

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF093025\  
 Data File : BF143838.D  
 Acq On : 30 Sep 2025 17:36  
 Operator : RC/JU  
 Sample : Q3179-01  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 EME-TS14-01

Quant Time: Sep 30 17:58:19 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF092325.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Sep 24 05:00:00 2025  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.881  | 152  | 143060   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.163  | 136  | 546052   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.922  | 164  | 302570   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.410 | 188  | 531398   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.057 | 240  | 566979   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.533 | 264  | 571887   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.493  | 112  | 513443   | 58.311  | ng    | 0.00     |
| 7) Phenol-d6                | 6.498  | 99   | 391015   | 38.433  | ng    | -0.01    |
| 23) Nitrobenzene-d5         | 7.446  | 82   | 827470   | 92.758  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.710 | 330  | 621128   | 138.644 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.245  | 172  | 1660843  | 85.657  | ng    | 0.00     |
| 79) Terphenyl-d14           | 13.004 | 244  | 2769182  | 99.330  | ng    | 0.00     |

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

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

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