

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP121823\  
 Data File : BP018944.D  
 Acq On : 19 Dec 2023 12:51  
 Operator : MA/JU  
 Sample : 05818-02DL 5X  
 Misc :  
 ALS Vial : 48 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BGRG7DL

Quant Time: Dec 20 00:16:00 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120123.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Dec 19 05:28:03 2023  
 Response via : Initial Calibration

### Manual Integrations APPROVED

Reviewed By :Yogesh Patel 12/20/2023  
 Supervised By :mohammad ahmed 12/20/2023

12/20/2023  
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 ahmed

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.828  | 152  | 215051   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.622 | 136  | 838772   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.463 | 164  | 540096   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.210 | 188  | 1274902  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.303 | 240  | 1260254  | 20.000 | ng/u1 | -0.01    |
| 88) Perylene-d12          | 23.674 | 264  | 1519439  | 20.000 | ng/u1 | -0.01    |

|                               |        |     |        |       |       |       |
|-------------------------------|--------|-----|--------|-------|-------|-------|
| System Monitoring Compounds   |        |     |        |       |       |       |
| 3) 1,4-Dioxane-d8             | 3.263  | 96  | 3512   | 0.557 | ng/uL | 0.00  |
| 4) Pyridine-d5                | 0.000  | 84  | 0d     | 0.000 | ng/u1 |       |
| 7) Phenol-d5                  | 7.004  | 99  | 62812  | 2.911 | ng/u1 | -0.01 |
| 9) Bis-(2-Chloroethyl)eth...  | 7.169  | 67  | 37281  | 2.915 | ng/u1 | 0.00  |
| 11) 2-Chlorophenol-d4         | 7.363  | 132 | 51324  | 3.048 | ng/u1 | 0.00  |
| 15) 4-Methylphenol-d8         | 8.545  | 113 | 48569  | 2.781 | ng/u1 | -0.01 |
| 21) Nitrobenzene-d5           | 8.992  | 128 | 23664  | 3.170 | ng/u1 | 0.00  |
| 24) 2-Nitrophenol-d4          | 9.710  | 143 | 24615  | 3.180 | ng/u1 | 0.00  |
| 28) 2,4-Dichlorophenol-d3     | 10.251 | 165 | 51130  | 3.231 | ng/u1 | 0.00  |
| 31) 4-Chloroaniline-d4        | 10.775 | 131 | 32552  | 1.574 | ng/u1 | 0.00  |
| 46) Dimethylphthalate-d6      | 13.874 | 166 | 157317 | 3.264 | ng/u1 | 0.00  |
| 49) Acenaphthylene-d8         | 14.157 | 160 | 184345 | 3.473 | ng/u1 | 0.00  |
| 54) 4-Nitrophenol-d4          | 14.692 | 143 | 10385m | 1.312 | ng/u1 | 0.00  |
| 60) Fluorene-d10              | 15.451 | 176 | 151749 | 3.750 | ng/u1 | 0.00  |
| 65) 4,6-Dinitro-2-methylph... | 15.586 | 200 | 4408   | 0.577 | ng/u1 | 0.00  |
| 73) Anthracene-d10            | 17.310 | 188 | 248790 | 3.712 | ng/u1 | 0.00  |
| 81) Pyrene-d10                | 19.551 | 212 | 310523 | 3.165 | ng/u1 | 0.00  |
| 92) Benzo(a)pyrene-d12        | 23.515 | 264 | 309666 | 3.495 | ng/u1 | -0.02 |

|                            |        |     |         |        |        |     |
|----------------------------|--------|-----|---------|--------|--------|-----|
| Target Compounds           |        |     |         | Qvalue |        |     |
| 16) Acetophenone           | 8.657  | 105 | 44961   | 1.727  | ng/u1  | 98  |
| 50) Acenaphthylene         | 14.186 | 152 | 75569   | 1.266  | ng/u1  | 99  |
| 72) Phenanthrene           | 17.257 | 178 | 727618  | 9.450  | ng/u1  | 100 |
| 74) Anthracene             | 17.345 | 178 | 216941  | 2.811  | ng/u1  | 99  |
| 80) Fluoranthene           | 19.227 | 202 | 2200386 | 18.975 | ng/u1  | 96  |
| 82) Pyrene                 | 19.580 | 202 | 2566806 | 21.942 | ng/u1  | 94  |
| 85) Benzo(a)anthracene     | 21.292 | 228 | 1771656 | 17.319 | ng/u1  | 98  |
| 87) Chrysene               | 21.339 | 228 | 1803067 | 19.141 | ng/u1  | 98  |
| 90) Benzo(b)fluoranthene   | 22.950 | 252 | 2439409 | 21.775 | ng/u1  | 97  |
| 91) Benzo(k)fluoranthene   | 22.997 | 252 | 712463m | 6.573  | ng/u1  |     |
| 93) Benzo(a)pyrene         | 23.568 | 252 | 1812317 | 17.704 | ng/u1  | 97  |
| 94) Indeno(1,2,3-cd)pyrene | 26.133 | 276 | 1348323 | 11.024 | ng/u1# | 94  |
| 95) Dibenzo(a,h)anthracene | 26.139 | 278 | 390244  | 3.836  | ng/u1  | 94  |
| 96) Benzo(g,h,i)perylene   | 26.886 | 276 | 465998  | 4.734  | ng/u1# | 94  |

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

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