

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF052822\  
 Data File : BF128542.D  
 Acq On : 28 May 2022 12:02  
 Operator : CG\JU  
 Sample : N2934-06  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 P025-SS001-2430-01

Quant Time: May 28 22:41:45 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF052622.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri May 27 01:51:59 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.840  | 152  | 192570   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.116  | 136  | 737376   | 20.000  | ng    | -0.01    |
| 39) Acenaphthene-d10        | 9.875  | 164  | 417956   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.363 | 188  | 772621   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.004 | 240  | 552534   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.468 | 264  | 426734   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.457  | 112  | 1174598  | 101.034 | ng    | 0.00     |
| 7) Phenol-d6                | 6.463  | 99   | 1544832  | 104.217 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.398  | 82   | 1042934  | 81.593  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 10.669 | 330  | 457805   | 115.077 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.198  | 172  | 1858103  | 72.227  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.957 | 244  | 1928272  | 67.423  | ng    | 0.00     |

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

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

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