 79) Terphenyl-d14           | 12.939 | 244  | 405048   | 99.139  | ng    | 0.00     |

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

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

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