

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042916\  
 Data File : BE091737.D  
 Acq On : 29 Apr 2016 17:17  
 Operator : UM/SJ  
 Sample : H2769-04  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_E  
 ClientSampleId :  
 KY038PS038-160427

Manual Integrations  
 APPROVED

sohil  
 5/3/2016 6:30:32 PM

Quant Time: May 03 11:54:43 2016  
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE042616.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Apr 26 16:09:46 2016  
 Response via : Initial Calibration

| Internal Standards            | R.T.  | QIon | Response | Conc | Units | Dev(Min) |
|-------------------------------|-------|------|----------|------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4     | 7.80  | 152  | 37243    | 5.00 | ng    | 0.00     |
| 7) Naphthalene-d8             | 10.57 | 136  | 167780   | 5.00 | ng    | 0.00     |
| 13) Acenaphthene-d10          | 14.41 | 164  | 95600    | 5.00 | ng    | -0.01    |
| 19) Phenanthrene-d10          | 17.15 | 188  | 244751   | 5.00 | ng    | 0.00     |
| 26) Chrysene-d12              | 21.31 | 240  | 268070   | 5.00 | ng    | 0.00     |
| 35) Perylene-d12              | 23.74 | 264  | 270807   | 5.00 | ng    | -0.01    |
| System Monitoring Compounds   |       |      |          |      |       |          |
| 4) 2-Fluorophenol             | 5.39  | 112  | 17564    | 1.92 | ng    | 0.00     |
| 5) Phenol-d6                  | 6.95  | 99   | 12305    | 1.01 | ng    | 0.00     |
| 8) Nitrobenzene-d5            | 8.95  | 82   | 33910    | 3.13 | ng    | 0.00     |
| 14) 2,4,6-Tribromophenol      | 15.90 | 330  | 20434    | 5.79 | ng    | 0.00     |
| 15) 2-Fluorobiphenyl          | 13.03 | 172  | 101681   | 3.76 | ng    | 0.00     |
| 29) Terphenyl-d14             | 19.76 | 244  | 189148   | 4.52 | ng    | 0.00     |
| Target Compounds              |       |      |          |      |       | Qvalue   |
| 20) 4-Bromophenyl-phenylether | 16.33 | 248  | 224      | 0.03 | ng    | 95       |
| 21) Hexachlorobenzene         | 16.45 | 284  | 238      | 0.03 | ng    | # 84     |
| 22) Pentachlorophenol         | 16.79 | 266  | 107      | 0.12 | ng    | # 82     |
| 23) Phenanthrene              | 17.18 | 178  | 1777     | 0.03 | ng    | # 82     |
| 24) Anthracene                | 17.28 | 178  | 1295     | 0.03 | ng    | 99       |
| 25) Fluoranthene              | 19.19 | 202  | 1855     | 0.03 | ng    | 98       |
| 28) Pyrene                    | 19.55 | 202  | 2034     | 0.03 | ng    | 98       |
| 30) Benzo(a)anthracene        | 21.29 | 228  | 2720m    | 0.04 | ng    |          |
| 32) Chrysene                  | 21.33 | 228  | 2268m    | 0.03 | ng    |          |
| 34) Indeno(1,2,3-cd)pyrene    | 26.31 | 276  | 1344     | 0.02 | ng    | 97       |
| 36) Benzo(b)fluoranthene      | 23.00 | 252  | 1579     | 0.02 | ng    | # 56     |
| 37) Benzo(k)fluoranthene      | 23.04 | 252  | 1396     | 0.02 | ng    | # 56     |
| 40) Benzo(a,h,i)perylene      | 27.09 | 276  | 1202     | 0.02 | ng    | # 77     |

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

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042916\  
Data File : BE091737.D  
Acq On : 29 Apr 2016 17:17  
Operator : UM/SJ  
Sample : H2769-04  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_E  
ClientSampleId :  
KY038PS038-160427

Manual Integrations  
APPROVED

sohil  
5/3/2016 6:30:32 PM

Quant Time: May 03 11:54:43 2016  
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE042616.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Tue Apr 26 16:09:46 2016  
Response via : Initial Calibration

