

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP051921\  
 Data File : BP005607.D  
 Acq On : 19 May 2021 12:12  
 Operator : CG/JU  
 Sample : PB136517BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 PB136517BL

Quant Time: May 19 13:20:52 2021  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP050521.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon May 17 15:35:47 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.657  | 152  | 9653     | 20.00  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.451 | 136  | 28877    | 20.00  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 14.316 | 164  | 22138    | 20.00  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.075 | 188  | 53388    | 20.00  | ng    | # 0.00   |
| 76) Chrysene-d12            | 21.180 | 240  | 79969    | 20.00  | ng    | # 0.00   |
| 86) Perylene-d12            | 23.474 | 264  | 89380    | 20.00  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.252  | 112  | 83478    | 165.96 | ng    | 0.00     |
| 7) Phenol-d6                | 6.858  | 99   | 81618    | 130.13 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 8.828  | 82   | 66943    | 91.24  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 15.816 | 330  | 76762    | 122.19 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 12.928 | 172  | 162996   | 83.05  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.651 | 244  | 439573   | 93.35  | ng    | 0.00     |

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

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

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