

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112423\
 Data File : BP018347.D
 Acq On : 25 Nov 2023 02:22
 Operator : MA/JU
 Sample : 05336-11
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCKE0

Quant Time: Nov 25 02:54:52 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 24 22:02:00 2023
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| Internal Standards | | | | | | |
| 1) 1,4-Dichlorobenzene-d4 | 7.869 | 152 | 371027 | 20.000 | ng/ul | 0.00 |
| 20) Naphthalene-d8 | 10.669 | 136 | 1482160 | 20.000 | ng/ul | 0.00 |
| 38) Acenaphthene-d10 | 14.498 | 164 | 845271 | 20.000 | ng/ul | 0.00 |
| 64) Phenanthrene-d10 | 17.245 | 188 | 1485241 | 20.000 | ng/ul | 0.00 |
| 79) Chrysene-d12 | 21.351 | 240 | 1125977 | 20.000 | ng/ul | 0.00 |
| 88) Perylene-d12 | 23.786 | 264 | 1394328 | 20.000 | ng/ul | 0.00 |

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| System Monitoring Compounds | | | | | | |
| 3) 1,4-Dioxane-d8 | 3.269 | 96 | 37334 | 3.769 | ng/uL | 0.00 |
| 4) Pyridine-d5 | 3.687 | 84 | 248419 | 9.125 | ng/ul | 0.00 |
| 7) Phenol-d5 | 7.028 | 99 | 180003 | 5.309 | ng/ul | 0.00 |
| 9) Bis-(2-Chloroethyl)eth... | 7.205 | 67 | 567761 | 28.001 | ng/ul | 0.00 |
| 11) 2-Chlorophenol-d4 | 7.393 | 132 | 578100 | 20.306 | ng/ul | 0.00 |
| 15) 4-Methylphenol-d8 | 8.575 | 113 | 353215 | 12.908 | ng/ul | 0.00 |
| 21) Nitrobenzene-d5 | 9.028 | 128 | 367165 | 30.499 | ng/ul | 0.00 |
| 24) 2-Nitrophenol-d4 | 9.752 | 143 | 324470 | 29.184 | ng/ul | 0.00 |
| 28) 2,4-Dichlorophenol-d3 | 10.287 | 165 | 662745 | 24.708 | ng/ul | 0.00 |
| 31) 4-Chloroaniline-d4 | 10.804 | 131 | 885688 | 24.482 | ng/ul | 0.00 |
| 46) Dimethylphthalate-d6 | 13.916 | 166 | 2357549 | 31.345 | ng/ul | 0.00 |
| 49) Acenaphthylene-d8 | 14.192 | 160 | 2615680 | 30.521 | ng/ul | 0.00 |
| 54) 4-Nitrophenol-d4 | 14.681 | 143 | 33873 | 3.262 | ng/ul | 0.00 |
| 60) Fluorene-d10 | 15.492 | 176 | 1967156 | 31.270 | ng/ul | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.604 | 200 | 158306 | 25.038 | ng/ul | 0.00 |
| 73) Anthracene-d10 | 17.345 | 188 | 2819085 | 35.511 | ng/ul | 0.00 |
| 81) Pyrene-d10 | 19.586 | 212 | 2849942 | 31.802 | ng/ul | 0.00 |
| 92) Benzo(a)pyrene-d12 | 23.627 | 264 | 3002245 | 37.139 | ng/ul | 0.00 |

