

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071520\  
 Data File : BP002758.D  
 Acq On : 15 Jul 2020 17:21  
 Operator : CG/JU  
 Sample : L3303-01  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 FK-GF-03-014-A

Manual Integrations  
 APPROVED

mohammad  
 7/16/2020 11:07:38 AM

Quant Time: Jul 15 18:04:23 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Jul 13 10:54:03 2020  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units | Dev(Min)   |
|-----------------------------|-------|------|----------|--------|-------|------------|
| 1) 1,4-Dichlorobenzene-d4   | 7.55  | 152  | 150910   | 20.00  | ng    | 0.00       |
| 21) Naphthalene-d8          | 10.33 | 136  | 608776   | 20.00  | ng    | 0.00       |
| 39) Acenaphthene-d10        | 14.20 | 164  | 371963   | 20.00  | ng    | 0.00       |
| 64) Phenanthrene-d10        | 16.96 | 188  | 788371   | 20.00  | ng    | 0.00       |
| 76) Chrysene-d12            | 21.07 | 240  | 648891   | 20.00  | ng    | 0.00       |
| 86) Perylene-d12            | 23.26 | 264  | 617840   | 20.00  | ng    | 0.00       |
| System Monitoring Compounds |       |      |          |        |       |            |
| 5) 2-Fluorophenol           | 5.18  | 112  | 541619   | 60.55  | ng    | 0.00       |
| 7) Phenol-d6                | 6.76  | 99   | 729815   | 57.17  | ng    | 0.00       |
| 23) Nitrobenzene-d5         | 8.70  | 82   | 431372   | 37.90  | ng    | 0.00       |
| 42) 2,4,6-Tribromophenol    | 15.71 | 330  | 202689   | 52.20  | ng    | 0.00       |
| 45) 2-Fluorobiphenyl        | 12.82 | 172  | 904503   | 37.62  | ng    | 0.00       |
| 79) Terphenyl-d14           | 19.55 | 244  | 1136287  | 40.15  | ng    | 0.00       |
| Target Compounds            |       |      |          |        |       |            |
| 50) Dimethylphthalate       | 13.67 | 163  | 508164   | 17.914 | ng    | Qvalue 100 |
| 71) Phenanthrene            | 17.00 | 178  | 436349   | 10.222 | ng    | 99         |
| 72) Anthracene              | 17.10 | 178  | 123937   | 2.959  | ng    | 98         |
| 75) Fluoranthene            | 18.99 | 202  | 571763   | 11.560 | ng    | 100        |
| 78) Pyrene                  | 19.35 | 202  | 540947   | 12.074 | ng    | 99         |
| 81) Benzo(a)anthracene      | 21.06 | 228  | 239288   | 5.697  | ng    | 93         |
| 83) Chrysene                | 21.11 | 228  | 216970   | 5.426  | ng    | 97         |
| 87) Indeno(1,2,3-cd)pyrene  | 25.46 | 276  | 114347   | 2.561  | ng    | 97         |
| 88) Benzo(b)fluoranthene    | 22.61 | 252  | 237944m  | 6.202  | ng    |            |
| 89) Benzo(k)fluoranthene    | 22.64 | 252  | 103187m  | 2.731  | ng    |            |
| 90) Benzo(a)pyrene          | 23.16 | 252  | 185910   | 5.114  | ng    | # 93       |
| 92) Benzo(a,h,i)perylene    | 26.13 | 276  | 124300   | 3.404  | ng    | # 90       |

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

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071520\  
Data File : BP002758.D  
Acq On : 15 Jul 2020 17:21  
Operator : CG/JU  
Sample : L3303-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
FK-GF-03-014-A

Manual Integrations  
APPROVED

mohammad  
7/16/2020 11:07:38 AM

Quant Time: Jul 15 18:04:23 2020  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Mon Jul 13 10:54:03 2020  
Response via : Initial Calibration

