

Data Path : Z:\HPCHEM1\BNA M\DATA\BM112816\  
 Data File : BM008090.D  
 Acq On : 28 Nov 2016 17:22  
 Operator : UM/SJ  
 Sample : H5701-07DL 2X  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 A4WD2DL

Manual Integrations  
 APPROVED

UMANGI  
 11/29/2016 2:22:38 PM

Quant Time: Nov 28 17:56:26 2016  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM112116.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Nov 25 01:27:01 2016  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units  | Dev(Min) |
|-----------------------------|-------|------|----------|--------|--------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 7.71  | 152  | 652      | 0.40   | ng/ul  | -0.02    |
| 2) Naphthalene-d8           | 10.50 | 136  | 2277     | 0.40   | ng/ul  | 0.00     |
| 6) Acenaphthene-d10         | 14.35 | 164  | 1364     | 0.40   | ng/ul  | 0.00     |
| 10) Phenanthrene-d10        | 17.09 | 188  | 2540     | 0.40   | ng/ul  | -0.02    |
| 16) Chrysene-d12            | 21.30 | 240  | 2338     | 0.40   | ng/ul  | 0.00     |
| 20) Perylene-d12            | 23.54 | 264  | 2358     | 0.40   | ng/ul  | 0.00     |
| System Monitoring Compounds |       |      |          |        |        |          |
| 4) 2-Methylnaphthalene-d10  | 12.10 | 152  | 257      | 0.08   | ng/ul  | 0.00     |
| 14) Fluoranthene-d10        | 19.14 | 212  | 788      | 0.11   | ng/ul  | 0.00     |
| Target Compounds            |       |      |          | Ovalue |        |          |
| 7) Acenaphthylene           | 14.06 | 152  | 841      | 0.14   | ng/ul# | 84       |
| 8) Acenaphthene             | 14.40 | 153  | 126      | 0.03   | ng/ul# | 85       |
| 9) Fluorene                 | 15.41 | 166  | 199      | 0.04   | ng/ul# | 81       |
| 12) Phenanthrene            | 17.14 | 178  | 3136     | 0.40   | ng/ul# | 93       |
| 13) Anthracene              | 17.23 | 178  | 1310m    | 0.18   | ng/ul  |          |
| 15) Fluoranthene            | 19.16 | 202  | 9297     | 0.91   | ng/ul  | 97       |
| 17) Pyrene                  | 19.52 | 202  | 7956     | 0.98   | ng/ul  | 99       |
| 18) Benzo(a)anthracene      | 21.28 | 228  | 5492     | 0.77   | ng/ul# | 80       |
| 19) Chrysene                | 21.33 | 228  | 5362     | 0.60   | ng/ul# | 78       |
| 21) Benzo(b)fluoranthene    | 22.86 | 252  | 8301m    | 0.88   | ng/ul  |          |
| 22) Benzo(k)fluoranthene    | 22.90 | 252  | 2565m    | 0.26   | ng/ul  |          |
| 23) Benzo(a)pyrene          | 23.43 | 252  | 5716     | 0.65   | ng/ul# | 75       |
| 24) Indeno(1,2,3-cd)pyrene  | 25.79 | 276  | 3876     | 0.36   | ng/ul# | 90       |
| 25) Dibenzo(a,h)anthracene  | 25.80 | 278  | 908      | 0.10   | ng/ul# | 1        |
| 26) Benzo(g,h,i)perylene    | 26.49 | 276  | 3253     | 0.34   | ng/ul# | 64       |

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

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