

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF041922\  
 Data File : BF127847.D  
 Acq On : 19 Apr 2022 18:13  
 Operator : CG\JU  
 Sample : PB144163TB  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB144163TB

Quant Time: Apr 20 03:54:09 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF041922.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Apr 19 17:31:18 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.940  | 152  | 88043    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.222  | 136  | 323292   | 20.000  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 9.986  | 164  | 197947   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.480 | 188  | 375194   | 20.000  | ng    | # 0.00   |
| 76) Chrysene-d12            | 14.121 | 240  | 334039   | 20.000  | ng    | -0.01    |
| 86) Perylene-d12            | 15.639 | 264  | 381550   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.581  | 112  | 748046   | 133.106 | ng    | 0.01     |
| 7) Phenol-d6                | 6.569  | 99   | 1194217  | 170.163 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.510  | 82   | 684385   | 98.177  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.781 | 330  | 338362   | 140.321 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.304  | 172  | 1275289  | 84.150  | ng    | 0.00     |
| 79) Terphenyl-d14           | 13.074 | 244  | 1613065  | 75.797  | ng    | 0.00     |

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

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

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