

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060921\  
 Data File : BP005811.D  
 Acq On : 09 Jun 2021 16:15  
 Operator : CG/JU  
 Sample : PB136947BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 PB136947BL

Quant Time: Jun 09 18:12:14 2021  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP060821.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Jun 09 11:54:55 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.952  | 152  | 85358    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.763 | 136  | 301819   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.581 | 164  | 182716   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.322 | 188  | 369267   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.392 | 240  | 406080   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 23.816 | 264  | 446955   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.517  | 112  | 730781   | 137.378 | ng    | 0.00     |
| 7) Phenol-d6                | 7.122  | 99   | 825987   | 119.798 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.116  | 82   | 517644   | 84.085  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.069 | 330  | 388866   | 124.019 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.210 | 172  | 1165551  | 83.765  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.869 | 244  | 1812283  | 83.567  | ng    | 0.00     |

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

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

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