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|--------|------|----------|-------|--------|--------|
| 2) 1,4-Dioxane | 3.317 | 88 | 843 | 0.079 | ng/uL# | 37 |
| 5) Pyridine | 3.711 | 79 | 194 | 0.007 | ng/ul# | 1 |
| 6) Benzaldehyde | 6.999 | 77 | 155 | 0.009 | ng/ul# | 17 |
| 8) Phenol | 7.052 | 94 | 1344 | 0.040 | ng/ul# | 23 |
| 10) Bis(2-Chloroethyl)ether | 7.293 | 93 | 175 | 0.006 | ng/ul# | 80 |
| 12) 2-Chlorophenol | 7.516 | 128 | 62 | 0.002 | ng/ul | 94 |
| 13) 2-Methylphenol | 8.322 | 108 | 34 | 0.001 | ng/ul | 100 |
| 14) 2,2'-oxybis(1-Chloropr... | 8.393 | 45 | 178 | 0.005 | ng/ul# | 1 |
| 16) Acetophenone | 8.693 | 105 | 685 | 0.016 | ng/ul# | 1 |
| 17) N-Nitroso-di-n-propyla... | 8.699 | 70 | 72 | 0.003 | ng/ul# | 1 |
| 18) 4-Methylphenol | 8.581 | 108 | 118 | 0.004 | ng/ul | 96 |
| 19) Hexachloroethane | 8.946 | 117 | 143 | 0.012 | ng/ul# | 7 |
| 22) Nitrobenzene | 9.075 | 77 | 209 | 0.007 | ng/ul# | 48 |
| 23) Isophorone | 9.593 | 82 | 278 | 0.004 | ng/ul# | 33 |
| 25) 2-Nitrophenol | 9.804 | 139 | 28 | 0.002 | ng/ul# | 1 |
| 26) 2,4-Dimethylphenol | 9.816 | 107 | 61 | 0.002 | ng/ul | 86 |
| 27) Bis(2-Chloroethoxy)met... | 10.093 | 93 | 47 | 0.001 | ng/ul# | 22 |
| 29) 2,4-Dichlorophenol | 10.316 | 162 | 89 | 0.003 | ng/ul# | 1 |
| 30) Naphthalene | 10.728 | 128 | 241 | 0.003 | ng/ul# | 22 |
| 32) 4-Chloroaniline | 10.893 | 127 | 113 | 0.003 | ng/ul | 93 |
| 33) Hexachlorobutadiene | 10.998 | 225 | 34 | 0.002 | ng/ul# | 67 |
| 34) Caprolactam | 11.598 | 113 | 290 | 0.039 | ng/ul# | 87 |
| 35) 4-Chloro-3-methylphenol | 11.945 | 107 | 151 | 0.005 | ng/ul# | 57 |
| 36) 2-Methylnaphthalene | 12.287 | 142 | 34 | 0.001 | ng/ul | 91 |

