

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP070323\  
 Data File : BP015948.D  
 Acq On : 03 Jul 2023 13:21  
 Operator : MA/JU  
 Sample : 03243-12  
 Misc :  
 ALS Vial : 28 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 DCGD2

Manual Integrations  
 APPROVED

Reviewed By :Sohil Jodhani 07/04/2023  
 Supervised By :mohammad ahmed 07/04/2023

Quant Time: Jul 03 23:11:06 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP062623.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sun Jul 02 06:12:16 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.052  | 152  | 152049   | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.881 | 136  | 609449   | 20.000 | ng/u1  | # 0.01   |
| 38) Acenaphthene-d10          | 14.704 | 164  | 350746   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.463 | 188  | 720688   | 20.000 | ng/u1  | # 0.00   |
| 79) Chrysene-d12              | 21.551 | 240  | 671490   | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 24.092 | 264  | 802891   | 20.000 | ng/u1  | 0.00     |
|                               |        |      |          |        |        |          |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.370  | 96   | 16374    | 3.875  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.823  | 84   | 149251   | 14.010 | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 7.252  | 99   | 254372   | 19.553 | ng/u1  | 0.03     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.375  | 67   | 168616   | 20.949 | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.587  | 132  | 222357   | 22.642 | ng/u1  | 0.01     |
| 15) 4-Methylphenol-d8         | 8.816  | 113  | 166306   | 16.182 | ng/u1  | 0.04     |
| 21) Nitrobenzene-d5           | 9.258  | 128  | 100625   | 21.604 | ng/u1  | 0.02     |
| 24) 2-Nitrophenol-d4          | 9.993  | 143  | 106764   | 19.640 | ng/u1  | 0.04     |
| 28) 2,4-Dichlorophenol-d3     | 10.581 | 165  | 177625m  | 18.336 | ng/u1  | 0.07     |
| 31) 4-Chloroaniline-d4        | 11.099 | 131  | 129644   | 10.418 | ng/u1  | 0.07     |
| 46) Dimethylphthalate-d6      | 14.116 | 166  | 638135   | 23.539 | ng/u1  | 0.01     |
| 49) Acenaphthylene-d8         | 14.398 | 160  | 748017   | 24.545 | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 15.034 | 143  | 65063m   | 18.186 | ng/u1  | 0.08     |
| 60) Fluorene-d10              | 15.704 | 176  | 559445   | 25.062 | ng/u1  | 0.01     |
| 65) 4,6-Dinitro-2-methylph... | 15.898 | 200  | 49758m   | 11.227 | ng/u1  | 0.04     |
| 73) Anthracene-d10            | 17.569 | 188  | 818124   | 24.852 | ng/u1  | 0.01     |
| 81) Pyrene-d10                | 19.792 | 212  | 1027373  | 23.431 | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.927 | 264  | 1051183  | 25.954 | ng/u1  | 0.00     |
|                               |        |      |          |        |        |          |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 72) Phenanthrene              | 17.504 | 178  | 74242    | 1.896  | ng/u1  | 96       |
| 80) Fluoranthene              | 19.469 | 202  | 241081   | 4.424  | ng/u1# | 92       |
| 82) Pyrene                    | 19.822 | 202  | 207345   | 3.730  | ng/u1# | 91       |
| 85) Benzo(a)anthracene        | 21.539 | 228  | 149489   | 3.165  | ng/u1  | 96       |
| 87) Chrysene                  | 21.592 | 228  | 157575   | 3.516  | ng/u1  | 98       |
| 90) Benzo(b)fluoranthene      | 23.316 | 252  | 239499   | 4.642  | ng/u1# | 90       |
| 91) Benzo(k)fluoranthene      | 23.357 | 252  | 86437    | 1.658  | ng/u1# | 86       |
| 93) Benzo(a)pyrene            | 23.980 | 252  | 151806   | 3.368  | ng/u1# | 93       |
| 94) Indeno(1,2,3-cd)pyrene    | 26.810 | 276  | 121623   | 1.987  | ng/u1# | 94       |
| 96) Benzo(g,h,i)perylene      | 27.651 | 276  | 106670   | 2.119  | ng/u1# | 91       |
| -----                         |        |      |          |        |        |          |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP070323\  
Data File : BP015948.D  
Acq On : 03 Jul 2023 13:21  
Operator : MA/JU  
Sample : 03243-12  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCGD2

Quant Time: Jul 03 23:11:06 2023  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP062623.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Sun Jul 02 06:12:16 2023  
Response via : Initial Calibration

Manual Integrations  
APPROVED

Reviewed By :Sohil Jodhani 07/04/2023  
Supervised By :mohammad ahmed 07/04/2023

