

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM061021\  
 Data File : BM030386.D  
 Acq On : 11 Jun 2021 01:21  
 Operator : CG/JU  
 Sample : PB136995BL  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 PB136995BL

Quant Time: Jun 11 07:54:16 2021  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM060921.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Jun 08 20:51:08 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.839  | 152  | 129317   | 20.000  | ng    | -0.01    |
| 21) Naphthalene-d8          | 10.633 | 136  | 533366   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.462 | 164  | 391250   | 20.000  | ng    | -0.01    |
| 64) Phenanthrene-d10        | 17.198 | 188  | 1012925  | 20.000  | ng    | -0.01    |
| 76) Chrysene-d12            | 21.356 | 240  | 1280046  | 20.000  | ng    | -0.01    |
| 86) Perylene-d12            | 23.638 | 264  | 1240597  | 20.000  | ng    | -0.01    |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.428  | 112  | 1303127  | 163.817 | ng    | 0.00     |
| 7) Phenol-d6                | 7.004  | 99   | 1679231  | 145.823 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 8.998  | 82   | 869550   | 79.425  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol    | 15.951 | 330  | 125312   | 25.475  | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.086 | 172  | 2447580  | 82.267  | ng    | -0.01    |
| 79) Terphenyl-d14           | 19.815 | 244  | 5853605  | 80.321  | ng    | 0.00     |

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

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

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