

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN120821\  
 Data File : BN017828.D  
 Acq On : 09 Dec 2021 09:58  
 Operator : CG/JU  
 Sample : M4922-16  
 Misc :  
 ALS Vial : 33 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 RE139D1-20211130

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 12/10/2021  
 Supervised By :mohammad ahmed 12/15/2021

Quant Time: Dec 10 01:53:01 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\8270-SIM-BN120721.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Dec 08 01:18:09 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc  | Units | Dev(Min) |
|-------------------------------|--------|------|----------|-------|-------|----------|
| -----                         |        |      |          |       |       |          |
| Internal Standards            |        |      |          |       |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.715  | 152  | 20192    | 0.400 | ng    | 0.00     |
| 7) Naphthalene-d8             | 10.494 | 136  | 79085    | 0.400 | ng    | 0.00     |
| 13) Acenaphthene-d10          | 14.347 | 164  | 57271    | 0.400 | ng    | 0.00     |
| 19) Phenanthrene-d10          | 17.092 | 188  | 122846   | 0.400 | ng    | 0.00     |
| 29) Chrysene-d12              | 21.293 | 240  | 108235   | 0.400 | ng    | # 0.01   |
| 35) Perylene-d12              | 23.589 | 264  | 79338    | 0.400 | ng    | 0.00     |
| System Monitoring Compounds   |        |      |          |       |       |          |
| 4) 2-Fluorophenol             | 5.336  | 112  | 6513     | 0.134 | ng    | 0.00     |
| 5) Phenol-d6                  | 6.903  | 99   | 3802     | 0.082 | ng    | 0.00     |
| 8) Nitrobenzene-d5            | 8.859  | 82   | 18265    | 0.347 | ng    | 0.00     |
| 11) 2-Methylnaphthalene-d10   | 12.090 | 152  | 43125    | 0.273 | ng    | 0.00     |
| 14) 2,4,6-Tribromophenol      | 15.844 | 330  | 12652    | 0.423 | ng    | 0.00     |
| 15) 2-Fluorobiphenyl          | 12.964 | 172  | 75272    | 0.359 | ng    | 0.00     |
| 27) Fluoranthene-d10          | 19.129 | 212  | 155695   | 0.379 | ng    | 0.00     |
| 31) Terphenyl-d14             | 19.730 | 244  | 127879   | 0.550 | ng    | 0.00     |
| Target Compounds              |        |      |          |       |       |          |
| 2) 1,4-Dioxane                | 3.270  | 88   | 8272m    | 1.113 | ng    | Qvalue   |
| 25) Phenanthrene              | 17.129 | 178  | 9046     | 0.028 | ng    | 98       |
| 34) Bis(2-ethylhexyl)phtha... | 21.219 | 149  | 18855    | 0.125 | ng    | 99       |
| -----                         |        |      |          |       |       |          |

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

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