

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG092625\  
 Data File : BG064457.D  
 Acq On : 26 Sep 2025 17:04  
 Operator : RC/JU  
 Sample : PB169873BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 PB169873BL

Quant Time: Sep 26 18:07:39 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG092525.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Sep 26 06:30:03 2025  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.836  | 152  | 15014    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.627 | 136  | 61808    | 20.000  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 14.463 | 164  | 50020    | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.201 | 188  | 133410   | 20.000  | ng    | # 0.00   |
| 76) Chrysene-d12            | 21.432 | 240  | 146948   | 20.000  | ng    | -0.01    |
| 86) Perylene-d12            | 24.428 | 264  | 166371   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.427  | 112  | 114037   | 145.144 | ng    | 0.00     |
| 7) Phenol-d6                | 7.013  | 99   | 146962   | 136.534 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 8.993  | 82   | 101843   | 89.299  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 15.956 | 330  | 122611   | 140.411 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.088 | 172  | 300545   | 88.724  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.828 | 244  | 714875   | 95.900  | ng    | 0.00     |

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

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

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