

Data Path : Z:\HPCHEM1\BNA G\DATA\BG061716\  
 Data File : BG022436.D  
 Acq On : 18 Jun 2016 23:18  
 Operator : UM/SJ  
 Sample : H3471-22  
 Misc :  
 ALS Vial : 45 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 STA-1141-(0-4)

Manual Integrations  
 APPROVED

sohil  
 6/20/2016 6:55:22 PM

Quant Time: Jun 20 03:57:13 2016  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060616.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Jun 13 17:40:23 2016  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|-------|------|----------|--------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 8.04  | 152  | 36951    | 20.00  | ng    | -0.05    |
| 21) Naphthalene-d8          | 10.86 | 136  | 189465   | 20.00  | ng    | -0.04    |
| 38) Acenaphthene-d10        | 14.68 | 164  | 103461   | 20.00  | ng    | -0.04    |
| 63) Phenanthrene-d10        | 17.43 | 188  | 192153   | 20.00  | ng    | -0.04    |
| 75) Chrysene-d12            | 21.69 | 240  | 155173   | 20.00  | ng    | -0.05    |
| 86) Perylene-d12            | 24.93 | 264  | 150609   | 20.00  | ng    | -0.08    |
| System Monitoring Compounds |       |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.61  | 112  | 276963   | 144.92 | ng    | -0.03    |
| 7) Phenol-d6                | 7.23  | 99   | 425452   | 141.59 | ng    | -0.02    |
| 23) Nitrobenzene-d5         | 9.22  | 82   | 260731   | 95.94  | ng    | -0.04    |
| 41) 2,4,6-Tribromophenol    | 16.18 | 330  | 106142   | 90.79  | ng    | -0.04    |
| 44) 2-Fluorobiphenyl        | 13.30 | 172  | 584043   | 86.03  | ng    | -0.04    |
| 78) Terphenyl-d14           | 20.02 | 244  | 497279   | 76.08  | ng    | -0.04    |
| Target Compounds            |       |      |          |        |       | Qvalue   |
| 31) Naphthalene             | 10.91 | 128  | 19453    | 2.06   | ng    | 96       |
| 49) Dimethylphthalate       | 14.13 | 163  | 137687   | 16.05  | ng    | 98       |
| 70) Phenanthrene            | 17.47 | 178  | 39035    | 3.74   | ng    | 97       |
| 74) Fluoranthene            | 19.47 | 202  | 143791   | 11.98  | ng    | 97       |
| 77) Pyrene                  | 19.84 | 202  | 130637   | 13.54  | ng    | 97       |
| 80) Benzo(a)anthracene      | 21.67 | 228  | 70648    | 8.12   | ng    | 97       |
| 82) Chrysene                | 21.73 | 228  | 86619    | 10.23  | ng    | 97       |
| 85) Indeno(1,2,3-cd)pyrene  | 28.65 | 276  | 38526    | 4.06   | ng    | # 76     |
| 87) Benzo(b)fluoranthene    | 23.90 | 252  | 115462   | 13.14  | ng    | # 95     |
| 88) Benzo(k)fluoranthene    | 23.96 | 252  | 40840m   | 4.72   | ng    |          |
| 89) Benzo(a)pyrene          | 24.78 | 252  | 55345    | 6.59   | ng    | # 90     |
| 91) Benzo(a,h,i)perylene    | 29.82 | 276  | 37396m   | 4.54   | ng    |          |

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

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