

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP072524\  
 Data File : BP021153.D  
 Acq On : 25 Jul 2024 18:49  
 Operator : MA/JU  
 Sample : P3263-17 5X  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 A4CK5

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 07/26/2024  
 Supervised By :mohammad ahmed 07/27/2024

Quant Time: Jul 25 23:26:02 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP071024.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Jul 25 00:15:24 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.652  | 152  | 207042   | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.416 | 136  | 787193   | 20.000 | ng/u1  | -0.01    |
| 38) Acenaphthene-d10          | 14.299 | 164  | 451852   | 20.000 | ng/u1  | 0.01     |
| 64) Phenanthrene-d10          | 17.110 | 188  | 803571   | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.563 | 240  | 798767   | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 24.898 | 264  | 1060990  | 20.000 | ng/u1  | 0.03     |
|                               |        |      |          |        |        |          |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.181  | 96   | 3731     | 0.820  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 0.000  | 84   | 0d       | 0.000  | ng/u1  |          |
| 7) Phenol-d5                  | 6.875  | 99   | 72580    | 4.300  | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.005  | 67   | 43945    | 4.296  | ng/u1  | 0.01     |
| 11) 2-Chlorophenol-d4         | 7.205  | 132  | 62260    | 4.477  | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.393  | 113  | 56297    | 4.016  | ng/u1  | 0.01     |
| 21) Nitrobenzene-d5           | 8.816  | 128  | 27149    | 4.440  | ng/u1  | 0.01     |
| 24) 2-Nitrophenol-d4          | 9.528  | 143  | 30928    | 4.419  | ng/u1  | 0.01     |
| 28) 2,4-Dichlorophenol-d3     | 10.093 | 165  | 53406    | 4.220  | ng/u1  | 0.02     |
| 31) 4-Chloroaniline-d4        | 10.605 | 131  | 7633m    | 0.455  | ng/u1  | 0.04     |
| 46) Dimethylphthalate-d6      | 13.699 | 166  | 165916   | 5.051  | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 13.987 | 160  | 184642   | 5.006  | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.640 | 143  | 15049m   | 3.220  | ng/u1  | 0.04     |
| 60) Fluorene-d10              | 15.304 | 176  | 137596   | 5.106  | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.487 | 200  | 10705    | 2.202  | ng/u1  | 0.02     |
| 73) Anthracene-d10            | 17.210 | 188  | 187760   | 5.261  | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.604 | 212  | 188881   | 3.825  | ng/u1  | 0.02     |
| 92) Benzo(a)pyrene-d12        | 24.657 | 264  | 186800   | 3.570  | ng/u1  | 0.00     |
|                               |        |      |          |        |        |          |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 6) Benzaldehyde               | 6.828  | 77   | 25452    | 2.991  | ng/u1  | 97       |
| 30) Naphthalene               | 10.469 | 128  | 76524    | 1.848  | ng/u1  | 97       |
| 50) Acenaphthylene            | 14.016 | 152  | 95028    | 2.276  | ng/u1  | 98       |
| 56) Dibenzofuran              | 14.704 | 168  | 39947    | 1.052  | ng/u1  | 98       |
| 72) Phenanthrene              | 17.151 | 178  | 430886   | 10.336 | ng/u1  | 100      |
| 74) Anthracene                | 17.251 | 178  | 151781   | 3.572  | ng/u1  | 99       |
| 80) Fluoranthene              | 19.257 | 202  | 796236   | 13.351 | ng/u1  | 100      |
| 82) Pyrene                    | 19.634 | 202  | 676187   | 11.142 | ng/u1  | 99       |
| 85) Benzo(a)anthracene        | 21.545 | 228  | 518762m  | 9.271  | ng/u1  |          |
| 87) Chrysene                  | 21.610 | 228  | 597028   | 11.483 | ng/u1  | 95       |
| 90) Benzo(b)fluoranthene      | 23.845 | 252  | 907665   | 14.097 | ng/u1# | 92       |
| 91) Benzo(k)fluoranthene      | 23.910 | 252  | 292199   | 4.547  | ng/u1# | 86       |
| 93) Benzo(a)pyrene            | 24.727 | 252  | 353640   | 5.812  | ng/u1# | 90       |
| 94) Indeno(1,2,3-cd)pyrene    | 28.692 | 276  | 544969m  | 6.951  | ng/u1  |          |
| 95) Dibenzo(a,h)anthracene    | 28.715 | 278  | 177010m  | 2.726  | ng/u1  |          |
| 96) Benzo(g,h,i)perylene      | 29.868 | 276  | 81036m   | 1.269  | ng/u1  |          |

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

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