

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM060122\  
 Data File : BM035341.D  
 Acq On : 01 Jun 2022 11:29  
 Operator : CG/JU  
 Sample : PB145220BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 PB145220BL

Quant Time: Jun 02 05:47:34 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM052722.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue May 31 15:35:47 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.863  | 152  | 99708    | 20.000  | ng    | -0.01    |
| 21) Naphthalene-d8          | 10.663 | 136  | 386626   | 20.000  | ng    | #-0.01   |
| 39) Acenaphthene-d10        | 14.504 | 164  | 269833   | 20.000  | ng    | -0.01    |
| 64) Phenanthrene-d10        | 17.245 | 188  | 624556   | 20.000  | ng    | -0.01    |
| 76) Chrysene-d12            | 21.433 | 240  | 660358   | 20.000  | ng    | # 0.00   |
| 86) Perylene-d12            | 23.797 | 264  | 666843   | 20.000  | ng    | -0.01    |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.452  | 112  | 790357   | 146.357 | ng    | 0.00     |
| 7) Phenol-d6                | 7.046  | 99   | 991672   | 141.819 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.022  | 82   | 596550   | 84.937  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 15.998 | 330  | 604839   | 172.785 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.127 | 172  | 1801438  | 92.581  | ng    | -0.01    |
| 79) Terphenyl-d14           | 19.874 | 244  | 3460238  | 92.852  | ng    | 0.00     |

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

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

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