

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE110624\
 Data File : BE101498.D
 Acq On : 6 Nov 2024 16:50
 Operator : RC/JU
 Sample : SSTDICC050
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
ClientSampleId :
 SSTDICC050

Manual Integrations
APPROVED
 Reviewed By :Jagrut Upadhyay 11/07/2024
 Supervised By :mohammad ahmed 11/08/2024

Quant Time: Nov 06 23:53:13 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_E\Methods\8270-BE110624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 06 23:45:11 2024
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) | Qvalue |
|------------------------------------|--------|------|----------|---------|-------|----------|--------|
| Internal Standards | | | | | | | |
| 1) 1,4-Dichlorobenzene-d4 | 7.561 | 152 | 91095 | 20.000 | ng | 0.00 | |
| 21) Naphthalene-d8 | 10.334 | 136 | 420546 | 20.000 | ng | 0.00 | |
| 39) Acenaphthene-d10 | 14.177 | 164 | 294714 | 20.000 | ng | 0.00 | |
| 64) Phenanthrene-d10 | 16.915 | 188 | 651829 | 20.000 | ng | 0.00 | |
| 76) Chrysene-d12 | 21.075 | 240 | 698124 | 20.000 | ng | 0.00 | |
| 86) Perylene-d12 | 23.360 | 264 | 907393 | 20.000 | ng | 0.00 | |
| System Monitoring Compounds | | | | | | | |
| 5) 2-Fluorophenol | 5.317 | 112 | 507391 | 98.486 | ng | 0.00 | |
| 7) Phenol-d6 | 6.950 | 99 | 715571 | 101.303 | ng | 0.00 | |
| 23) Nitrobenzene-d5 | 8.777 | 82 | 669237 | 100.518 | ng | 0.00 | |
| 42) 2,4,6-Tribromophenol | 15.687 | 330 | 535740 | 96.606 | ng | 0.00 | |
| 45) 2-Fluorobiphenyl | 12.802 | 172 | 1709745 | 94.725 | ng | 0.00 | |
| 79) Terphenyl-d14 | 19.506 | 244 | 2857301 | 90.597 | ng | 0.00 | |
| Target Compounds | | | | | | | |
| 2) 1,4-Dioxane | 3.190 | 88 | 86138 | 44.928 | ng | | 99 |
| 3) Pyridine | 3.613 | 79 | 271363m | 50.734 | ng | | |
| 4) n-Nitrosodimethylamine | 3.537 | 42 | 104883 | 49.532 | ng | | 97 |
| 6) Aniline | 7.003 | 93 | 233770 | 44.094 | ng | | 100 |
| 8) 2-Chlorophenol | 7.209 | 128 | 304216 | 49.841 | ng | | 99 |
| 9) Benzaldehyde | 6.768 | 77 | 149018 | 43.128 | ng | | 98 |
| 10) Phenol | 6.974 | 94 | 389135 | 50.610 | ng | | 99 |