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 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 24 22:02:00 2023
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-------------------------------|--------|------|----------|-------|--------|----------|
| 37) 1-Methylnaphthalene | 12.545 | 142 | 119 | 0.002 | ng/ul# | 88 |
| 39) 1,2,4,5-Tetrachloroben... | 12.681 | 216 | 26 | 0.001 | ng/ul# | 1 |
| 40) Hexachlorocyclopentadiene | 12.669 | 237 | 29 | 0.002 | ng/ul# | 44 |
| 41) 2,4,6-Trichlorophenol | 12.934 | 196 | 79 | 0.004 | ng/ul# | 65 |
| 42) 2,4,5-Trichlorophenol | 12.987 | 196 | 74 | 0.004 | ng/ul# | 52 |
| 43) 1,1'-Biphenyl | 13.340 | 154 | 567 | 0.008 | ng/ul# | 87 |
| 44) 2-Chloronaphthalene | 13.369 | 162 | 306 | 0.005 | ng/ul# | 88 |
| 45) 2-Nitroaniline | 13.581 | 65 | 28 | 0.002 | ng/ul# | 60 |
| 47) Dimethylphthalate | 13.963 | 163 | 1918 | 0.026 | ng/ul# | 93 |
| 48) 2,6-Dinitrotoluene | 14.010 | 165 | 85 | 0.007 | ng/ul | 99 |
| 50) Acenaphthylene | 14.222 | 152 | 1310 | 0.014 | ng/ul# | 87 |
| 51) 3-Nitroaniline | 14.387 | 138 | 25 | 0.002 | ng/ul# | 14 |
| 52) Acenaphthene | 14.563 | 153 | 1253 | 0.020 | ng/ul | 95 |
| 53) 2,4-Dinitrophenol | 14.598 | 184 | 2 | 0.000 | ng/ul# | 1 |
| 55) 4-Nitrophenol | 14.687 | 109 | 70 | 0.009 | ng/ul# | 1 |
| 56) Dibenzofuran | 14.892 | 168 | 1406 | 0.016 | ng/ul | 100 |
| 57) 2,4-Dinitrotoluene | 14.851 | 165 | 153 | 0.009 | ng/ul# | 87 |
| 58) 2,3,4,6-Tetrachlorophenol | 15.116 | 232 | 54 | 0.004 | ng/ul# | 1 |
| 59) Diethylphthalate | 15.322 | 149 | 2752 | 0.037 | ng/ul# | 66 |
| 61) Fluorene | 15.551 | 166 | 1006 | 0.014 | ng/ul# | 92 |
| 62) 4-Chlorophenyl-phenyle... | 15.557 | 204 | 260 | 0.007 | ng/ul# | 67 |
| 63) 4-Nitroaniline | 15.551 | 138 | 124 | 0.011 | ng/ul# | 2 |
| 66) 4,6-Dinitro-2-methylph... | 15.622 | 198 | 83 | 0.012 | ng/ul# | 1 |
| 67) N-Nitrosodiphenylamine | 15.751 | 169 | 314 | 0.006 | ng/ul# | 27 |
| 68) 4-Bromophenyl-phenylether | 16.433 | 248 | 56 | 0.003 | ng/ul# | 47 |
| 69) Hexachlorobenzene | 16.545 | 284 | 220 | 0.010 | ng/ul# | 67 |
| 70) Atrazine | 16.545 | 200 | 48 | 0.003 | ng/ul | 93 |
| 71) Pentachlorophenol | 16.892 | 266 | 49 | 0.005 | ng/ul# | 30 |
| 72) Phenanthrene | 17.292 | 178 | 1697 | 0.019 | ng/ul# | 91 |
| 74) Anthracene | 17.381 | 178 | 1847 | 0.020 | ng/ul# | 80 |
| 75) 1,2,3,4-Tetrachloroben... | 13.298 | 216 | 46 | 0.002 | ng/ul# | 56 |
| 76) Pentachlorobenzene | 14.816 | 250 | 113 | 0.005 | ng/ul# | 72 |
| 77) Carbazole | 17.657 | 167 | 1334 | 0.018 | ng/ul# | 82 |
| 78) Di-n-butylphthalate | 18.216 | 149 | 5686 | 0.055 | ng/ul# | 96 |
| 80) Fluoranthene | 19.257 | 202 | 2038 | 0.019 | ng/ul# | 83 |
| 82) Pyrene | 19.616 | 202 | 1947 | 0.018 | ng/ul# | 77 |
| 83) Butylbenzylphthalate | 20.510 | 149 | 658 | 0.016 | ng/ul# | 95 |
| 84) 3,3'-Dichlorobenzidine | 21.286 | 252 | 557 | 0.020 | ng/ul# | 71 |
| 85) Benzo(a)anthracene | 21.351 | 228 | 5098 | 0.056 | ng/ul# | 75 |
| 86) Bis(2-ethylhexyl)phtha... | 21.292 | 149 | 3609 | 0.060 | ng/ul# | 97 |
| 87) Chrysene | 21.386 | 228 | 3126 | 0.037 | ng/ul# | 88 |
| 89) Di-n-octyl phthalate | 22.245 | 149 | 483 | 0.005 | ng/ul | 100 |
| 90) Benzo(b)fluoranthene | 23.033 | 252 | 1891 | 0.019 | ng/ul# | 19 |
| 91) Benzo(k)fluoranthene | 23.086 | 252 | 902 | 0.009 | ng/ul# | 38 |
| 93) Benzo(a)pyrene | 23.663 | 252 | 314 | 0.003 | ng/ul# | 9 |
| 94) Indeno(1,2,3-cd)pyrene | 26.339 | 276 | 475 | 0.004 | ng/ul# | 77 |
| 95) Dibenzo(a,h)anthracene | 26.386 | 278 | 1028 | 0.012 | ng/ul# | 75 |
| 96) Benzo(g,h,i)perylene | 27.127 | 276 | 161 | 0.002 | ng/ul# | 75 |

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

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