

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP091225\  
 Data File : BP025743.D  
 Acq On : 12 Sep 2025 14:30  
 Operator : CG/JU  
 Sample : PB169499BL  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 PB169499BL

Quant Time: Sep 12 15:00:14 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP091225.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Sep 11 16:45:04 2025  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.754  | 152  | 134957   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.531 | 136  | 525832   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.378 | 164  | 350062   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.183 | 188  | 778197   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.618 | 240  | 870891   | 20.000  | ng    | -0.02    |
| 86) Perylene-d12            | 24.971 | 264  | 895941   | 20.000  | ng    | -0.03    |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.384  | 112  | 1215331  | 145.305 | ng    | 0.00     |
| 7) Phenol-d6                | 6.943  | 99   | 1534418  | 138.021 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 8.901  | 82   | 1013516  | 85.022  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 15.889 | 330  | 602486   | 137.985 | ng    | -0.02    |
| 45) 2-Fluorobiphenyl        | 12.984 | 172  | 2234692  | 84.999  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.907 | 244  | 3998630  | 82.599  | ng    | 0.00     |

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

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

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