| 11) bis(2-Chloroethyl)ether | 7.032 | 93 | 329221 | 51.732 | ng | | 98 |
| 12) 1,3-Dichlorobenzene | 7.450 | 146 | 321129 | 48.192 | ng | | 99 |
| 13) 1,4-Dichlorobenzene | 7.596 | 146 | 327207 | 48.480 | ng | | 99 |
| 14) 1,2-Dichlorobenzene | 7.926 | 146 | 322412 | 48.664 | ng | | 99 |
| 15) Benzyl Alcohol | 7.902 | 79 | 218205 | 52.883 | ng | | 98 |
| 16) 2,2'-oxybis(1-Chloropr... | 8.072 | 45 | 369966 | 49.103 | ng | | 99 |
| 17) 2-Methylphenol | 8.155 | 107 | 255685 | 51.364 | ng | | 98 |
| 18) Hexachloroethane | 8.619 | 117 | 112465 | 49.341 | ng | | 98 |
| 19) n-Nitroso-di-n-propyla... | 8.372 | 70 | 238326 | 51.171 | ng | | 99 |
| 20) 3+4-Methylphenols | 8.490 | 107 | 354774 | 51.400 | ng | | 99 |
| 22) Acetophenone | 8.419 | 105 | 469557 | 49.302 | ng | # | 99 |
| 24) Nitrobenzene | 8.819 | 77 | 345554 | 49.900 | ng | | 100 |
| 25) Isophorone | 9.271 | 82 | 649442 | 50.124 | ng | | 100 |
| 26) 2-Nitrophenol | 9.500 | 139 | 190923 | 51.807 | ng | | 99 |
| 27) 2,4-Dimethylphenol | 9.624 | 122 | 215719 | 50.589 | ng | | 99 |
| 28) bis(2-Chloroethoxy)met... | 9.765 | 93 | 394162 | 49.697 | ng | | 100 |
| 29) 2,4-Dichlorophenol | 10.058 | 162 | 308633 | 50.986 | ng | | 99 |
| 30) 1,2,4-Trichlorobenzene | 10.176 | 180 | 328828 | 48.661 | ng | | 99 |
| 31) Naphthalene | 10.381 | 128 | 1033552 | 48.682 | ng | | 100 |
| 32) Benzoic acid | 9.935 | 122 | 200828 | 50.078 | ng | | 99 |
| 33) 4-Chloroaniline | 10.611 | 127 | 374148 | 50.114 | ng | | 98 |
| 34) Hexachlorobutadiene | 10.628 | 225 | 205479 | 49.334 | ng | | 98 |
| 35) Caprolactam | 11.421 | 113 | 111070m | 50.028 | ng | | |
| 36) 4-Chloro-3-methylphenol | 11.792 | 107 | 334516 | 50.606 | ng | | 99 |
| 37) 2-Methylnaphthalene | 11.974 | 142 | 741321 | 49.161 | ng | | 98 |
| 38) 1-Methylnaphthalene | 12.197 | 142 | 733235 | 48.758 | ng | | 99 |
| 40) 1,2,4,5-Tetrachloroben... | 12.332 | 216 | 382507 | 49.228 | ng | | 100 |
| 41) Hexachlorocyclopentadiene | 12.309 | 237 | 128897 | 56.847 | ng | | 98 |
| 43) 2,4,6-Trichlorophenol | 12.649 | 196 | 265972 | 49.842 | ng | | 99 |

