

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM110423\  
 Data File : BM042592.D  
 Acq On : 04 Nov 2023 15:29  
 Operator : MA/JU  
 Sample : 05234-03  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SP-B

Quant Time: Nov 06 01:45:21 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM103023.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Nov 03 15:52:20 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.910  | 152  | 45619    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.733 | 136  | 170892   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.563 | 164  | 116429   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.321 | 188  | 300016   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.503 | 240  | 279199   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 23.921 | 264  | 314206   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.446  | 112  | 334247   | 118.248 | ng    | 0.00     |
| 7) Phenol-d6                | 7.081  | 99   | 433249   | 111.758 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.116  | 82   | 299719   | 81.758  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.062 | 330  | 234908   | 154.537 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.186 | 172  | 668184   | 77.224  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.939 | 244  | 1376346  | 83.587  | ng    | 0.00     |

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

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

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