

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN040825\  
 Data File : BN036861.D  
 Acq On : 08 Apr 2025 15:31  
 Operator : RC/JU  
 Sample : Q1502-08DL2 100X  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 PT-ACIDS-WPDL2

Quant Time: Apr 08 15:57:52 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\8270-SIM-BN031025.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Apr 04 17:30:43 2025  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc  | Units | Dev(Min) |
|-------------------------------|--------|------|----------|-------|-------|----------|
| -----                         |        |      |          |       |       |          |
| Internal Standards            |        |      |          |       |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.681  | 152  | 1573     | 0.400 | ng    | 0.00     |
| 7) Naphthalene-d8             | 10.466 | 136  | 3831     | 0.400 | ng    | -0.01    |
| 13) Acenaphthene-d10          | 14.323 | 164  | 2049     | 0.400 | ng    | -0.01    |
| 19) Phenanthrene-d10          | 17.074 | 188  | 3999     | 0.400 | ng    | # 0.00   |
| 29) Chrysene-d12              | 21.277 | 240  | 3589     | 0.400 | ng    | 0.00     |
| 35) Perylene-d12              | 23.513 | 264  | 3878     | 0.400 | ng    | 0.00     |
| System Monitoring Compounds   |        |      |          |       |       |          |
| 4) 2-Fluorophenol             | 5.276  | 112  | 6208     | 1.693 | ng    | 0.00     |
| 5) Phenol-d6                  | 6.858  | 99   | 7361     | 1.626 | ng    | 0.00     |
| 8) Nitrobenzene-d5            | 8.832  | 82   | 4095     | 0.983 | ng    | -0.01    |
| 11) 2-Methylnaphthalene-d10   | 0.000  | 152  | 0d       | 0.000 | ng    |          |
| 14) 2,4,6-Tribromophenol      | 15.820 | 330  | 1306     | 1.405 | ng    | -0.01    |
| 15) 2-Fluorobiphenyl          | 12.947 | 172  | 11192    | 0.939 | ng    | -0.01    |
| 27) Fluoranthene-d10          | 0.000  | 212  | 0d       | 0.000 | ng    |          |
| 31) Terphenyl-d14             | 19.712 | 244  | 8852     | 1.029 | ng    | 0.00     |
| Target Compounds              |        |      |          |       |       |          |
| 20) 4,6-Dinitro-2-methylph... | 15.467 | 198  | 801      | 0.919 | ng    | # 67     |
| 24) Pentachlorophenol         | 16.726 | 266  | 1170     | 0.848 | ng    | 96       |
| -----                         |        |      |          |       |       |          |

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

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