

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF022323\  
 Data File : BF132519.D  
 Acq On : 23 Feb 2023 14:22  
 Operator : CG\JU  
 Sample : PB151028BL  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB151028BL

Quant Time: Feb 24 00:36:40 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF022223.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Feb 23 00:59:42 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.007  | 152  | 189419   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.295  | 136  | 702216   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 10.054 | 164  | 388042   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.548 | 188  | 765069   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.207 | 240  | 588176   | 20.000  | ng    | -0.01    |
| 86) Perylene-d12            | 15.759 | 264  | 488967   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.648  | 112  | 1450536  | 124.284 | ng    | 0.01     |
| 7) Phenol-d6                | 6.637  | 99   | 1900010  | 124.739 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.572  | 82   | 1288045  | 87.025  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.854 | 330  | 540400   | 128.953 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.372  | 172  | 2018047  | 79.187  | ng    | 0.00     |
| 79) Terphenyl-d14           | 13.148 | 244  | 2638905  | 75.857  | ng    | 0.00     |

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

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

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