

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP090825\  
 Data File : BP025684.D  
 Acq On : 08 Sep 2025 14:41  
 Operator : CG/JU  
 Sample : SSTD00510  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTD005639

Manual Integrations  
 APPROVED

Reviewed By :Rahul  
 Chavli

09/09/2025  
 Supervised By :Jagrut  
 Upadhyay

Quant Time: Sep 08 17:20:03 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP090825.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Sep 08 16:53:59 2025  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.761  | 152  | 125875   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.525 | 136  | 495930   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.378 | 164  | 356358   | 20.000 | ng/u1 | -0.01    |
| 64) Phenanthrene-d10      | 17.178 | 188  | 782622   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.625 | 240  | 869821   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12          | 24.983 | 264  | 950861   | 20.000 | ng/u1 | 0.02     |

|                               |        |     |        |       |       |      |
|-------------------------------|--------|-----|--------|-------|-------|------|
| System Monitoring Compounds   |        |     |        |       |       |      |
| 3) 1,4-Dioxane-d8             | 3.278  | 96  | 5994   | 2.098 | ng/uL | 0.00 |
| 4) Pyridine-d5                | 0.000  | 84  | 0d     | 0.000 | ng/u1 |      |
| 7) Phenol-d5                  | 0.000  | 99  | 0d     | 0.000 | ng/u1 |      |
| 9) Bis-(2-Chloroethyl)eth...  | 0.000  | 67  | 0d     | 0.000 | ng/u1 |      |
| 11) 2-Chlorophenol-d4         | 7.308  | 132 | 35263  | 4.482 | ng/u1 | 0.00 |
| 15) 4-Methylphenol-d8         | 0.000  | 113 | 0d     | 0.000 | ng/u1 |      |
| 21) Nitrobenzene-d5           | 8.908  | 128 | 16446  | 4.790 | ng/u1 | 0.00 |
| 24) 2-Nitrophenol-d4          | 9.619  | 143 | 18724  | 5.815 | ng/u1 | 0.00 |
| 28) 2,4-Dichlorophenol-d3     | 10.184 | 165 | 33060  | 4.428 | ng/u1 | 0.02 |
| 31) 4-Chloroaniline-d4        | 0.000  | 131 | 0d     | 0.000 | ng/u1 |      |
| 46) Dimethylphthalate-d6      | 13.778 | 166 | 131614 | 5.110 | ng/u1 | 0.00 |
| 49) Acenaphthylene-d8         | 14.066 | 160 | 138909 | 4.672 | ng/u1 | 0.00 |
| 54) 4-Nitrophenol-d4          | 0.000  | 143 | 0d     | 0.000 | ng/u1 |      |
| 60) Fluorene-d10              | 15.390 | 176 | 108545 | 4.929 | ng/u1 | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 0.000  | 200 | 0d     | 0.000 | ng/u1 |      |
| 73) Anthracene-d10            | 17.284 | 188 | 181335 | 4.848 | ng/u1 | 0.00 |
| 81) Pyrene-d10                | 19.648 | 212 | 229064 | 5.071 | ng/u1 | 0.00 |
| 92) Benzo(a)pyrene-d12        | 24.754 | 264 | 231883 | 4.874 | ng/u1 | 0.02 |

| Target Compounds              |        |     |        |       | Qvalue    |
|-------------------------------|--------|-----|--------|-------|-----------|
| 2) 1,4-Dioxane                | 3.314  | 88  | 6131   | 1.992 | ng/uL# 95 |
| 12) 2-Chlorophenol            | 7.343  | 128 | 34672  | 4.308 | ng/u1 94  |
| 17) N-Nitroso-di-n-propyla... | 8.555  | 70  | 30828  | 5.214 | ng/u1 99  |
| 19) Hexachloroethane          | 8.825  | 117 | 17258  | 5.423 | ng/u1 95  |
| 22) Nitrobenzene              | 8.943  | 77  | 39745  | 4.984 | ng/u1 98  |
| 23) Isophorone                | 9.466  | 82  | 87227  | 5.286 | ng/u1 100 |
| 25) 2-Nitrophenol             | 9.649  | 139 | 19233  | 5.314 | ng/u1 94  |
| 26) 2,4-Dimethylphenol        | 9.719  | 107 | 42179  | 5.063 | ng/u1 91  |
| 27) Bis(2-Chloroethoxy)met... | 9.943  | 93  | 47648  | 4.732 | ng/u1 98  |
| 29) 2,4-Dichlorophenol        | 10.207 | 162 | 32602  | 4.407 | ng/u1 96  |
| 30) Naphthalene               | 10.578 | 128 | 126526 | 4.886 | ng/u1 99  |
| 33) Hexachlorobutadiene       | 10.860 | 225 | 27890  | 5.515 | ng/u1 99  |
| 35) 4-Chloro-3-methylphenol   | 11.831 | 107 | 39850  | 5.184 | ng/u1 98  |
| 36) 2-Methylnaphthalene       | 12.190 | 142 | 88330  | 4.764 | ng/u1 96  |
| 37) 1-Methylnaphthalene       | 12.401 | 142 | 88456  | 4.891 | ng/u1 99  |
| 39) 1,2,4,5-Tetrachloroben... | 12.560 | 216 | 48726  | 4.619 | ng/u1 97  |
| 41) 2,4,6-Trichlorophenol     | 12.801 | 196 | 30141m | 4.736 | ng/u1     |
| 42) 2,4,5-Trichlorophenol     | 12.901 | 196 | 34581m | 4.901 | ng/u1     |
| 43) 1,1'-Biphenyl             | 13.201 | 154 | 118698 | 4.548 | ng/u1 99  |
| 44) 2-Chloronaphthalene       | 13.243 | 162 | 90049  | 4.423 | ng/u1 94  |
| 45) 2-Nitroaniline            | 13.460 | 65  | 26052  | 6.015 | ng/u1 98  |
| 47) Dimethylphthalate         | 13.825 | 163 | 131635 | 5.182 | ng/u1 99  |
| 48) 2,6-Dinitrotoluene        | 13.954 | 165 | 23396  | 5.579 | ng/u1 96  |
| 50) Acenaphthylene            | 14.095 | 152 | 158658 | 4.712 | ng/u1 99  |

