

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM072021\  
 Data File : BM031027.D  
 Acq On : 20 Jul 2021 17:46  
 Operator : CG/JU  
 Sample : PB137646BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 PB137646BL

Quant Time: Jul 20 18:21:46 2021  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM072021.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Jul 20 13:21:40 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.640  | 152  | 71281    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.404 | 136  | 306229   | 20.000  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 14.268 | 164  | 214802   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.015 | 188  | 484293   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.203 | 240  | 526169   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 23.380 | 264  | 550907   | 20.000  | ng    | -0.01    |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.263  | 112  | 647650   | 163.155 | ng    | 0.00     |
| 7) Phenol-d6                | 6.822  | 99   | 840963   | 146.143 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 8.786  | 82   | 541287   | 88.750  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 15.762 | 330  | 362656   | 129.292 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 12.886 | 172  | 1275484  | 88.598  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.656 | 244  | 2567015  | 96.527  | ng    | 0.00     |

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

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

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