

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM060222\  
 Data File : BM035368.D  
 Acq On : 02 Jun 2022 18:34  
 Operator : CG/JU  
 Sample : PB145245BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 PB145245BL

Quant Time: Jun 03 01:43:11 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM060222.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Jun 02 16:16:23 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.851  | 152  | 153323   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.651 | 136  | 624040   | 20.000  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 14.492 | 164  | 478660   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.239 | 188  | 1165684  | 20.000  | ng    | # 0.00   |
| 76) Chrysene-d12            | 21.421 | 240  | 1266935  | 20.000  | ng    | # 0.00   |
| 86) Perylene-d12            | 23.780 | 264  | 1274337  | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.440  | 112  | 1156739  | 158.533 | ng    | 0.00     |
| 7) Phenol-d6                | 7.040  | 99   | 1495774  | 150.076 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.016  | 82   | 884174   | 89.965  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 15.986 | 330  | 1141421  | 148.623 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.122 | 172  | 2974194  | 89.244  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.868 | 244  | 6014190  | 94.574  | ng    | 0.00     |

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

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

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