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|-------------------------------|--------|------|----------|-------|--------|----------|
| -----                         |        |      |          |       |        |          |
| 52) Acenaphthene              | 14.437 | 153  | 103088   | 4.760 | ng/u1  | 99       |
| 56) Dibenzofuran              | 14.784 | 168  | 148985   | 4.831 | ng/u1  | 100      |
| 57) 2,4-Dinitrotoluene        | 14.754 | 165  | 34255    | 5.512 | ng/u1  | 97       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.031 | 232  | 30579    | 5.283 | ng/u1# | 97       |
| 59) Diethylphthalate          | 15.207 | 149  | 135279   | 5.393 | ng/u1  | 99       |
| 61) Fluorene                  | 15.442 | 166  | 123777   | 4.961 | ng/u1  | 98       |
| 62) 4-Chlorophenyl-phenyle... | 15.442 | 204  | 64396    | 5.180 | ng/u1  | 94       |
| 67) N-Nitrosodiphenylamine    | 15.654 | 169  | 104502   | 4.564 | ng/u1  | 97       |
| 68) 4-Bromophenyl-phenylether | 16.354 | 248  | 38552    | 4.716 | ng/u1  | 97       |
| 69) Hexachlorobenzene         | 16.478 | 284  | 45271    | 4.475 | ng/u1  | 98       |
| 72) Phenanthrene              | 17.225 | 178  | 208086   | 4.820 | ng/u1  | 99       |
| 74) Anthracene                | 17.319 | 178  | 208183   | 4.801 | ng/u1  | 99       |
| 75) 1,2,3,4-Tetrachloroben... | 13.172 | 216  | 51651    | 4.439 | ng/uL  | 96       |
| 76) Pentachlorobenzene        | 14.707 | 250  | 52190    | 4.452 | ng/uL  | 98       |
| 78) Di-n-butylphthalate       | 18.189 | 149  | 242046   | 5.480 | ng/u1  | 99       |
| 80) Fluoranthene              | 19.307 | 202  | 259098   | 4.964 | ng/u1  | 99       |
| 82) Pyrene                    | 19.683 | 202  | 278626   | 5.042 | ng/u1  | 99       |
| 83) Butylbenzylphthalate      | 20.630 | 149  | 123234   | 6.899 | ng/u1  | 95       |
| 85) Benzo(a)anthracene        | 21.601 | 228  | 273601   | 4.895 | ng/u1  | 99       |
| 86) Bis(2-ethylhexyl)phtha... | 21.536 | 149  | 182378   | 6.326 | ng/u1  | 97       |
| 87) Chrysene                  | 21.666 | 228  | 259881   | 4.968 | ng/u1  | 99       |
| 90) Benzo(b)fluoranthene      | 23.913 | 252  | 257177   | 4.631 | ng/u1  | 99       |
| 91) Benzo(k)fluoranthene      | 23.989 | 252  | 276208   | 4.951 | ng/u1  | 98       |
| 93) Benzo(a)pyrene            | 24.824 | 252  | 256031   | 4.892 | ng/u1  | 98       |
| 94) Indeno(1,2,3-cd)pyrene    | 28.806 | 276  | 309705m  | 4.672 | ng/u1  |          |
| 95) Dibenzo(a,h)anthracene    | 28.871 | 278  | 240130   | 4.474 | ng/u1  | 99       |
| 96) Benzo(g,h,i)perylene      | 29.977 | 276  | 253885m  | 4.704 | ng/u1  |          |
| -----                         |        |      |          |       |        |          |

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

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