

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080824\  
 Data File : BP021477.D  
 Acq On : 09 Aug 2024 19:37  
 Operator : RC/JU  
 Sample : P3329-05MEDL 4X  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 F7E07MEDL

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 08/12/2024  
 Supervised By :mohammad ahmed 08/12/2024

Quant Time: Aug 09 23:17:33 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP071024.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Aug 08 18:23:21 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.616  | 152  | 152888   | 20.000 | ng/u1  | 0.02     |
| 20) Naphthalene-d8            | 10.369 | 136  | 666890   | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.245 | 164  | 458056   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.075 | 188  | 993168   | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.521 | 240  | 999479   | 20.000 | ng/u1  | 0.01     |
| 88) Perylene-d12              | 24.792 | 264  | 1123581  | 20.000 | ng/u1  | 0.01     |
|                               |        |      |          |        |        |          |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.164  | 96   | 1100     | 0.327  | ng/uL  | 0.01     |
| 4) Pyridine-d5                | 0.000  | 84   | 0d       | 0.000  | ng/u1  |          |
| 7) Phenol-d5                  | 6.993  | 99   | 8624m    | 0.692  | ng/u1  | 0.16     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.040  | 67   | 15818m   | 2.094  | ng/u1  | 0.08     |
| 11) 2-Chlorophenol-d4         | 7.269  | 132  | 17977m   | 1.750  | ng/u1  | 0.11     |
| 15) 4-Methylphenol-d8         | 8.493  | 113  | 20327m   | 1.964  | ng/u1  | 0.15     |
| 21) Nitrobenzene-d5           | 8.869  | 128  | 8711m    | 1.682  | ng/u1  | 0.09     |
| 24) 2-Nitrophenol-d4          | 9.551  | 143  | 8401m    | 1.417  | ng/u1  | 0.06     |
| 28) 2,4-Dichlorophenol-d3     | 10.204 | 165  | 20222m   | 1.886  | ng/u1  | 0.16     |
| 31) 4-Chloroaniline-d4        | 10.651 | 131  | 23634m   | 1.664  | ng/u1  | 0.11     |
| 46) Dimethylphthalate-d6      | 13.692 | 166  | 92850    | 2.788  | ng/u1  | 0.04     |
| 49) Acenaphthylene-d8         | 13.951 | 160  | 88871    | 2.377  | ng/u1  | 0.02     |
| 54) 4-Nitrophenol-d4          | 0.000  | 143  | 0d       | 0.000  | ng/u1  |          |
| 60) Fluorene-d10              | 15.292 | 176  | 73468    | 2.689  | ng/u1  | 0.03     |
| 65) 4,6-Dinitro-2-methylph... | 0.000  | 200  | 0d       | 0.000  | ng/u1  |          |
| 73) Anthracene-d10            | 17.198 | 188  | 132945   | 3.014  | ng/u1  | 0.02     |
| 81) Pyrene-d10                | 19.569 | 212  | 165367   | 2.676  | ng/u1  | 0.01     |
| 92) Benzo(a)pyrene-d12        | 24.580 | 264  | 161592   | 2.916  | ng/u1  | 0.02     |
|                               |        |      |          |        |        |          |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 30) Naphthalene               | 10.422 | 128  | 1510286  | 43.054 | ng/u1  | 99       |
| 36) 2-Methylnaphthalene       | 12.051 | 142  | 287032   | 11.920 | ng/u1  | 98       |
| 37) 1-Methylnaphthalene       | 12.275 | 142  | 133822   | 5.404  | ng/u1  | 97       |
| 43) 1,1'-Biphenyl             | 13.087 | 154  | 66712    | 1.993  | ng/u1  | 95       |
| 52) Acenaphthene              | 14.322 | 153  | 221441   | 7.881  | ng/u1  | 99       |
| 56) Dibenzofuran              | 14.681 | 168  | 228808   | 5.947  | ng/u1  | 96       |
| 61) Fluorene                  | 15.345 | 166  | 225209   | 7.173  | ng/u1  | 99       |
| 72) Phenanthrene              | 17.122 | 178  | 1138462  | 22.097 | ng/u1  | 100      |
| 74) Anthracene                | 17.233 | 178  | 236203m  | 4.497  | ng/u1  |          |
| 77) Carbazole                 | 17.551 | 167  | 131571   | 3.064  | ng/u1  | 99       |
| 80) Fluoranthene              | 19.222 | 202  | 705535   | 9.455  | ng/u1  | 98       |
| 82) Pyrene                    | 19.598 | 202  | 478161   | 6.297  | ng/u1  | 98       |
| 85) Benzo(a)anthracene        | 21.510 | 228  | 150160m  | 2.145  | ng/u1  |          |
| 87) Chrysene                  | 21.568 | 228  | 125281   | 1.926  | ng/u1  | 96       |
| 90) Benzo(b)fluoranthene      | 23.780 | 252  | 76328    | 1.119  | ng/u1  | 98       |
| -----                         |        |      |          |        |        |          |

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

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