

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG111725\  
 Data File : BG064791.D  
 Acq On : 17 Nov 2025 19:26  
 Operator : RC/JU  
 Sample : Q3647-02  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 VNJ-212

Quant Time: Nov 18 00:23:13 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG111225.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Nov 12 17:10:56 2025  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.165  | 152  | 37326    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.979 | 136  | 132061   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.775 | 164  | 102733   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.501 | 188  | 258106   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.749 | 240  | 351977   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 25.010 | 264  | 357694   | 20.000  | ng    | -0.01    |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.703  | 112  | 269462   | 126.062 | ng    | 0.00     |
| 7) Phenol-d6                | 7.307  | 99   | 323328   | 113.727 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.328  | 82   | 241596   | 90.452  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.250 | 330  | 270765   | 135.139 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.406 | 172  | 641608   | 87.145  | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.092 | 244  | 1634436  | 91.344  | ng    | 0.00     |

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

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

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