

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF072522\  
 Data File : BF129338.D  
 Acq On : 25 Jul 2022 18:24  
 Operator : CG\JU  
 Sample : PB146401BL  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB146401BL

Quant Time: Jul 26 03:10:34 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF072522.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Jul 25 17:24:17 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.881  | 152  | 272878   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.163  | 136  | 1134040  | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.922  | 164  | 610048   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.416 | 188  | 1057393  | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.057 | 240  | 831449   | 20.000  | ng    | #-0.01   |
| 86) Perylene-d12            | 15.551 | 264  | 707101   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.510  | 112  | 1977992  | 118.080 | ng    | 0.01     |
| 7) Phenol-d6                | 6.510  | 99   | 2502996  | 118.877 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.445  | 82   | 1644575  | 79.164  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.716 | 330  | 674667   | 115.030 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.245  | 172  | 2769483  | 70.228  | ng    | 0.00     |
| 79) Terphenyl-d14           | 13.004 | 244  | 3093996  | 70.204  | ng    | 0.00     |

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

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

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