

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM102723\  
 Data File : BM042479.D  
 Acq On : 27 Oct 2023 19:10  
 Operator : MA/JU  
 Sample : PB156597TB  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 PB156597TB

Quant Time: Oct 28 00:27:59 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM102123.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Sat Oct 28 00:26:13 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.986  | 152  | 145457   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.810 | 136  | 528407   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.627 | 164  | 290556   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.380 | 188  | 534174   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.562 | 240  | 363930   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 24.021 | 264  | 348157   | 20.000  | ng    | # 0.00   |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.534  | 112  | 1322703  | 128.519 | ng    | 0.00     |
| 7) Phenol-d6                | 7.163  | 99   | 1615158  | 117.105 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.192  | 82   | 1044986  | 87.449  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.127 | 330  | 473314   | 146.260 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.251 | 172  | 2046182  | 94.832  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.992 | 244  | 2434638  | 118.644 | ng    | 0.00     |

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

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

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