

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF012924\  
 Data File : BF137187.D  
 Acq On : 29 Jan 2024 12:02  
 Operator : CG\JU  
 Sample : PB158518BL  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB158518BL

Quant Time: Jan 29 12:23:54 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF012724.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Sat Jan 27 02:14:33 2024  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.798  | 152  | 59873    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.075  | 136  | 236959   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.834  | 164  | 141438   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.328 | 188  | 286184   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 13.980 | 240  | 218478   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.468 | 264  | 166158   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.440  | 112  | 358853   | 104.127 | ng    | 0.01     |
| 7) Phenol-d6                | 6.428  | 99   | 472818   | 96.142  | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.363  | 82   | 457651   | 87.615  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.628 | 330  | 127921   | 80.951  | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.157  | 172  | 841808   | 83.563  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.933 | 244  | 1146365  | 73.595  | ng    | 0.00     |

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

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

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