

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF031723\  
 Data File : BF132839.D  
 Acq On : 18 Mar 2023 00:07  
 Operator : CG\JU  
 Sample : 01954-08  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SV-TP4-0316

Quant Time: Mar 18 03:05:31 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF031623.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Mar 17 03:35:59 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.957  | 152  | 222992   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.240  | 136  | 832687   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.998  | 164  | 236507   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.492 | 188  | 749040   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.145 | 240  | 181578   | 20.000  | ng    | # 0.00   |
| 86) Perylene-d12            | 15.680 | 264  | 4540     | 20.000  | ng    | # 0.00   |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.616  | 112  | 1815393  | 137.372 | ng    | 0.03     |
| 7) Phenol-d6                | 6.604  | 99   | 1492234  | 86.996  | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.528  | 82   | 1534208  | 94.289  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.798 | 330  | 709048   | 282.022 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.322  | 172  | 2437915  | 169.328 | ng    | 0.00     |
| 79) Terphenyl-d14           | 13.086 | 244  | 2643329  | 262.864 | ng    | 0.00     |

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

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

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