

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF111925\  
 Data File : BF144282.D  
 Acq On : 19 Nov 2025 17:20  
 Operator : RC/JU  
 Sample : PB170637BL  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB170637BL

Quant Time: Nov 19 18:02:11 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF110525.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Nov 05 18:37:33 2025  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.869  | 152  | 51808    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.145  | 136  | 195866   | 20.000  | ng    | -0.01    |
| 39) Acenaphthene-d10        | 9.904  | 164  | 101180   | 20.000  | ng    | -0.01    |
| 64) Phenanthrene-d10        | 11.398 | 188  | 150413   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.045 | 240  | 144217   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.521 | 264  | 166107   | 20.000  | ng    | -0.02    |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.498  | 112  | 412517   | 124.602 | ng    | 0.00     |
| 7) Phenol-d6                | 6.498  | 99   | 447040   | 119.170 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.428  | 82   | 290178   | 85.565  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.698 | 330  | 131756   | 103.770 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.222  | 172  | 599854   | 87.561  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.986 | 244  | 586441   | 64.591  | ng    | 0.00     |

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

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

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