

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM070821\  
 Data File : BM030886.D  
 Acq On : 08 Jul 2021 15:54  
 Operator : CG/JU  
 Sample : M2825-07  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BGDR1

Manual Integrations  
 APPROVED

mohammad  
 7/9/2021 2:48:03 PM

Quant Time: Jul 08 16:33:00 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-SIM-BM062821.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Jul 08 10:44:22 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-----------------------------|--------|------|----------|-------|--------|----------|
| -----                       |        |      |          |       |        |          |
| Internal Standards          |        |      |          |       |        |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.767  | 152  | 1945     | 0.400 | ng/ul  | 0.00     |
| 4) Naphthalene-d8           | 10.550 | 136  | 7587     | 0.400 | ng/ul  | 0.00     |
| 9) Acenaphthene-d10         | 14.394 | 164  | 4120     | 0.400 | ng/ul  | -0.01    |
| 13) Phenanthrene-d10        | 17.131 | 188  | 6789     | 0.400 | ng/ul  | -0.02    |
| 17) Chrysene-d12            | 21.304 | 240  | 5362     | 0.400 | ng/ul  | -0.01    |
| 23) Perylene-d12            | 23.549 | 264  | 5388     | 0.400 | ng/ul  | #-0.02   |
|                             |        |      |          |       |        |          |
| System Monitoring Compounds |        |      |          |       |        |          |
| 3) 1,4-Dioxane-d8           | 3.265  | 96   | 2721     | 1.284 | ng/ul  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 12.151 | 152  | 2176     | 0.183 | ng/ul  | 0.00     |
| 18) Fluoranthene-d10        | 19.153 | 212  | 3231     | 0.210 | ng/ul  | -0.02    |
|                             |        |      |          |       |        |          |
| Target Compounds            |        |      |          |       |        |          |
|                             |        |      |          |       | Qvalue |          |
| 5) Naphthalene              | 10.599 | 128  | 5451     | 0.252 | ng/ul  | 96       |
| 7) 2-Methylnaphthalene      | 12.227 | 142  | 997      | 0.069 | ng/ul  | 98       |
| 8) 1-Methylnaphthalene      | 12.439 | 142  | 468      | 0.033 | ng/ul  | 94       |
| 10) Acenaphthylene          | 14.114 | 152  | 2499     | 0.136 | ng/ul# | 87       |
| 11) Acenaphthene            | 14.455 | 153  | 19074    | 1.280 | ng/ul  | 99       |
| 12) Fluorene                | 15.445 | 166  | 10785    | 0.656 | ng/ul  | 98       |
| 15) Phenanthrene            | 17.173 | 178  | 99020    | 4.339 | ng/ul  | 100      |
| 16) Anthracene              | 17.263 | 178  | 34545    | 1.728 | ng/ul  | 100      |
| 19) Fluoranthene            | 19.184 | 202  | 110419   | 4.666 | ng/ul  | 99       |
| 20) Pyrene                  | 19.546 | 202  | 85894    | 3.609 | ng/ul  | 99       |
| 21) Benzo(a)anthracene      | 21.287 | 228  | 45350    | 2.122 | ng/ul  | 97       |
| 22) Chrysene                | 21.340 | 228  | 40180    | 1.678 | ng/ul  | 98       |
| 24) Benzo(b)fluoranthene    | 22.873 | 252  | 48755m   | 2.043 | ng/ul  |          |
| 25) Benzo(k)fluoranthene    | 22.912 | 252  | 21254m   | 0.864 | ng/ul  |          |
| 26) Benzo(a)pyrene          | 23.447 | 252  | 31303    | 1.542 | ng/ul# | 93       |
| 27) Indeno(1,2,3-cd)pyrene  | 25.817 | 276  | 21188    | 0.801 | ng/ul# | 89       |
| 28) Dibenzo(a,h)anthracene  | 25.817 | 278  | 5720     | 0.270 | ng/ul# | 86       |
| 29) Benzo(g,h,i)perylene    | 26.512 | 276  | 2264     | 0.099 | ng/ul# | 81       |
| -----                       |        |      |          |       |        |          |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM070821\  
Data File : BM030886.D  
Acq On : 08 Jul 2021 15:54  
Operator : CG/JU  
Sample : M2825-07  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BGDR1

Manual Integrations  
APPROVED

mohammad  
7/9/2021 2:48:03 PM

Quant Time: Jul 08 16:33:00 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-SIM-BM062821.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Thu Jul 08 10:44:22 2021  
Response via : Initial Calibration

