

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP070524\  
 Data File : BP020808.D  
 Acq On : 06 Jul 2024 05:33  
 Operator : MA/JU  
 Sample : P2964-11  
 Misc :  
 ALS Vial : 27 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 A4CS5

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 07/08/2024  
 Supervised By :mohammad ahmed 07/08/2024

Quant Time: Jul 06 06:01:06 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     | 7.728  | 152  | 232895   | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.498 | 136  | 958659   | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.357 | 164  | 615816   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.175 | 188  | 1236087  | 20.000 | ng/u1  | 0.01     |
| 79) Chrysene-d12              | 21.633 | 240  | 1037479  | 20.000 | ng/u1  | 0.01     |
| 88) Perylene-d12              | 25.004 | 264  | 1254684  | 20.000 | ng/u1  | 0.03     |
|                               |        |      |          |        |        |          |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.264  | 96   | 16278    | 3.001  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 0.000  | 84   | 0d       | 0.000  | ng/u1  |          |
| 7) Phenol-d5                  | 6.922  | 99   | 375022   | 19.633 | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.075  | 67   | 224083   | 18.504 | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.269  | 132  | 321813   | 20.614 | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.434  | 113  | 298449   | 19.237 | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 8.881  | 128  | 158118   | 22.340 | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.587  | 143  | 180302   | 25.456 | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.134 | 165  | 310466   | 21.280 | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.645 | 131  | 78770    | 3.852  | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.769 | 166  | 963988   | 22.514 | ng/u1  | 0.01     |
| 49) Acenaphthylene-d8         | 14.045 | 160  | 1101546  | 21.560 | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.598 | 143  | 123236   | 18.838 | ng/u1  | 0.02     |
| 60) Fluorene-d10              | 15.369 | 176  | 829671   | 23.027 | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.504 | 200  | 96822    | 15.543 | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.269 | 188  | 1202024  | 21.668 | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.663 | 212  | 1169141  | 18.758 | ng/u1  | 0.01     |
| 92) Benzo(a)pyrene-d12        | 24.768 | 264  | 904599   | 14.822 | ng/u1  | 0.03     |
|                               |        |      |          |        |        |          |
| Target Compounds              |        |      |          |        | Qvalue |          |
| 30) Naphthalene               | 10.551 | 128  | 90169    | 1.757  | ng/u1  | 99       |
| 36) 2-Methylnaphthalene       | 12.151 | 142  | 35087    | 1.021  | ng/u1  | 98       |
| 50) Acenaphthylene            | 14.081 | 152  | 144737   | 2.508  | ng/u1  | 99       |
| 56) Dibenzofuran              | 14.763 | 168  | 53919    | 1.045  | ng/u1  | 96       |
| 61) Fluorene                  | 15.422 | 166  | 45040    | 1.087  | ng/u1# | 99       |
| 72) Phenanthrene              | 17.216 | 178  | 903266   | 13.955 | ng/u1  | 99       |
| 74) Anthracene                | 17.304 | 178  | 238387   | 3.580  | ng/u1  | 99       |
| 77) Carbazole                 | 17.598 | 167  | 79639    | 1.449  | ng/u1  | 95       |
| 80) Fluoranthene              | 19.310 | 202  | 1713111  | 22.501 | ng/u1  | 99       |
| 82) Pyrene                    | 19.692 | 202  | 1379191  | 17.664 | ng/u1  | 98       |
| 85) Benzo(a)anthracene        | 21.615 | 228  | 920051   | 12.652 | ng/u1# | 90       |
| 87) Chrysene                  | 21.686 | 228  | 1000328  | 14.553 | ng/u1  | 96       |
| 90) Benzo(b)fluoranthene      | 23.939 | 252  | 1352171  | 17.789 | ng/u1# | 95       |
| 91) Benzo(k)fluoranthene      | 24.004 | 252  | 464223m  | 6.038  | ng/u1  |          |
| 93) Benzo(a)pyrene            | 24.833 | 252  | 544042   | 7.582  | ng/u1# | 92       |
| 94) Indeno(1,2,3-cd)pyrene    | 28.821 | 276  | 647989   | 7.041  | ng/u1  | 96       |
| 95) Dibenzo(a,h)anthracene    | 28.886 | 278  | 229223m  | 3.030  | ng/u1  |          |
| 96) Benzo(g,h,i)perylene      | 29.997 | 276  | 92956m   | 1.237  | ng/u1  |          |
| -----                         |        |      |          |        |        |          |

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

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