

Data Path : Z:\HPCHEM1\BNA M\DATA\BM112717\  
 Data File : BM013125.D  
 Acq On : 27 Nov 2017 23:45  
 Operator : SJ/JU  
 Sample : I6496-01  
 Misc :  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 C0AA0

Manual Integrations  
 APPROVED

Sohil  
 11/28/2017 2:54:09 PM

Quant Time: Nov 28 03:30:03 2017  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM111617.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Nov 28 01:49:10 2017  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc  | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|-------|--------|----------|
| 1) 1,4-Dichlorobenzene-d4      | 7.70  | 152  | 187918   | 20.00 | ng/ul  | 0.00     |
| 18) Naphthalene-d8             | 10.49 | 136  | 839121   | 20.00 | ng/ul  | 0.00     |
| 35) Acenaphthene-d10           | 14.35 | 164  | 537988   | 20.00 | ng/ul  | 0.00     |
| 61) Phenanthrene-d10           | 17.09 | 188  | 1274991  | 20.00 | ng/ul  | 0.00     |
| 75) Chrysene-d12               | 21.27 | 240  | 1533138  | 20.00 | ng/ul  | 0.00     |
| 83) Perylene-d12               | 23.50 | 264  | 1476547  | 20.00 | ng/ul  | 0.00     |
| System Monitoring Compounds    |       |      |          |       |        |          |
| 3) 1,4-Dioxane-d8              | 3.25  | 96   | 16453    | 4.06  | ng/uL  | 0.00     |
| 5) Phenol-d5                   | 6.89  | 99   | 539629   | 33.15 | ng/ul  | 0.00     |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.05  | 67   | 280102   | 30.33 | ng/ul  | 0.00     |
| 9) 2-Chlorophenol-d4           | 7.25  | 132  | 413939   | 31.82 | ng/ul  | 0.00     |
| 13) 4-Methylphenol-d8          | 8.42  | 113  | 471321   | 35.35 | ng/ul  | -0.01    |
| 19) Nitrobenzene-d5            | 8.86  | 128  | 208738   | 34.90 | ng/ul  | 0.00     |
| 22) 2-Nitrophenol-d4           | 9.57  | 143  | 234867   | 41.29 | ng/ul  | 0.00     |
| 26) 2,4-Dichlorophenol-d3      | 10.12 | 165  | 491717   | 34.05 | ng/ul  | -0.01    |
| 29) 4-Chloroaniline-d4         | 10.63 | 131  | 469822   | 31.11 | ng/ul  | 0.00     |
| 43) Dimethylphthalate-d6       | 13.76 | 166  | 1844389  | 39.08 | ng/ul  | 0.00     |
| 46) Acenaphthylene-d8          | 14.04 | 160  | 2075380  | 34.52 | ng/ul  | 0.00     |
| 51) 4-Nitrophenol-d4           | 14.57 | 143  | 295974   | 41.35 | ng/ul  | -0.01    |
| 57) Fluorene-d10               | 15.34 | 176  | 1476802  | 36.83 | ng/ul  | 0.00     |
| 62) 4,6-Dinitro-2-methylphenol | 15.46 | 200  | 275639   | 49.86 | ng/ul  | 0.00     |
| 70) Anthracene-d10             | 17.19 | 188  | 2481786  | 38.02 | ng/ul  | 0.00     |
| 76) Pyrene-d10                 | 19.49 | 212  | 3052037  | 38.97 | ng/ul  | 0.00     |
| 87) Benzo(a)pyrene-d12         | 23.37 | 264  | 3150483  | 42.08 | ng/ul  | 0.00     |
| Target Compounds               |       |      |          |       |        |          |
| 6) Phenol                      | 6.92  | 94   | 76983    | 4.792 | ng/ul  | 90       |
| 44) Dimethylphthalate          | 13.81 | 163  | 236644   | 5.244 | ng/ul  | 99       |
| 69) Phenanthrene               | 17.13 | 178  | 108593   | 1.514 | ng/ul  | 99       |
| 74) Fluoranthene               | 19.15 | 202  | 403745   | 4.658 | ng/ul# | 94       |
| 77) Pyrene                     | 19.52 | 202  | 378243   | 3.962 | ng/ul# | 93       |
| 80) Benzo(a)anthracene         | 21.26 | 228  | 252944   | 2.578 | ng/ul  | 99       |
| 82) Chrysene                   | 21.31 | 228  | 225081   | 2.474 | ng/ul  | 99       |
| 85) Benzo(b)fluoranthene       | 22.83 | 252  | 316098   | 3.313 | ng/ul# | 92       |
| 86) Benzo(k)fluoranthene       | 22.87 | 252  | 117677m  | 1.340 | ng/ul  |          |
| 88) Benzo(a)pyrene             | 23.41 | 252  | 233705   | 2.630 | ng/ul# | 96       |

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

Data Path : Z:\HPCHEM1\BNA M\DATA\BM112717\  
Data File : BM013125.D  
Acq On : 27 Nov 2017 23:45  
Operator : SJ/JU  
Sample : I6496-01  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C0AA0

Manual Integrations  
APPROVED

Sohil  
11/28/2017 2:54:09 PM

Quant Time: Nov 28 03:30:03 2017  
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM111617.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Tue Nov 28 01:49:10 2017  
Response via : Initial Calibration

