

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF060425\  
 Data File : BF142612.D  
 Acq On : 04 Jun 2025 17:48  
 Operator : RC/JU  
 Sample : Q2160-08  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 TP05-MHH-WC

Quant Time: Jun 04 18:21:12 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF052025.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue May 20 16:26:47 2025  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.898  | 152  | 122107   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.181  | 136  | 455092   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.933  | 164  | 229398   | 20.000  | ng    | -0.01    |
| 64) Phenanthrene-d10        | 11.422 | 188  | 334337   | 20.000  | ng    | -0.01    |
| 76) Chrysene-d12            | 14.063 | 240  | 219434   | 20.000  | ng    | -0.01    |
| 86) Perylene-d12            | 15.557 | 264  | 271463   | 20.000  | ng    | -0.01    |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.528  | 112  | 815161   | 112.471 | ng    | 0.01     |
| 7) Phenol-d6                | 6.522  | 99   | 894051   | 102.475 | ng    | -0.01    |
| 23) Nitrobenzene-d5         | 7.457  | 82   | 646508   | 77.464  | ng    | -0.02    |
| 42) 2,4,6-Tribromophenol    | 10.727 | 330  | 290299   | 114.020 | ng    | -0.01    |
| 45) 2-Fluorobiphenyl        | 9.251  | 172  | 1279984  | 74.873  | ng    | -0.02    |
| 79) Terphenyl-d14           | 13.010 | 244  | 969529   | 60.378  | ng    | 0.00     |

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

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

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