

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP041025\  
 Data File : BP024250.D  
 Acq On : 10 Apr 2025 13:15  
 Operator : RC/JU  
 Sample : Q1740-04  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 TP-20

Quant Time: Apr 10 13:46:45 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP031225.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Apr 10 11:43:37 2025  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.740  | 152  | 279443   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.516 | 136  | 1104872  | 20.000  | ng    | -0.01    |
| 39) Acenaphthene-d10        | 14.375 | 164  | 652312   | 20.000  | ng    | -0.02    |
| 64) Phenanthrene-d10        | 17.175 | 188  | 1268299  | 20.000  | ng    | -0.02    |
| 76) Chrysene-d12            | 21.622 | 240  | 1378986  | 20.000  | ng    | -0.02    |
| 86) Perylene-d12            | 24.998 | 264  | 1514466  | 20.000  | ng    | -0.01    |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.358  | 112  | 2349415  | 134.207 | ng    | 0.00     |
| 7) Phenol-d6                | 6.922  | 99   | 2731577  | 116.685 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 8.887  | 82   | 1869573  | 95.471  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 15.887 | 330  | 1339446  | 163.015 | ng    | -0.01    |
| 45) 2-Fluorobiphenyl        | 12.987 | 172  | 4054292  | 91.646  | ng    | -0.01    |
| 79) Terphenyl-d14           | 19.898 | 244  | 6331251  | 94.773  | ng    | -0.02    |

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

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

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