

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF052322\  
 Data File : BF128425.D  
 Acq On : 24 May 2022 03:32  
 Operator : CG\JU  
 Sample : N2875-16 5X  
 Misc :  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 P007-SS009-0006-01

Quant Time: May 24 04:22:36 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF052022.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Sat May 21 05:28:09 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.857  | 152  | 106101   | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.139  | 136  | 362425   | 20.000 | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.898  | 164  | 164604   | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.386 | 188  | 279470   | 20.000 | ng    | 0.00     |
| 76) Chrysene-d12            | 14.039 | 240  | 231686   | 20.000 | ng    | -0.01    |
| 86) Perylene-d12            | 15.527 | 264  | 171481   | 20.000 | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.481  | 112  | 91900    | 13.784 | ng    | 0.00     |
| 7) Phenol-d6                | 6.492  | 99   | 114469   | 13.038 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.422  | 82   | 66360    | 8.204  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 10.692 | 330  | 19216    | 11.320 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.210  | 172  | 118213   | 11.744 | ng    | -0.01    |
| 79) Terphenyl-d14           | 12.968 | 244  | 125391   | 10.448 | ng    | -0.01    |

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

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

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