

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF112322\  
 Data File : BF131380.D  
 Acq On : 24 Nov 2022 01:24  
 Operator : CG\JU  
 Sample : N5641-11  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SOIL-PILE-11

Quant Time: Nov 24 05:59:44 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF112122.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Nov 24 05:53:32 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.957  | 152  | 103304   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.245  | 136  | 341400   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 10.010 | 164  | 159981   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.504 | 188  | 242504   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.162 | 240  | 210412   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.698 | 264  | 155861   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.569  | 112  | 636239   | 117.217 | ng    | 0.01     |
| 7) Phenol-d6                | 6.569  | 99   | 774749   | 111.872 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.522  | 82   | 489578   | 72.287  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.804 | 330  | 151153   | 112.304 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.327  | 172  | 798366   | 72.601  | ng    | 0.00     |
| 79) Terphenyl-d14           | 13.104 | 244  | 703791   | 54.697  | ng    | 0.00     |

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

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

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