

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080824\  
 Data File : BP021476.D  
 Acq On : 09 Aug 2024 18:50  
 Operator : RC/JU  
 Sample : P3329-04MEDL 5X  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 F7E03MEDL

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 08/10/2024  
 Supervised By :mohammad ahmed 08/12/2024

Quant Time: Aug 09 23:12:05 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     | 7.634  | 152  | 121087   | 20.000 | ng/u1  | 0.03     |
| 20) Naphthalene-d8            | 10.375 | 136  | 463651   | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.263 | 164  | 349685   | 20.000 | ng/u1  | 0.02     |
| 64) Phenanthrene-d10          | 17.075 | 188  | 679037   | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.521 | 240  | 759754   | 20.000 | ng/u1  | 0.01     |
| 88) Perylene-d12              | 24.798 | 264  | 878018   | 20.000 | ng/u1  | 0.02     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.158  | 96   | 1324     | 0.498  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 0.000  | 84   | 0d       | 0.000  | ng/u1  |          |
| 7) Phenol-d5                  | 7.005  | 99   | 5063m    | 0.513  | ng/u1  | 0.18     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.069  | 67   | 15580m   | 2.604  | ng/u1  | 0.11     |
| 11) 2-Chlorophenol-d4         | 7.287  | 132  | 16395m   | 2.016  | ng/u1  | 0.13     |
| 15) 4-Methylphenol-d8         | 8.510  | 113  | 16394m   | 1.999  | ng/u1  | 0.16     |
| 21) Nitrobenzene-d5           | 8.875  | 128  | 8384m    | 2.328  | ng/u1  | 0.10     |
| 24) 2-Nitrophenol-d4          | 9.557  | 143  | 8360m    | 2.028  | ng/u1  | 0.07     |
| 28) 2,4-Dichlorophenol-d3     | 10.246 | 165  | 14872m   | 1.995  | ng/u1  | 0.20     |
| 31) 4-Chloroaniline-d4        | 10.693 | 131  | 18158m   | 1.839  | ng/u1  | 0.15     |
| 46) Dimethylphthalate-d6      | 13.704 | 166  | 71382    | 2.808  | ng/u1  | 0.05     |
| 49) Acenaphthylene-d8         | 13.963 | 160  | 68304    | 2.393  | ng/u1  | 0.03     |
| 54) 4-Nitrophenol-d4          | 0.000  | 143  | 0d       | 0.000  | ng/u1  |          |
| 60) Fluorene-d10              | 15.287 | 176  | 61993    | 2.973  | ng/u1  | 0.02     |
| 65) 4,6-Dinitro-2-methylph... | 0.000  | 200  | 0d       | 0.000  | ng/u1  |          |
| 73) Anthracene-d10            | 17.192 | 188  | 81085    | 2.689  | ng/u1  | 0.02     |
| 81) Pyrene-d10                | 19.580 | 212  | 110619   | 2.355  | ng/u1  | 0.02     |
| 92) Benzo(a)pyrene-d12        | 24.592 | 264  | 121186   | 2.799  | ng/u1  | 0.03     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 30) Naphthalene               | 10.428 | 128  | 727358   | 29.824 | ng/u1  | 100      |
| 36) 2-Methylnaphthalene       | 12.081 | 142  | 138566   | 8.277  | ng/u1  | 95       |
| 37) 1-Methylnaphthalene       | 12.298 | 142  | 64418    | 3.741  | ng/u1# | 91       |
| 43) 1,1'-Biphenyl             | 13.116 | 154  | 32805    | 1.284  | ng/u1  | 91       |
| 52) Acenaphthene              | 14.328 | 153  | 101677   | 4.740  | ng/u1  | 99       |
| 56) Dibenzofuran              | 14.704 | 168  | 121812   | 4.147  | ng/u1  | 98       |
| 61) Fluorene                  | 15.363 | 166  | 122238   | 5.100  | ng/u1  | 97       |
| 72) Phenanthrene              | 17.116 | 178  | 466059   | 13.231 | ng/u1  | 99       |
| 74) Anthracene                | 17.239 | 178  | 275843   | 7.681  | ng/u1  | 98       |
| 77) Carbazole                 | 17.545 | 167  | 117145   | 3.990  | ng/u1  | 99       |
| 80) Fluoranthene              | 19.227 | 202  | 275939   | 4.865  | ng/u1  | 100      |
| 82) Pyrene                    | 19.610 | 202  | 194455   | 3.369  | ng/u1  | 97       |
| 85) Benzo(a)anthracene        | 21.510 | 228  | 64006    | 1.203  | ng/u1# | 94       |
| 87) Chrysene                  | 21.574 | 228  | 54595    | 1.104  | ng/u1  | 96       |
| -----                         |        |      |          |        |        |          |

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

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