

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101023\  
 Data File : BF135680.D  
 Acq On : 10 Oct 2023 12:17  
 Operator : CG\JU  
 Sample : PB156142TB  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB156142TB

Quant Time: Oct 11 02:46:42 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF091123.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Oct 11 02:43:17 2023  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.740  | 152  | 95922    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.016  | 136  | 363923   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 9.775  | 164  | 192881   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.269 | 188  | 394655   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 13.933 | 240  | 248659   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.404 | 264  | 207084   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.381  | 112  | 835169   | 141.611 | ng    | 0.01     |
| 7) Phenol-d6                | 6.393  | 99   | 1034802  | 137.988 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.304  | 82   | 659543   | 98.462  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.575 | 330  | 289666   | 151.122 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.098  | 172  | 1208885  | 94.559  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.869 | 244  | 1374332  | 85.679  | ng    | 0.00     |

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

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

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