

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF070121\  
 Data File : BF124509.D  
 Acq On : 1 Jul 2021 13:43  
 Operator : JU/CG  
 Sample : PB137447TB  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB137447TB

Quant Time: Jul 01 14:17:43 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF062321.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Jul 01 14:15:13 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.681  | 152  | 43447    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 7.963  | 136  | 165848   | 20.000  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 9.716  | 164  | 92556    | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.204 | 188  | 175884   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 13.851 | 240  | 123831   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.280 | 264  | 92635    | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.316  | 112  | 353959   | 128.126 | ng    | 0.01     |
| 7) Phenol-d6                | 6.339  | 99   | 415192   | 115.104 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.251  | 82   | 308017   | 77.005  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.510 | 330  | 141429   | 127.897 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.045  | 172  | 642952   | 85.095  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.792 | 244  | 730127   | 88.312  | ng    | 0.00     |

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

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

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