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE110624\
 Data File : BE101498.D
 Acq On : 6 Nov 2024 16:50
 Operator : RC/JU
 Sample : SSTDICC050
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDICC050

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 11/07/2024
 Supervised By :mohammad ahmed 11/08/2024

Quant Time: Nov 06 23:53:13 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_E\Methods\8270-BE110624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 06 23:45:11 2024
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 44) 2,4,5-Trichlorophenol | 12.755 | 196 | 302473 | 50.879 | ng | 98 |
| 46) 1,1'-Biphenyl | 13.002 | 154 | 990577 | 48.727 | ng | 99 |
| 47) 2-Chloronaphthalene | 13.049 | 162 | 784561 | 48.931 | ng | 99 |
| 48) 2-Nitroaniline | 13.384 | 65 | 220977 | 52.065 | ng | 99 |
| 49) Acenaphthylene | 13.907 | 152 | 1168645 | 48.909 | ng | 99 |
| 50) Dimethylphthalate | 13.689 | 163 | 1007892 | 48.026 | ng | 99 |
| 51) 2,6-Dinitrotoluene | 13.824 | 165 | 241242 | 49.804 | ng | 98 |
| 52) Acenaphthene | 14.242 | 154 | 733889 | 48.131 | ng | 100 |
| 53) 3-Nitroaniline | 14.230 | 138 | 238076 | 50.629 | ng | 97 |
| 54) 2,4-Dinitrophenol | 14.371 | 184 | 153156 | 53.956 | ng | 98 |
| 55) Dibenzofuran | 14.582 | 168 | 1172466 | 47.751 | ng | 98 |
| 56) 4-Nitrophenol | 14.659 | 139 | 201601 | 51.919 | ng | 98 |
| 57) 2,4-Dinitrotoluene | 14.606 | 165 | 340766 | 50.076 | ng | 99 |
| 58) Fluorene | 15.223 | 166 | 978319 | 47.743 | ng | 99 |
| 59) 2,3,4,6-Tetrachlorophenol | 14.853 | 232 | 272447 | 49.527 | ng | 99 |
| 60) Diethylphthalate | 15.011 | 149 | 1062117 | 47.998 | ng | 99 |
| 61) 4-Chlorophenyl-phenyle... | 15.205 | 204 | 500724 | 48.121 | ng | 99 |
| 62) 4-Nitroaniline | 15.423 | 138 | 261365 | 51.589 | ng | 99 |
| 63) Azobenzene | 15.505 | 77 | 871780 | 48.291 | ng | 99 |
| 65) 4,6-Dinitro-2-methylph... | 15.352 | 198 | 214210 | 55.361 | ng | 99 |
| 66) n-Nitrosodiphenylamine | 15.464 | 169 | 857158 | 49.551 | ng | 99 |
| 67) 4-Bromophenyl-phenylether | 16.092 | 248 | 359667 | 50.029 | ng | 98 |
| 68) Hexachlorobenzene | 16.204 | 284 | 468408 | 49.506 | ng | 99 |
| 69) Atrazine | 16.392 | 200 | 197336 | 38.639 | ng | 99 |
| 70) Pentachlorophenol | 16.609 | 266 | 272040 | 52.991 | ng | 99 |
| 71) Phenanthrene | 16.956 | 178 | 1526029 | 48.135 | ng | 100 |
| 72) Anthracene | 17.044 | 178 | 1518459 | 48.501 | ng | 99 |
| 73) Carbazole | 17.373 | 167 | 1504975 | 47.935 | ng | 98 |
| 74) Di-n-butylphthalate | 17.820 | 149 | 1837620 | 47.396 | ng | 99 |
| 75) Fluoranthene | 18.960 | 202 | 1885849 | 46.395 | ng | 99 |
| 77) Benzidine | 19.224 | 184 | 1144629 | 73.321 | ng | 100 |
| 78) Pyrene | 19.330 | 202 | 1962930 | 49.416 | ng | 98 |
| 80) Butylbenzylphthalate | 20.176 | 149 | 873276 | 48.920 | ng | 98 |
| 81) Benzo(a)anthracene | 21.057 | 228 | 1985312 | 47.883 | ng | 98 |
| 82) 3,3'-Dichlorobenzidine | 21.034 | 252 | 833578 | 51.072 | ng | 99 |
| 83) Chrysene | 21.116 | 228 | 1875257 | 47.584 | ng | 99 |
| 84) Bis(2-ethylhexyl)phtha... | 20.881 | 149 | 1287632 | 47.710 | ng | # 97 |
| 85) Di-n-octyl phthalate | 21.727 | 149 | 2140362 | 46.878 | ng | 100 |
| 87) Indeno(1,2,3-cd)pyrene | 25.711 | 276 | 3044286 | 48.821 | ng | 100 |
| 88) Benzo(b)fluoranthene | 22.655 | 252 | 2388668 | 47.982 | ng | 99 |
| 89) Benzo(k)fluoranthene | 22.702 | 252 | 2191941 | 48.502 | ng | 99 |
| 90) Benzo(a)pyrene | 23.260 | 252 | 2093708 | 48.714 | ng | 99 |
| 91) Dibenzo(a,h)anthracene | 25.699 | 278 | 2519176 | 48.510 | ng | 100 |
| 92) Benzo(g,h,i)perylene | 26.457 | 276 | 2604721 | 49.369 | ng | 100 |

