

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061421\  
 Data File : BP005918.D  
 Acq On : 14 Jun 2021 22:03  
 Operator : CG/JU  
 Sample : M2625-28RE  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 18-B12(166-170)RE

Quant Time: Jun 15 04:18:19 2021  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP060821.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Jun 14 13:15:03 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.934  | 152  | 65085    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.740 | 136  | 230021   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.563 | 164  | 138462   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.304 | 188  | 299999   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.374 | 240  | 374973   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 23.786 | 264  | 426978   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.505  | 112  | 539659   | 133.049 | ng    | 0.00     |
| 7) Phenol-d6                | 7.111  | 99   | 545576   | 103.776 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.099  | 82   | 449734   | 95.857  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.051 | 330  | 372341   | 156.702 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.193 | 172  | 977820   | 92.734  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.851 | 244  | 1231888  | 61.516  | ng    | 0.00     |

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

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

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