

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092021\  
 Data File : BP007059.D  
 Acq On : 20 Sep 2021 20:17  
 Operator : CG/JU  
 Sample : M3786-07  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 N71924

Quant Time: Sep 21 01:03:17 2021  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP091621.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Sep 17 10:50:16 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.904  | 152  | 304649   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.704 | 136  | 1185662  | 20.000  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 14.534 | 164  | 734237   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.280 | 188  | 1532291  | 20.000  | ng    | # 0.00   |
| 76) Chrysene-d12            | 21.363 | 240  | 1527407  | 20.000  | ng    | # 0.00   |
| 86) Perylene-d12            | 23.774 | 264  | 1641242  | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.481  | 112  | 1026833  | 55.039  | ng    | 0.00     |
| 7) Phenol-d6                | 7.069  | 99   | 736551   | 28.925  | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.069  | 82   | 1837332  | 87.034  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.022 | 330  | 1215523  | 137.658 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.163 | 172  | 4534029  | 85.248  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.839 | 244  | 6867809  | 86.268  | ng    | 0.00     |

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

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

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