

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF070925\  
 Data File : BF143051.D  
 Acq On : 09 Jul 2025 12:24  
 Operator : RC/JU  
 Sample : PB168750BL  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB168750BL

Quant Time: Jul 09 12:45:41 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF061925.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Jun 24 05:28:21 2025  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.869  | 152  | 58608    | 20.000  | ng    | -0.01    |
| 21) Naphthalene-d8          | 8.157  | 136  | 223429   | 20.000  | ng    | -0.01    |
| 39) Acenaphthene-d10        | 9.916  | 164  | 120074   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.404 | 188  | 209574   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.051 | 240  | 109296   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.545 | 264  | 119678   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.504  | 112  | 441344   | 125.376 | ng    | 0.01     |
| 7) Phenol-d6                | 6.504  | 99   | 522633   | 125.842 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.434  | 82   | 322776   | 79.240  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 10.710 | 330  | 159473   | 129.061 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.233  | 172  | 719775   | 78.747  | ng    | -0.01    |
| 79) Terphenyl-d14           | 12.992 | 244  | 683043   | 86.349  | ng    | 0.00     |

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

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

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