

Data Path : Z:\HPCHEM1\BNA\_E\DATA\BE080316\  
 Data File : BE092176.D  
 Acq On : 4 Aug 2016 10:36  
 Operator : UM/SJ  
 Sample : H4221-23MSD  
 Misc :  
 ALS Vial : 26 Sample Multiplier: 1

Instrument :  
 BNA\_E  
 ClientSampleId :  
 GW-44-8.6-15.0MSD

Quant Time: Aug 04 14:13:33 2016  
 Quant Method : Z:\HPCHEM1\BNA\_E\METHODS\8270-PAH-BE080216.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Aug 02 16:59:37 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc | Units | Dev(Min) |
|---------------------------|-------|------|----------|------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.17  | 152  | 11956    | 5.00 | ng    | 0.00     |
| 6) Naphthalene-d8         | 10.98 | 136  | 54782    | 5.00 | ng    | -0.02    |
| 10) Acenaphthene-d10      | 14.83 | 164  | 29638    | 5.00 | ng    | -0.02    |
| 16) Phenanthrene-d10      | 17.56 | 188  | 63872    | 5.00 | ng    | -0.02    |
| 22) Chrysene-d12          | 21.75 | 240  | 83661    | 5.00 | ng    | 0.00     |
| 28) Perylene-d12          | 24.14 | 264  | 82926    | 5.00 | ng    | 0.00     |

| System Monitoring Compounds | R.T.  | QIon | Response | Conc | Units | Dev(Min) |
|-----------------------------|-------|------|----------|------|-------|----------|
| 4) 2-Fluorophenol           | 5.69  | 112  | 15031    | 4.90 | ng    | 0.00     |
| 5) Phenol-d6                | 7.34  | 99   | 15436    | 3.77 | ng    | -0.01    |
| 7) Nitrobenzene-d5          | 9.34  | 82   | 15743    | 3.64 | ng    | -0.01    |
| 11) 2,4,6-Tribromophenol    | 16.32 | 330  | 7417     | 7.71 | ng    | -0.01    |
| 12) 2-Fluorobiphenyl        | 13.46 | 172  | 34371    | 3.66 | ng    | 0.00     |
| 24) Terphenyl-d14           | 20.19 | 244  | 53834    | 4.48 | ng    | -0.02    |

| Target Compounds           | R.T.  | QIon | Response | Conc | Units | Qvalue |
|----------------------------|-------|------|----------|------|-------|--------|
| 2) 1,4-Dioxane             | 3.44  | 88   | 3868     | 2.38 | ng    | # 89   |
| 3) n-Nitrosodimethylamine  | 3.78  | 42   | 6074     | 2.73 | ng    | # 97   |
| 8) Naphthalene             | 11.03 | 128  | 40876    | 3.47 | ng    | 95     |
| 9) 2-Methylnaphthalene     | 12.64 | 142  | 28419    | 3.65 | ng    | # 84   |
| 13) Acenaphthylene         | 14.54 | 152  | 187553   | 3.59 | ng    | 98     |
| 14) Acenaphthene           | 14.90 | 154  | 31180    | 3.74 | ng    | # 93   |
| 15) Fluorene               | 15.88 | 166  | 37758    | 3.57 | ng    | 99     |
| 17) Hexachlorobenzene      | 16.89 | 284  | 10616    | 3.40 | ng    | # 38   |
| 18) Pentachlorophenol      | 17.24 | 266  | 9015     | 7.68 | ng    | # 82   |
| 19) Phenanthrene           | 17.60 | 178  | 69100    | 3.97 | ng    | 99     |
| 20) Anthracene             | 17.70 | 178  | 64298    | 3.78 | ng    | 99     |
| 21) Fluoranthene           | 19.63 | 202  | 302836   | 3.88 | ng    | 97     |
| 23) Pyrene                 | 19.99 | 202  | 327367   | 3.74 | ng    | 98     |
| 25) Benzo(a)anthracene     | 21.72 | 228  | 354227   | 3.85 | ng    | 97     |
| 26) Chrysene               | 21.77 | 228  | 316075   | 3.43 | ng    | # 87   |
| 27) Indeno(1,2,3-cd)pyrene | 26.63 | 276  | 377916   | 3.78 | ng    | 100    |
| 29) Benzo(b)fluoranthene   | 23.40 | 252  | 86210    | 3.80 | ng    | 97     |
| 30) Benzo(k)fluoranthene   | 23.45 | 252  | 84485    | 3.89 | ng    | 95     |
| 31) Benzo(a)pyrene         | 24.02 | 252  | 81337    | 3.90 | ng    | 95     |
| 32) Dibenzo(a,h)anthracene | 26.67 | 278  | 78060    | 3.92 | ng    | # 92   |
| 33) Benzo(g,h,i)perylene   | 27.42 | 276  | 321415   | 3.78 | ng    | # 94   |

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

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