

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF111622\  
 Data File : BF131255.D  
 Acq On : 16 Nov 2022 09:48  
 Operator : CG\JU  
 Sample : PB148901BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB148901BL

Quant Time: Nov 16 22:10:44 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF111522.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Nov 16 06:35:36 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.975  | 152  | 86946    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.263  | 136  | 323900   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 10.027 | 164  | 192011   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.527 | 188  | 355096   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 14.180 | 240  | 264018   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.727 | 264  | 194916   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.592  | 112  | 721169   | 146.874 | ng    | 0.02     |
| 7) Phenol-d6                | 6.586  | 99   | 907343   | 145.898 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.539  | 82   | 588734   | 102.024 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.821 | 330  | 256958   | 140.100 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.345  | 172  | 1154292  | 87.974  | ng    | 0.00     |
| 79) Terphenyl-d14           | 13.127 | 244  | 1449594  | 94.287  | ng    | 0.00     |

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

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

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