

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF021323\  
 Data File : BF132423.D  
 Acq On : 13 Feb 2023 22:23  
 Operator : CG\JU  
 Sample : 01328-06  
 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 COMP-2-BORING-1-7-14

Quant Time: Feb 14 01:41:22 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF021323.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Feb 14 00:51:00 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.031  | 152  | 162393   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.319  | 136  | 783904   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 10.083 | 164  | 402278   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.578 | 188  | 718507   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.236 | 240  | 489032   | 20.000  | ng    | # 0.00   |
| 86) Perylene-d12            | 15.801 | 264  | 449337   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.666  | 112  | 1228362  | 138.033 | ng    | 0.02     |
| 7) Phenol-d6                | 6.654  | 99   | 1225323  | 110.481 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.595  | 82   | 991615   | 75.463  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.878 | 330  | 515926   | 96.375  | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.395  | 172  | 1809413  | 65.116  | ng    | 0.00     |
| 79) Terphenyl-d14           | 13.172 | 244  | 1985422  | 68.281  | ng    | 0.00     |

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

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

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