

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP021623\  
 Data File : BP013831.D  
 Acq On : 16 Feb 2023 16:59  
 Operator : CG/JU  
 Sample : PB150892BL  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 PB150892BL

Quant Time: Feb 17 05:08:54 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP021623.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Feb 17 01:08:52 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.111  | 152  | 156579   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.940 | 136  | 580698   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.745 | 164  | 349353   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.498 | 188  | 781641   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.575 | 240  | 788537   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 24.127 | 264  | 911240   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.652  | 112  | 1449579  | 152.579 | ng    | 0.00     |
| 7) Phenol-d6                | 7.263  | 99   | 1792495  | 147.317 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.287  | 82   | 1079150  | 100.352 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.234 | 330  | 734021   | 144.338 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.375 | 172  | 2537009  | 95.384  | ng    | 0.00     |
| 79) Terphenyl-d14           | 20.028 | 244  | 4278699  | 104.398 | ng    | 0.00     |

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

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

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