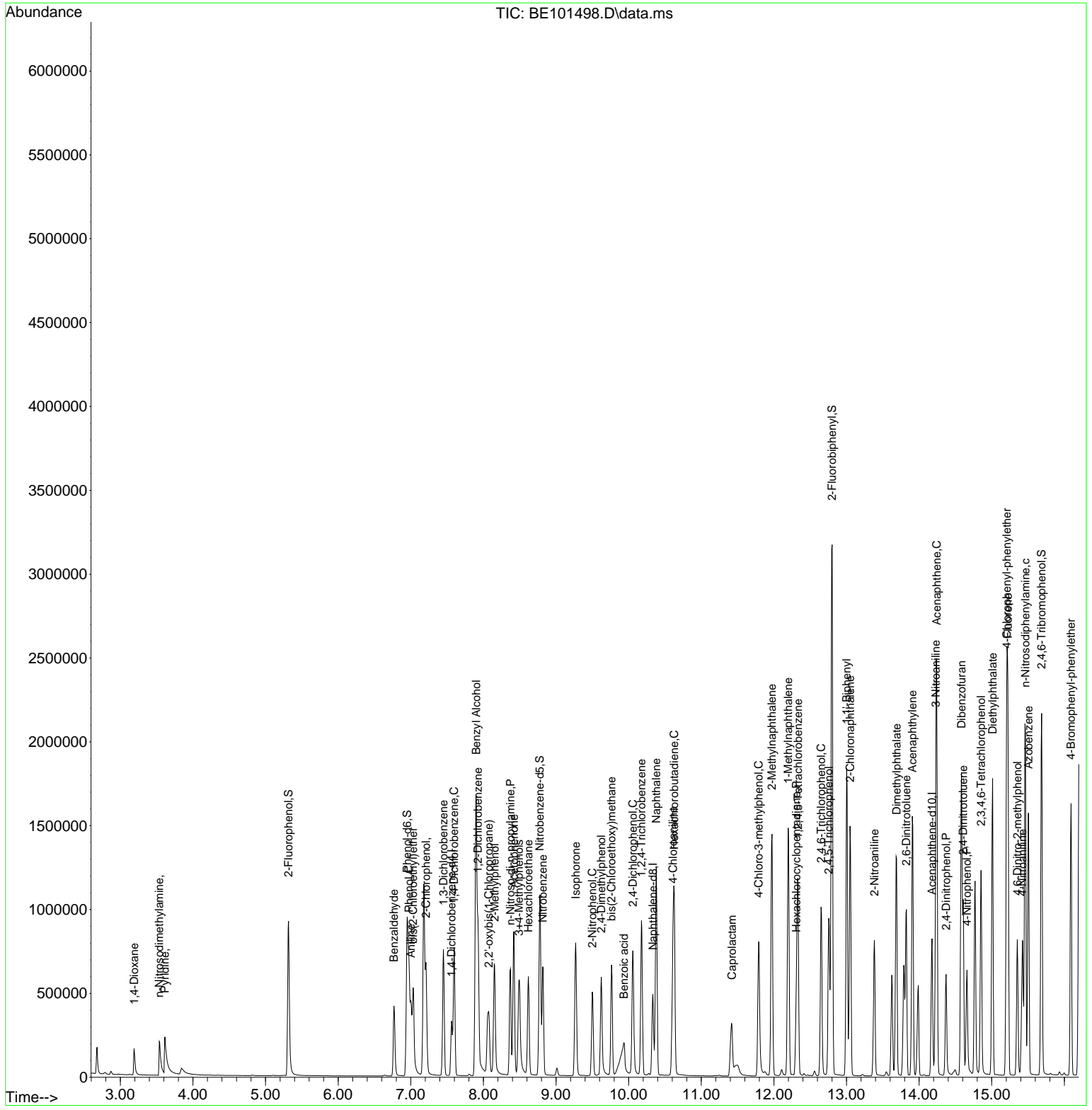
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE110624\
 Data File : BE101498.D
 Acq On : 6 Nov 2024 16:50
 Operator : RC/JU
 Sample : SSTDICC050
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
ClientSampleId :
 SSTDICC050

Quant Time: Nov 06 23:53:13 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_E\Methods\8270-BE110624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 06 23:45:11 2024
 Response via : Initial Calibration

Manual Integrations
APPROVED
 Reviewed By :Jagrut Upadhyay 11/07/2024
 Supervised By :mohammad ahmed 11/08/2024



Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE110624\
 Data File : BE101498.D
 Acq On : 6 Nov 2024 16:50
 Operator : RC/JU
 Sample : SSTDICC050
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :

BNA_E

ClientSampleId :

SSTDICC050

Manual Integrations

APPROVED

Reviewed By :Jagrut Upadhyay 11/07/2024
 Supervised By :mohammad ahmed 11/08/2024

Quant Time: Nov 06 23:53:13 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_E\Methods\8270-BE110624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 06 23:45:11 2024
 Response via : Initial Calibration

