

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF062121\  
 Data File : BF124388.D  
 Acq On : 21 Jun 2021 12:01  
 Operator : JU/CG  
 Sample : PB137220BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PB137220BL

Quant Time: Jun 21 13:38:29 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF060321.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Jun 15 20:09:26 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 6.786  | 152  | 37862    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 8.069  | 136  | 147894   | 20.000  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 9.822  | 164  | 84839    | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 11.310 | 188  | 155899   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 13.957 | 240  | 93087    | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 15.415 | 264  | 74122    | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.422  | 112  | 323817   | 128.745 | ng    | 0.00     |
| 7) Phenol-d6                | 6.439  | 99   | 397479   | 117.570 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 7.351  | 82   | 293255   | 78.643  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 10.622 | 330  | 131772   | 109.631 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 9.151  | 172  | 567711   | 84.205  | ng    | 0.00     |
| 79) Terphenyl-d14           | 12.904 | 244  | 590466   | 87.089  | ng    | 0.00     |

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

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

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