

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF032422\  
 Data File : BF127498.D  
 Acq On : 24 Mar 2022 13:02  
 Operator : CG\JU  
 Sample : N2062-10  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 EB-2022-03-21

Quant Time: Mar 24 15:21:06 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF032122.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Mar 22 02:16:33 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.846  | 152  | 89580    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.128  | 136  | 321047   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.886  | 164  | 184426   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.375 | 188  | 322559   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.022 | 240  | 207258   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.504 | 264  | 197081   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.469  | 112  | 376472   | 67.083  | ng    | 0.00     |
| 7) Phenol-d6                | 6.481  | 99   | 323616   | 42.186  | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.416  | 82   | 703795   | 93.707  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.681 | 330  | 257311   | 130.973 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.204  | 172  | 1162941  | 92.078  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.963 | 244  | 1202558  | 94.984  | ng    | 0.00     |

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

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

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