

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG042622\  
 Data File : BG053322.D  
 Acq On : 26 Apr 2022 16:12  
 Operator : CG/JU  
 Sample : PB144364TB  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 PB144364TB

Quant Time: Apr 27 03:37:24 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG040822.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Sun Apr 24 01:13:30 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.107  | 152  | 26402    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.915 | 136  | 107739   | 20.000  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 14.727 | 164  | 82946    | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.476 | 188  | 211940   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.753 | 240  | 231993   | 20.000  | ng    | -0.01    |
| 86) Perylene-d12            | 25.054 | 264  | 221883   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.675  | 112  | 219433   | 142.470 | ng    | 0.00     |
| 7) Phenol-d6                | 7.267  | 99   | 313674   | 132.064 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.270  | 82   | 225314   | 84.914  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.219 | 330  | 161613   | 122.468 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.352 | 172  | 505559   | 84.045  | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.079 | 244  | 1136722  | 82.753  | ng    | 0.00     |

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

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

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