

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP070921\  
 Data File : BP006168.D  
 Acq On : 09 Jul 2021 19:15  
 Operator : CG/JU  
 Sample : M2990-07  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 TP-K

Quant Time: Jul 10 01:12:33 2021  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP070721.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Jul 07 16:08:13 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.869  | 152  | 183084   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.675 | 136  | 709335   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.510 | 164  | 414238   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.263 | 188  | 792823   | 20.000  | ng    | # 0.00   |
| 76) Chrysene-d12            | 21.345 | 240  | 761282   | 20.000  | ng    | # 0.00   |
| 86) Perylene-d12            | 23.739 | 264  | 873572   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.446  | 112  | 1179162  | 110.819 | ng    | -0.01    |
| 7) Phenol-d6                | 7.058  | 99   | 1394155  | 99.056  | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.040  | 82   | 777280   | 69.216  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.004 | 330  | 499032   | 101.187 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.134 | 172  | 1834676  | 65.476  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.816 | 244  | 2643612  | 68.466  | ng    | 0.00     |

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

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

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