

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN101221\  
 Data File : BN016894.D  
 Acq On : 11 Oct 2021 23:23  
 Operator : CG/JU  
 Sample : PB139657BL  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 PB139657BL

Quant Time: Oct 12 02:35:58 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\8270-SIM-BN101221.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Oct 12 02:31:22 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc | Units | Dev(Min) |
|-----------------------------|--------|------|----------|------|-------|----------|
| -----                       |        |      |          |      |       |          |
| Internal Standards          |        |      |          |      |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.698  | 152  | 23103    | 0.40 | ng    | 0.00     |
| 7) Naphthalene-d8           | 10.457 | 136  | 103199   | 0.40 | ng    | 0.00     |
| 13) Acenaphthene-d10        | 14.294 | 164  | 52863    | 0.40 | ng    | 0.00     |
| 19) Phenanthrene-d10        | 17.042 | 188  | 97314    | 0.40 | ng    | 0.00     |
| 29) Chrysene-d12            | 21.227 | 240  | 89703    | 0.40 | ng    | #-0.01   |
| 35) Perylene-d12            | 23.474 | 264  | 92144    | 0.40 | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |      |       |          |
| 4) 2-Fluorophenol           | 5.337  | 112  | 21750    | 0.37 | ng    | 0.04     |
| 5) Phenol-d6                | 6.886  | 99   | 20372    | 0.31 | ng    | 0.02     |
| 8) Nitrobenzene-d5          | 8.835  | 82   | 23969    | 0.34 | ng    | 0.00     |
| 11) 2-Methylnaphthalene-d10 | 12.043 | 152  | 60020    | 0.40 | ng    | 0.00     |
| 14) 2,4,6-Tribromophenol    | 15.791 | 330  | 4111     | 0.16 | ng    | 0.00     |
| 15) 2-Fluorobiphenyl        | 12.924 | 172  | 75602    | 0.37 | ng    | 0.00     |
| 27) Fluoranthene-d10        | 19.073 | 212  | 102655   | 0.40 | ng    | 0.00     |
| 31) Terphenyl-d14           | 19.678 | 244  | 85903    | 0.42 | ng    | 0.00     |

Target Compounds Qvalue

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

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