

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM111422\  
 Data File : BM037500.D  
 Acq On : 14 Nov 2022 11:41  
 Operator : CG/JU  
 Sample : PB148922BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 PB148922BL

Quant Time: Nov 15 00:01:31 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM111222.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Nov 14 01:42:00 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.792  | 152  | 185832   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.592 | 136  | 775364   | 20.000  | ng    | -0.01    |
| 39) Acenaphthene-d10        | 14.439 | 164  | 510359   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.186 | 188  | 1119886  | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.374 | 240  | 1212566  | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 23.709 | 264  | 1249925  | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.363  | 112  | 1829723  | 185.939 | ng    | 0.00     |
| 7) Phenol-d6                | 6.975  | 99   | 2415420  | 170.602 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 8.963  | 82   | 1422217  | 95.320  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 15.933 | 330  | 1184348  | 167.831 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.063 | 172  | 3561364  | 96.981  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.815 | 244  | 7004023  | 115.359 | ng    | 0.00     |

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

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

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