

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF121222\  
 Data File : BF131595.D  
 Acq On : 12 Dec 2022 16:31  
 Operator : CG\JU  
 Sample : N5964-01  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SU-518-COMP-01

Quant Time: Dec 13 04:16:07 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF120822.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Dec 13 04:12:13 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.893  | 152  | 60451    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.181  | 136  | 218296   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.945  | 164  | 125559   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.439 | 188  | 241871   | 20.000  | ng    | -0.01    |
| 76) Chrysene-d12            | 14.098 | 240  | 129411   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.610 | 264  | 124830   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.510  | 112  | 368104   | 101.587 | ng    | 0.00     |
| 7) Phenol-d6                | 6.522  | 99   | 483270   | 101.776 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.457  | 82   | 328119   | 56.827  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 10.739 | 330  | 142004   | 101.795 | ng    | -0.01    |
| 45) 2-Fluorobiphenyl        | 9.263  | 172  | 549422   | 55.173  | ng    | 0.00     |
| 79) Terphenyl-d14           | 13.039 | 244  | 571623   | 60.576  | ng    | 0.00     |

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

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

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