

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061825\  
 Data File : BP024997.D  
 Acq On : 18 Jun 2025 16:49  
 Operator : RC/JU  
 Sample : Q2325-04  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 TP-8

Quant Time: Jun 18 18:41:31 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP060625.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Jun 06 16:20:27 2025  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.602  | 152  | 201735   | 20.000  | ng    | -0.01    |
| 21) Naphthalene-d8          | 10.378 | 136  | 757706   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.254 | 164  | 455205   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.066 | 188  | 912049   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.513 | 240  | 1019432  | 20.000  | ng    | 0.03     |
| 86) Perylene-d12            | 24.789 | 264  | 1159619  | 20.000  | ng    | 0.07     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.237  | 112  | 1601228  | 132.489 | ng    | 0.00     |
| 7) Phenol-d6                | 6.819  | 99   | 1771928  | 110.810 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 8.760  | 82   | 1340161  | 85.947  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 15.790 | 330  | 916470   | 145.623 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 12.854 | 172  | 2828835  | 83.716  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.789 | 244  | 4978268  | 87.522  | ng    | 0.00     |

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

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

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