

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF040722\  
 Data File : BF127683.D  
 Acq On : 07 Apr 2022 14:47  
 Operator : CG\JU  
 Sample : PB143804BL  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB143804BL

Quant Time: Apr 07 16:54:50 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF040622.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Apr 06 13:55:27 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.963  | 152  | 145251   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.245  | 136  | 520948   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 10.010 | 164  | 294542   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.504 | 188  | 549203   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.151 | 240  | 445498   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.674 | 264  | 328252   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.593  | 112  | 936084   | 104.718 | ng    | 0.01     |
| 7) Phenol-d6                | 6.581  | 99   | 1207341  | 103.243 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.528  | 82   | 952204   | 81.487  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.804 | 330  | 391550   | 124.922 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.328  | 172  | 1613137  | 77.478  | ng    | 0.00     |
| 79) Terphenyl-d14           | 13.098 | 244  | 1913467  | 70.895  | ng    | 0.00     |

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

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

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