

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM011222\  
 Data File : BM033700.D  
 Acq On : 12 Jan 2022 15:50  
 Operator : CG/JU  
 Sample : N1099-09  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 VACLOT-BACKFILL-TILCON-09

Quant Time: Jan 13 07:50:59 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM010522.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Jan 13 07:49:01 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-----------------------------|--------|------|----------|---------|-------|----------|
| -----                       |        |      |          |         |       |          |
| Internal Standards          |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.966  | 152  | 83083    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8          | 10.783 | 136  | 366613   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10        | 14.607 | 164  | 248318   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 17.342 | 188  | 563361   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12            | 21.500 | 240  | 723604   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12            | 23.924 | 264  | 912841   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |         |       |          |
| 5) 2-Fluorophenol           | 5.513  | 112  | 590969   | 112.182 | ng    | 0.00     |
| 7) Phenol-d6                | 7.119  | 99   | 810953   | 114.219 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 9.131  | 82   | 511881   | 64.482  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 16.089 | 330  | 382578   | 147.508 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 13.236 | 172  | 1180262  | 62.254  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.953 | 244  | 2125870  | 52.882  | ng    | 0.00     |

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

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

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