

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF091025\  
 Data File : BF143698.D  
 Acq On : 10 Sep 2025 12:41  
 Operator : RC/JU  
 Sample : SP6869  
 Misc : 8270 SURROGATE  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SP6869

Quant Time: Sep 10 13:29:58 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF091025.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Sep 09 15:20:16 2025  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.886  | 152  | 43241    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.169  | 136  | 162291   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.922  | 164  | 96334    | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.410 | 188  | 160186   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.045 | 240  | 104506   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.515 | 264  | 95807    | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.510  | 112  | 361628   | 142.686 | ng    | 0.01     |
| 7) Phenol-d6                | 6.504  | 99   | 434497   | 140.800 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.445  | 82   | 255138   | 97.786  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.710 | 330  | 141277   | 145.522 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.245  | 172  | 600085   | 94.099  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.998 | 244  | 635346   | 98.520  | ng    | 0.00     |

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

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

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