

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF020624\  
 Data File : BF137293.D  
 Acq On : 06 Feb 2024 15:49  
 Operator : CG\JU  
 Sample : PB158779TB  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB158779TB

Quant Time: Feb 06 16:14:57 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF013024.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Jan 31 04:56:22 2024  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.787  | 152  | 63984    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.063  | 136  | 251830   | 20.000  | ng    | -0.01    |
| 39) Acenaphthene-d10        | 9.822  | 164  | 153163   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.316 | 188  | 315173   | 20.000  | ng    | # 0.00   |
| 76) Chrysene-d12            | 13.974 | 240  | 278617   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.457 | 264  | 255023   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.434  | 112  | 579759   | 149.514 | ng    | 0.02     |
| 7) Phenol-d6                | 6.428  | 99   | 762768   | 137.839 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.351  | 82   | 512594   | 91.132  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.616 | 330  | 257562   | 148.901 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.145  | 172  | 951615   | 85.409  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.921 | 244  | 1410467  | 68.820  | ng    | 0.00     |

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

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

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