

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN041621\  
 Data File : BN014313.D  
 Acq On : 16 Apr 2021 17:12  
 Operator : CG/JU  
 Sample : PB135776BL  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 PB135776BL

Quant Time: Apr 19 09:22:37 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\8270-SIM-BN041621.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Apr 16 16:29:41 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc | Units | Dev(Min) |
|-----------------------------|--------|------|----------|------|-------|----------|
| -----                       |        |      |          |      |       |          |
| Internal Standards          |        |      |          |      |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.700  | 152  | 7024     | 0.40 | ng    | 0.00     |
| 7) Naphthalene-d8           | 10.480 | 136  | 25456    | 0.40 | ng    | 0.00     |
| 13) Acenaphthene-d10        | 14.334 | 164  | 13141    | 0.40 | ng    | 0.00     |
| 19) Phenanthrene-d10        | 17.079 | 188  | 26747    | 0.40 | ng    | # 0.01   |
| 29) Chrysene-d12            | 21.269 | 240  | 24077    | 0.40 | ng    | # 0.00   |
| 36) Perylene-d12            | 23.504 | 264  | 18120    | 0.40 | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |      |       |          |
| 4) 2-Fluorophenol           | 5.300  | 112  | 6801     | 0.44 | ng    | 0.00     |
| 5) Phenol-d6                | 6.871  | 99   | 7898     | 0.40 | ng    | 0.00     |
| 8) Nitrobenzene-d5          | 8.845  | 82   | 6517     | 0.41 | ng    | 0.00     |
| 11) 2-Methylnaphthalene-d10 | 12.066 | 152  | 16412    | 0.42 | ng    | 0.00     |
| 14) 2,4,6-Tribromophenol    | 15.818 | 330  | 1567     | 0.33 | ng    | 0.00     |
| 15) 2-Fluorobiphenyl        | 12.951 | 172  | 22427    | 0.43 | ng    | 0.00     |
| 27) Fluoranthene-d10        | 19.111 | 212  | 28664    | 0.39 | ng    | 0.00     |
| 31) Terphenyl-d14           | 19.717 | 244  | 26030    | 0.52 | ng    | 0.00     |

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

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

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