

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF081823\  
 Data File : BF134852.D  
 Acq On : 18 Aug 2023 21:24  
 Operator : CG\JU  
 Sample : 04069-01  
 Misc :  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 CONC-WC-1

Quant Time: Aug 19 01:47:35 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF080123.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Aug 04 09:29:15 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min)    |
|-----------------------------|--------|------|----------|---------|-------|-------------|
| -----                       |        |      |          |         |       |             |
| Internal Standards          |        |      |          |         |       |             |
| 1) 1,4-Dichlorobenzene-d4   | 6.792  | 152  | 140399   | 20.000  | ng    | 0.00        |
| 21) Naphthalene-d8          | 8.069  | 136  | 542939   | 20.000  | ng    | #-0.01      |
| 39) Acenaphthene-d10        | 9.833  | 164  | 167769   | 20.000  | ng    | 0.00        |
| 64) Phenanthrene-d10        | 11.339 | 188  | 171235   | 20.000  | ng    | # 0.02      |
| 76) Chrysene-d12            | 13.980 | 240  | 245213   | 20.000  | ng    | # 0.00      |
| 86) Perylene-d12            | 15.439 | 264  | 186108   | 20.000  | ng    | 0.00        |
| System Monitoring Compounds |        |      |          |         |       |             |
| 5) 2-Fluorophenol           | 5.422  | 112  | 575079   | 62.264  | ng    | 0.00        |
| 7) Phenol-d6                | 6.428  | 99   | 981123   | 83.099  | ng    | -0.01       |
| 23) Nitrobenzene-d5         | 7.351  | 82   | 696067   | 80.118  | ng    | -0.02       |
| 42) 2,4,6-Tribromophenol    | 10.633 | 330  | 57500    | 33.621  | ng    | 0.00        |
| 45) 2-Fluorobiphenyl        | 9.151  | 172  | 1077039  | 106.741 | ng    | -0.01       |
| 79) Terphenyl-d14           | 12.933 | 244  | 718809   | 51.676  | ng    | 0.01        |
| Target Compounds            |        |      |          |         |       |             |
| 71) Phenanthrene            | 11.368 | 178  | 120123   | 13.207  | ng    | Qvalue # 60 |
| -----                       |        |      |          |         |       |             |

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

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