

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF051822\  
 Data File : BF128312.D  
 Acq On : 18 May 2022 15:30  
 Operator : CG\JU  
 Sample : N2824-12  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 COM-2

Quant Time: May 18 16:50:52 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF050922.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed May 18 01:31:11 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min)      |
|-----------------------------|--------|------|----------|---------|-------|---------------|
| -----                       |        |      |          |         |       |               |
| Internal Standards          |        |      |          |         |       |               |
| 1) 1,4-Dichlorobenzene-d4   | 6.869  | 152  | 75208    | 20.000  | ng    | -0.01         |
| 21) Naphthalene-d8          | 8.151  | 136  | 298371   | 20.000  | ng    | -0.01         |
| 39) Acenaphthene-d10        | 9.910  | 164  | 167847   | 20.000  | ng    | -0.01         |
| 64) Phenanthrene-d10        | 11.404 | 188  | 298081   | 20.000  | ng    | -0.01         |
| 76) Chrysene-d12            | 14.051 | 240  | 201744   | 20.000  | ng    | -0.01         |
| 86) Perylene-d12            | 15.539 | 264  | 159624   | 20.000  | ng    | -0.02         |
| System Monitoring Compounds |        |      |          |         |       |               |
| 5) 2-Fluorophenol           | 5.510  | 112  | 571797   | 121.200 | ng    | 0.01          |
| 7) Phenol-d6                | 6.510  | 99   | 706676   | 113.249 | ng    | 0.00          |
| 23) Nitrobenzene-d5         | 7.440  | 82   | 589316   | 88.873  | ng    | -0.01         |
| 42) 2,4,6-Tribromophenol    | 10.704 | 330  | 232266   | 120.658 | ng    | -0.01         |
| 45) 2-Fluorobiphenyl        | 9.228  | 172  | 937151   | 87.362  | ng    | -0.01         |
| 79) Terphenyl-d14           | 12.992 | 244  | 1014429  | 94.352  | ng    | -0.01         |
| Target Compounds            |        |      |          |         |       |               |
| 71) Phenanthrene            | 11.428 | 178  | 54351    | 3.662   | ng    | Qvalue<br>100 |
| -----                       |        |      |          |         |       |               |

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

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