

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF093022\  
 Data File : BF130490.D  
 Acq On : 30 Sep 2022 18:20  
 Operator : CG\JU  
 Sample : N4896-01RE  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 CRUSHED-MASONRY-PILE-9RE

Quant Time: Oct 01 01:36:44 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF091422.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Sep 28 02:59:31 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.634  | 152  | 98392    | 20.000 | ng    | -0.01    |
| 21) Naphthalene-d8          | 7.916  | 136  | 374327   | 20.000 | ng    | -0.01    |
| 39) Acenaphthene-d10        | 9.669  | 164  | 187656   | 20.000 | ng    | -0.01    |
| 64) Phenanthrene-d10        | 11.145 | 188  | 286335   | 20.000 | ng    | -0.01    |
| 76) Chrysene-d12            | 13.780 | 240  | 184716   | 20.000 | ng    | -0.01    |
| 86) Perylene-d12            | 15.168 | 264  | 221984   | 20.000 | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.251  | 112  | 9417     | 1.660  | ng    | 0.00     |
| 7) Phenol-d6                | 6.275  | 99   | 152234   | 21.448 | ng    | -0.02    |
| 23) Nitrobenzene-d5         | 7.204  | 82   | 455179   | 62.123 | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 0.000  | 330  | 0        | 0.000  | ng    |          |
| 45) 2-Fluorobiphenyl        | 8.998  | 172  | 871522   | 67.629 | ng    | -0.01    |
| 79) Terphenyl-d14           | 12.733 | 244  | 698192   | 62.612 | ng    | -0.01    |

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

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

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