

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM050515\  
 Data File : BM001217.D  
 Acq On : 05 May 2015 20:40  
 Operator : TP/IZ  
 Sample : PB83182BS  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 PB83182BS

Quant Time: May 06 13:10:16 2015  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SIMPAH-BM050515.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed May 06 12:46:35 2015  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc | Units | Dev(Min) |
|---------------------------|-------|------|----------|------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.52  | 152  | 13881    | 5.00 | ng    | 0.00     |
| 6) Naphthalene-d8         | 10.30 | 136  | 57293    | 5.00 | ng    | 0.00     |
| 12) Acenaphthene-d10      | 14.19 | 164  | 34700    | 5.00 | ng    | 0.00     |
| 18) Phenanthrene-d10      | 16.95 | 188  | 53236    | 5.00 | ng    | 0.00     |
| 24) Chrysene-d12          | 21.14 | 240  | 79141    | 5.00 | ng    | 0.00     |
| 30) Perylene-d12          | 23.29 | 264  | 68867    | 5.00 | ng    | 0.00     |

| System Monitoring Compounds | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|-----------------------------|-------|------|----------|-------|-------|----------|
| 4) 2-Fluorophenol           | 5.11  | 112  | 34073    | 11.74 | ng    | 0.00     |
| 5) Phenol-d6                | 6.71  | 99   | 47168    | 12.78 | ng    | -0.02    |
| 7) Nitrobenzene-d5          | 8.68  | 82   | 29945    | 9.24  | ng    | 0.00     |
| 13) 2,4,6-Tribromophenol    | 15.69 | 330  | 29748    | 11.89 | ng    | 0.00     |
| 14) 2-Fluorobiphenyl        | 12.80 | 172  | 89162    | 8.40  | ng    | -0.01    |
| 26) Terphenyl-d14           | 19.59 | 244  | 85976    | 7.59  | ng    | 0.00     |

| Target Compounds           | R.T.  | QIon | Response | Conc  | Units | Qvalue |
|----------------------------|-------|------|----------|-------|-------|--------|
| 2) 1,4-Dioxane             | 3.01  | 88   | 8922     | 8.31  | ng    | # 71   |
| 3) n-Nitrosodimethylamine  | 3.32  | 42   | 14191    | 8.43  | ng    | # 86   |
| 8) Nitrobenzene            | 8.72  | 77   | 34235    | 8.07  | ng    | 98     |
| 9) Naphthalene             | 10.35 | 128  | 102050   | 8.18  | ng    | 99     |
| 10) Hexachlorobutadiene    | 10.64 | 225  | 19094    | 7.91  | ng    | 98     |
| 11) 2-Methylnaphthalene    | 11.98 | 142  | 71471    | 8.16  | ng    | 97     |
| 15) Acenaphthylene         | 13.90 | 152  | 134233   | 8.98  | ng    | 100    |
| 16) Acenaphthene           | 14.25 | 154  | 76766    | 8.17  | ng    | 99     |
| 17) Fluorene               | 15.23 | 166  | 138746   | 9.08  | ng    | 99     |
| 19) Hexachlorobenzene      | 16.24 | 284  | 24957    | 8.21  | ng    | 96     |
| 20) Pentachlorophenol      | 16.61 | 266  | 30313    | 17.88 | ng    | 92     |
| 21) Phenanthrene           | 16.98 | 178  | 202937   | 7.83  | ng    | 100    |
| 22) Anthracene             | 17.07 | 178  | 173744   | 7.90  | ng    | 99     |
| 23) Fluoranthene           | 19.01 | 202  | 194978   | 9.29  | ng    | 99     |
| 25) Pyrene                 | 19.37 | 202  | 210517   | 8.32  | ng    | 99     |
| 27) Benzo(a)anthracene     | 21.12 | 228  | 196005   | 8.68  | ng    | 97     |
| 28) Chrysene               | 21.17 | 228  | 183146   | 8.41  | ng    | 98     |
| 29) Indeno(1,2,3-cd)pyrene | 25.43 | 276  | 166918   | 9.66  | ng    | 99     |
| 31) Benzo(b)fluoranthene   | 22.64 | 252  | 183250   | 8.32  | ng    | 98     |
| 32) Benzo(k)fluoranthene   | 22.68 | 252  | 180066   | 8.90  | ng    | 99     |
| 33) Benzo(a)pyrene         | 23.18 | 252  | 169299   | 9.31  | ng    | 95     |
| 34) Dibenzo(a,h)anthracene | 25.44 | 278  | 135477   | 9.16  | ng    | 98     |
| 35) Benzo(g,h,i)perylene   | 26.07 | 276  | 136521   | 8.90  | ng    | 99     |

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

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