

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF012422\  
 Data File : BF126642.D  
 Acq On : 24 Jan 2022 16:44  
 Operator : CG/JU  
 Sample : N1210-02  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 20220030-KEANSBURG-1-MODIFIED-ELUTRIATE

Quant Time: Jan 25 03:13:21 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF011822.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Jan 19 07:04:14 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.963  | 152  | 107859   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.245  | 136  | 417574   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 10.004 | 164  | 229190   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.498 | 188  | 399124   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.139 | 240  | 296244   | 20.000  | ng    | #-0.01   |
| 86) Perylene-d12            | 15.662 | 264  | 231336   | 20.000  | ng    | -0.01    |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.575  | 112  | 422309   | 51.523  | ng    | 0.00     |
| 7) Phenol-d6                | 6.569  | 99   | 397631   | 34.916  | ng    | -0.01    |
| 23) Nitrobenzene-d5         | 7.522  | 82   | 960648   | 90.046  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 10.792 | 330  | 232587   | 108.439 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.322  | 172  | 1227440  | 82.020  | ng    | 0.00     |
| 79) Terphenyl-d14           | 13.086 | 244  | 1208058  | 68.397  | ng    | 0.00     |

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

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

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