

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP051521\  
 Data File : BP005572.D  
 Acq On : 15 May 2021 17:13  
 Operator : CG/JU  
 Sample : PB136344TB  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 PB136344TB

Quant Time: May 16 02:45:47 2021  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP050521.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri May 14 12:01:17 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.669  | 152  | 9436     | 20.00  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.463 | 136  | 30343    | 20.00  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 14.328 | 164  | 24344    | 20.00  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.086 | 188  | 56890    | 20.00  | ng    | # 0.00   |
| 76) Chrysene-d12            | 21.186 | 240  | 76810    | 20.00  | ng    | # 0.00   |
| 86) Perylene-d12            | 23.486 | 264  | 100885   | 20.00  | ng    | -0.01    |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.264  | 112  | 74108    | 150.72 | ng    | 0.00     |
| 7) Phenol-d6                | 6.869  | 99   | 81490    | 132.91 | ng    | -0.01    |
| 23) Nitrobenzene-d5         | 8.840  | 82   | 70253    | 91.13  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 15.828 | 330  | 76098    | 110.16 | ng    | -0.01    |
| 45) 2-Fluorobiphenyl        | 12.945 | 172  | 181702   | 84.20  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.657 | 244  | 422072   | 93.32  | ng    | 0.00     |

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

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

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