

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP070824\  
 Data File : BP020848.D  
 Acq On : 09 Jul 2024 02:36  
 Operator : MA/JU  
 Sample : P3035-14  
 Misc :  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 A4CW9

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 07/09/2024  
 Supervised By :mohammad ahmed 07/10/2024

Quant Time: Jul 09 03:40:04 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP062524.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Jul 03 15:01:56 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.722  | 152  | 225833   | 20.000 | ng/u1  | -0.01    |
| 20) Naphthalene-d8            | 10.493 | 136  | 937204   | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.357 | 164  | 631616   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.169 | 188  | 1329315  | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.616 | 240  | 1210973  | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 24.957 | 264  | 1417183  | 20.000 | ng/u1  | -0.02    |
|                               |        |      |          |        |        |          |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.264  | 96   | 16965    | 3.226  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.693  | 84   | 20870m   | 1.362  | ng/u1  | 0.01     |
| 7) Phenol-d5                  | 6.916  | 99   | 395880   | 21.373 | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.069  | 67   | 240019   | 20.440 | ng/u1  | -0.01    |
| 11) 2-Chlorophenol-d4         | 7.269  | 132  | 330306   | 21.820 | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.434  | 113  | 326418   | 21.698 | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 8.881  | 128  | 162119   | 23.430 | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.587  | 143  | 186807   | 26.978 | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.128 | 165  | 328234   | 23.013 | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.640 | 131  | 317026   | 15.860 | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.757 | 166  | 1047566  | 23.854 | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 14.045 | 160  | 1161493  | 22.165 | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.598 | 143  | 146529   | 21.838 | ng/u1  | 0.02     |
| 60) Fluorene-d10              | 15.369 | 176  | 892565   | 24.153 | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.510 | 200  | 136684   | 20.404 | ng/u1  | 0.01     |
| 73) Anthracene-d10            | 17.275 | 188  | 1393902  | 23.364 | ng/u1  | 0.01     |
| 81) Pyrene-d10                | 19.651 | 212  | 1462484  | 20.103 | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.727 | 264  | 1112484  | 16.138 | ng/u1  | -0.01    |
|                               |        |      |          |        |        |          |
| Target Compounds              |        |      |          |        | Qvalue |          |
| 72) Phenanthrene              | 17.216 | 178  | 267621   | 3.845  | ng/u1  | 99       |
| 80) Fluoranthene              | 19.304 | 202  | 562018   | 6.324  | ng/u1  | 100      |
| 82) Pyrene                    | 19.686 | 202  | 421011   | 4.620  | ng/u1  | 98       |
| 85) Benzo(a)anthracene        | 21.592 | 228  | 268858   | 3.168  | ng/u1  | 96       |
| 86) Bis(2-ethylhexyl)phtha... | 21.522 | 149  | 175673   | 3.343  | ng/u1  | 99       |
| 87) Chrysene                  | 21.663 | 228  | 277535   | 3.459  | ng/u1  | 97       |
| 90) Benzo(b)fluoranthene      | 23.892 | 252  | 386018   | 4.496  | ng/u1  | 96       |
| 91) Benzo(k)fluoranthene      | 23.963 | 252  | 118748m  | 1.367  | ng/u1  |          |
| 93) Benzo(a)pyrene            | 24.798 | 252  | 145713   | 1.798  | ng/u1  | 94       |
| 94) Indeno(1,2,3-cd)pyrene    | 28.721 | 276  | 133231m  | 1.282  | ng/u1  |          |
| -----                         |        |      |          |        |        |          |

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

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