

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF042822\  
 Data File : BF127975.D  
 Acq On : 28 Apr 2022 12:19  
 Operator : CG\JU  
 Sample : PB144382BL  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB144382BL

Quant Time: Apr 28 17:23:20 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF042222.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Apr 28 05:13:03 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.928  | 152  | 142510   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.216  | 136  | 566768   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.975  | 164  | 322153   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.463 | 188  | 578047   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.110 | 240  | 393241   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.621 | 264  | 310262   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.569  | 112  | 1071776  | 119.789 | ng    | 0.02     |
| 7) Phenol-d6                | 6.557  | 99   | 1414124  | 121.806 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.498  | 82   | 1047313  | 86.894  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.769 | 330  | 379867   | 117.134 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.292  | 172  | 1660204  | 87.183  | ng    | 0.00     |
| 79) Terphenyl-d14           | 13.057 | 244  | 1785987  | 83.048  | ng    | 0.00     |

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

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

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