

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF072023\  
 Data File : BF134376.D  
 Acq On : 20 Jul 2023 12:02  
 Operator : CG\JU  
 Sample : PB154198BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB154198BL

Quant Time: Jul 20 13:29:33 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF062623.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Jul 14 22:51:27 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.834  | 152  | 76156    | 20.000  | ng    | -0.01    |
| 21) Naphthalene-d8          | 8.116  | 136  | 235064   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.875  | 164  | 126285   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.369 | 188  | 261976   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.033 | 240  | 207405   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.533 | 264  | 209803   | 20.000  | ng    | -0.01    |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.481  | 112  | 605665   | 133.034 | ng    | 0.01     |
| 7) Phenol-d6                | 6.481  | 99   | 761010   | 135.921 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.404  | 82   | 450769   | 103.371 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.675 | 330  | 261497   | 165.153 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.198  | 172  | 891582   | 102.646 | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.969 | 244  | 1188038  | 92.189  | ng    | 0.00     |

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

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

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