

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_M\DATA\BM080319\  
 Data File : BM021795.D  
 Acq On : 04 Aug 2019 04:28  
 Operator : HP/JU  
 Sample : K3850-21  
 Misc :  
 ALS Vial : 38 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BENS7

Manual Integrations  
 APPROVED

mohammad  
 8/5/2019 10:14:37 PM

Quant Time: Aug 04 05:20:58 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-SIM-BM080319.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Sun Aug 04 04:24:24 2019  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units  | Dev(Min) |
|-----------------------------|-------|------|----------|--------|--------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 7.64  | 152  | 1118     | 0.40   | ng/ul  | -0.01    |
| 2) Naphthalene-d8           | 10.42 | 136  | 4369     | 0.40   | ng/ul  | -0.03    |
| 6) Acenaphthene-d10         | 14.29 | 164  | 2528     | 0.40   | ng/ul  | -0.01    |
| 10) Phenanthrene-d10        | 17.04 | 188  | 6212m    | 0.40   | ng/ul  | -0.03    |
| 16) Chrysene-d12            | 21.24 | 240  | 6456m    | 0.40   | ng/ul  | -0.01    |
| 20) Perylene-d12            | 23.46 | 264  | 6868m    | 0.40   | ng/ul  | -0.02    |
| System Monitoring Compounds |       |      |          |        |        |          |
| 4) 2-Methylnaphthalene-d10  | 12.03 | 152  | 1282     | 0.18   | ng/ul  | -0.02    |
| 14) Fluoranthene-d10        | 19.08 | 212  | 4434m    | 0.21   | ng/ul  | -0.02    |
| Target Compounds            |       |      |          | Qvalue |        |          |
| 5) 2-Methylnaphthalene      | 12.11 | 142  | 229      | 0.028  | ng/ul  | 100      |
| 7) Acenaphthylene           | 14.01 | 152  | 626      | 0.053  | ng/ul# | 82       |
| 12) Phenanthrene            | 17.08 | 178  | 4057     | 0.224  | ng/ul  | 99       |
| 13) Anthracene              | 17.18 | 178  | 828m     | 0.054  | ng/ul  |          |
| 15) Fluoranthene            | 19.11 | 202  | 10852m   | 0.448  | ng/ul  |          |
| 17) Pyrene                  | 19.47 | 202  | 9961     | 0.375  | ng/ul  | 99       |
| 18) Benzo(a)anthracene      | 21.23 | 228  | 6002     | 0.260  | ng/ul  | 94       |
| 19) Chrysene                | 21.28 | 228  | 5464     | 0.184  | ng/ul# | 95       |
| 21) Benzo(b)fluoranthene    | 22.79 | 252  | 9532m    | 0.412  | ng/ul  |          |
| 22) Benzo(k)fluoranthene    | 22.83 | 252  | 3384m    | 0.129  | ng/ul  |          |
| 23) Benzo(a)pyrene          | 23.36 | 252  | 5685m    | 0.241  | ng/ul  |          |
| 24) Indeno(1,2,3-cd)pyrene  | 25.68 | 276  | 4872m    | 0.173  | ng/ul  |          |
| 25) Dibenzo(a,h)anthracene  | 25.68 | 278  | 1107m    | 0.049  | ng/ul  |          |
| 26) Benzo(g,h,i)perylene    | 26.36 | 276  | 5238m    | 0.207  | ng/ul  |          |

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

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_M\DATA\BM080319\  
Data File : BM021795.D  
Acq On : 04 Aug 2019 04:28  
Operator : HP/JU  
Sample : K3850-21  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BENS7

Manual Integrations  
APPROVED

mohammad  
8/5/2019 10:14:37 PM

Quant Time: Aug 04 05:20:58 2019  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-SIM-BM080319.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Sun Aug 04 04:24:24 2019  
Response via : Initial Calibration

