

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF081121\  
 Data File : BF125061.D  
 Acq On : 11 Aug 2021 17:54  
 Operator : JU/CG  
 Sample : M3316-02  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 GROUND-WATER

Quant Time: Aug 12 05:10:24 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF080421.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Aug 10 11:32:36 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.810  | 152  | 6616     | 20.000  | ng    | # 0.00   |
| 21) Naphthalene-d8          | 8.092  | 136  | 27124    | 20.000  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 9.845  | 164  | 15699    | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.333 | 188  | 28490    | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 13.969 | 240  | 10869    | 20.000  | ng    | # 0.00   |
| 86) Perylene-d12            | 15.421 | 264  | 10319    | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.434  | 112  | 26618    | 65.906  | ng    | 0.00     |
| 7) Phenol-d6                | 6.446  | 99   | 19587    | 38.461  | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.381  | 82   | 42645    | 82.247  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.639 | 330  | 19300    | 137.607 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.169  | 172  | 92716    | 86.417  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.922 | 244  | 67916    | 105.403 | ng    | 0.00     |

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

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

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