

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100121\  
 Data File : BN016683.D  
 Acq On : 02 Oct 2021 22:33  
 Operator : CG/JU  
 Sample : M3892-12  
 Misc :  
 ALS Vial : 51 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 S-825-N-SO-1-1.5-092221

Manual Integrations  
 APPROVED

mohammad  
 10/4/2021 11:48:56 AM

Quant Time: Oct 04 03:35:49 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\8270-SIM-BN100121.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Oct 01 13:35:26 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc | Units | Dev(Min) |
|-------------------------------|--------|------|----------|------|-------|----------|
| -----                         |        |      |          |      |       |          |
| Internal Standards            |        |      |          |      |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.642  | 152  | 31366    | 0.40 | ng    | 0.00     |
| 7) Naphthalene-d8             | 10.407 | 136  | 142266   | 0.40 | ng    | # 0.00   |
| 13) Acenaphthene-d10          | 14.269 | 164  | 78881    | 0.40 | ng    | 0.00     |
| 19) Phenanthrene-d10          | 17.018 | 188  | 153963   | 0.40 | ng    | #-0.01   |
| 29) Chrysene-d12              | 21.227 | 240  | 155667   | 0.40 | ng    | #-0.01   |
| 35) Perylene-d12              | 23.485 | 264  | 163583   | 0.40 | ng    | 0.00     |
| System Monitoring Compounds   |        |      |          |      |       |          |
| 4) 2-Fluorophenol             | 5.273  | 112  | 20804    | 0.26 | ng    | 0.02     |
| 5) Phenol-d6                  | 6.831  | 99   | 21281    | 0.24 | ng    | 0.00     |
| 8) Nitrobenzene-d5            | 8.784  | 82   | 22342    | 0.24 | ng    | 0.00     |
| 11) 2-Methylnaphthalene-d10   | 12.003 | 152  | 55045    | 0.26 | ng    | 0.00     |
| 14) 2,4,6-Tribromophenol      | 15.766 | 330  | 10787    | 0.30 | ng    | 0.00     |
| 15) 2-Fluorobiphenyl          | 12.886 | 172  | 67173    | 0.21 | ng    | -0.01    |
| 27) Fluoranthene-d10          | 19.063 | 212  | 92998    | 0.23 | ng    | 0.00     |
| 31) Terphenyl-d14             | 19.679 | 244  | 67395    | 0.20 | ng    | 0.00     |
| Target Compounds              |        |      |          |      |       |          |
|                               |        |      |          |      |       | Qvalue   |
| 25) Phenanthrene              | 17.066 | 178  | 16064m   | 0.03 | ng    |          |
| 26) Anthracene                | 17.151 | 178  | 8042     | 0.02 | ng    | # 56     |
| 28) Fluoranthene              | 19.093 | 202  | 52420    | 0.10 | ng    | 98       |
| 30) Pyrene                    | 19.455 | 202  | 43999    | 0.09 | ng    | # 96     |
| 32) Benzo(a)anthracene        | 21.217 | 228  | 35414    | 0.07 | ng    | 97       |
| 33) Chrysene                  | 21.270 | 228  | 37949    | 0.08 | ng    | 98       |
| 34) Bis(2-ethylhexyl)phtha... | 21.164 | 149  | 29475    | 0.15 | ng    | # 99     |
| 36) Indeno(1,2,3-cd)pyrene    | 25.764 | 276  | 27610    | 0.05 | ng    | # 85     |
| 37) Benzo(b)fluoranthene      | 22.807 | 252  | 51145    | 0.10 | ng    | # 87     |
| 38) Benzo(k)fluoranthene      | 22.848 | 252  | 30618m   | 0.06 | ng    |          |
| 39) Benzo(a)pyrene            | 23.385 | 252  | 29865    | 0.06 | ng    | # 79     |
| 40) Dibenzo(a,h)anthracene    | 25.785 | 278  | 13152    | 0.03 | ng    | # 49     |
| 41) Benzo(g,h,i)perylene      | 26.459 | 276  | 22646    | 0.04 | ng    | # 87     |
| -----                         |        |      |          |      |       |          |

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

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