

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN092021\  
 Data File : BN016356.D  
 Acq On : 20 Sep 2021 23:35  
 Operator : CG/JU  
 Sample : PB139121BL  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 PB139121BL

Quant Time: Sep 21 01:34:47 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\8270-SIM-BN091721.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Sep 20 11:09:23 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units | Dev(Min) |
|-----------------------------|--------|------|----------|-------|-------|----------|
| Internal Standards          |        |      |          |       |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.679  | 152  | 25370    | 0.400 | ng    | 0.00     |
| 7) Naphthalene-d8           | 10.445 | 136  | 115459   | 0.400 | ng    | #-0.01   |
| 13) Acenaphthene-d10        | 14.307 | 164  | 60448    | 0.400 | ng    | 0.00     |
| 19) Phenanthrene-d10        | 17.054 | 188  | 120231   | 0.400 | ng    | 0.00     |
| 29) Chrysene-d12            | 21.259 | 240  | 118859   | 0.400 | ng    | # 0.00   |
| 35) Perylene-d12            | 23.519 | 264  | 100790   | 0.400 | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |       |       |          |
| 4) 2-Fluorophenol           | 5.328  | 112  | 24466    | 0.384 | ng    | 0.02     |
| 5) Phenol-d6                | 6.868  | 99   | 27406    | 0.355 | ng    | 0.00     |
| 8) Nitrobenzene-d5          | 8.822  | 82   | 28943    | 0.422 | ng    | 0.00     |
| 11) 2-Methylnaphthalene-d10 | 12.043 | 152  | 72711    | 0.400 | ng    | 0.00     |
| 14) 2,4,6-Tribromophenol    | 15.804 | 330  | 5187     | 0.292 | ng    | 0.00     |
| 15) 2-Fluorobiphenyl        | 12.937 | 172  | 100262   | 0.424 | ng    | 0.00     |
| 27) Fluoranthene-d10        | 19.093 | 212  | 142916   | 0.411 | ng    | 0.00     |
| 31) Terphenyl-d14           | 19.703 | 244  | 126554   | 0.460 | ng    | 0.00     |

Target Compounds Qvalue

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

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