

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF080324\  
 Data File : BF138762.D  
 Acq On : 03 Aug 2024 11:02  
 Operator : RC/JU  
 Sample : P3379-03RX  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 OU4-PCS-02-072324RX

Quant Time: Aug 05 00:49:01 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   | 6.845  | 152  | 53746    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.128  | 136  | 242226   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.880  | 164  | 143809   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.369 | 188  | 209060   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.010 | 240  | 126686   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.480 | 264  | 130904   | 20.000  | ng    | 0.01     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.487  | 112  | 431008   | 123.791 | ng    | 0.02     |
| 7) Phenol-d6                | 6.492  | 99   | 603774   | 129.161 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.410  | 82   | 418614   | 84.494  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.674 | 330  | 167014   | 141.779 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.204  | 172  | 788398   | 82.371  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.957 | 244  | 727767   | 96.181  | ng    | 0.01     |

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

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

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