

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF053024\  
 Data File : BF138069.D  
 Acq On : 30 May 2024 11:59  
 Operator : CG\JU  
 Sample : PB161142TB  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB161142TB

Quant Time: May 30 12:26:50 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF051424.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed May 15 04:12:16 2024  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.034  | 152  | 93253    | 20.000  | ng    | -0.01    |
| 21) Naphthalene-d8          | 8.316  | 136  | 367067   | 20.000  | ng    | -0.01    |
| 39) Acenaphthene-d10        | 10.080 | 164  | 209402   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.580 | 188  | 347506   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.233 | 240  | 186339   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.804 | 264  | 181628   | 20.000  | ng    | 0.02     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.675  | 112  | 719210   | 131.007 | ng    | 0.01     |
| 7) Phenol-d6                | 6.657  | 99   | 853121   | 126.153 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.598  | 82   | 533829   | 85.587  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 10.875 | 330  | 298126   | 141.198 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.398  | 172  | 1211340  | 89.108  | ng    | 0.00     |
| 79) Terphenyl-d14           | 13.168 | 244  | 1178844  | 93.238  | ng    | 0.00     |

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

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

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