

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG113023\  
 Data File : BG059918.D  
 Acq On : 30 Nov 2023 20:38  
 Operator : MA/JU  
 Sample : SSTDICC050  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SSTDICC050

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 12/01/2023  
 Supervised By :mohammad ahmed 12/01/2023

Quant Time: Dec 01 03:30:35 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG113023.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Dec 01 03:17:30 2023  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc Units	Dev(Min)	
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.967	152	30637	20.000 ng	0.00	
21) Naphthalene-d8	10.769	136	144988	20.000 ng	0.00	
39) Acenaphthene-d10	14.606	164	102868	20.000 ng	0.00	
64) Phenanthrene-d10	17.350	188	234541	20.000 ng	0.00	
76) Chrysene-d12	21.627	240	216389	20.000 ng	0.00	
86) Perylene-d12	24.829	264	258997	20.000 ng	0.00	
System Monitoring Compounds						
5) 2-Fluorophenol	5.528	112	187930	96.084 ng	0.00	
7) Phenol-d6	7.115	99	273565	96.898 ng	0.00	
23) Nitrobenzene-d5	9.130	82	254752	102.186 ng	0.00	
42) 2,4,6-Tribromophenol	16.092	330	153018	96.926 ng	0.00	
45) 2-Fluorobiphenyl	13.231	172	711190	96.163 ng	0.00	
79) Terphenyl-d14	19.976	244	1195422	91.863 ng	0.00	
Target Compounds						
2) 1,4-Dioxane	3.442	88	24699m	40.499 ng		Qvalue
3) Pyridine	3.830	79	71238m	45.329 ng		
4) n-Nitrosodimethylamine	3.748	42	42114	46.859 ng	#	97
6) Aniline	7.291	93	146571	48.456 ng		99
8) 2-Chlorophenol	7.526	128	104427	47.398 ng		92
9) Benzaldehyde	7.097	77	72328	40.396 ng		95
10) Phenol	7.144	94	141279	47.801 ng		96
11) bis(2-Chloroethyl)ether	7.385	93	110890	48.665 ng		98
12) 1,3-Dichlorobenzene	7.855	146	116074	48.097 ng		95
13) 1,4-Dichlorobenzene	8.002	146	118699	47.729 ng		97
14) 1,2-Dichlorobenzene	8.319	146	115310	47.059 ng		98
15) Benzyl Alcohol	8.196	79	105662	48.417 ng		94
16) 2,2'-oxybis(1-Chloropr...	8.495	45	207374	47.760 ng		97
17) 2-Methylphenol	8.396	107	98537	48.154 ng		97
18) Hexachloroethane	9.054	117	42073	47.091 ng		99
19) n-Nitroso-di-n-propyla...	8.772	70	96611	48.426 ng		99
20) 3+4-Methylphenols	8.730	107	139086	48.249 ng		96
22) Acetophenone	8.789	105	191696	47.758 ng		99
24) Nitrobenzene	9.171	77	137695	50.714 ng		97
25) Isophorone	9.694	82	285904	47.553 ng		99
26) 2-Nitrophenol	9.876	139	50257	57.082 ng		97
27) 2,4-Dimethylphenol	9.935	122	82741	47.462 ng		94
28) bis(2-Chloroethoxy)met...	10.176	93	161615	47.482 ng		94
29) 2,4-Dichlorophenol	10.411	162	114809	48.242 ng		98
30) 1,2,4-Trichlorobenzene	10.628	180	132059	47.594 ng		96
31) Naphthalene	10.822	128	389793	47.656 ng		98
32) Benzoic acid	10.064	122	77996	47.316 ng		97
33) 4-Chloroaniline	10.922	127	159915	48.269 ng		98
34) Hexachlorobutadiene	11.110	225	88761	47.686 ng		97
35) Caprolactam	11.703	113	39306	46.480 ng		93
36) 4-Chloro-3-methylphenol	12.044	107	136244	48.026 ng		98
37) 2-Methylnaphthalene	12.426	142	272896	47.135 ng		98
38) 1-Methylnaphthalene	12.649	142	272253	46.474 ng		94
40) 1,2,4,5-Tetrachloroben...	12.796	216	159273	49.688 ng		99
41) Hexachlorocyclopentadiene	12.779	237	63306	51.842 ng		98
43) 2,4,6-Trichlorophenol	13.031	196	103780	49.898 ng		98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.102	196	117225	49.729	ng	99
46) 1,1'-Biphenyl	13.443	154	368961	47.818	ng	98
47) 2-Chloronaphthalene	13.478	162	310529	48.672	ng	98
48) 2-Nitroaniline	13.678	65	91398	54.265	ng	99
49) Acenaphthylene	14.324	152	468904	47.848	ng	98
50) Dimethylphthalate	14.059	163	393863	47.722	ng	99
51) 2,6-Dinitrotoluene	14.177	165	76472	48.846	ng	97
52) Acenaphthene	14.671	154	293253	48.170	ng	99
53) 3-Nitroaniline	14.506	138	75789	52.598	ng	97
54) 2,4-Dinitrophenol	14.706	184	29566	47.373	ng	98
55) Dibenzofuran	15.000	168	446218	47.100	ng	100
56) 4-Nitrophenol	14.800	139	61619	52.584	ng	98
57) 2,4-Dinitrotoluene	14.958	165	98988	47.623	ng	98
58) Fluorene	15.652	166	353823	46.061	ng	96
59) 2,3,4,6-Tetrachlorophenol	15.223	232	103112	48.123	ng	98
60) Diethylphthalate	15.428	149	382766	45.988	ng	97
61) 4-Chlorophenyl-phenyle...	15.646	204	195557	47.191	ng	98
62) 4-Nitroaniline	15.669	138	76327	49.490	ng	95
63) Azobenzene	15.940	77	367468	46.968	ng	99
65) 4,6-Dinitro-2-methylph...	15.722	198	45136	46.975	ng	98
66) n-Nitrosodiphenylamine	15.863	169	309383	48.964	ng	99
67) 4-Bromophenyl-phenylether	16.545	248	136667	48.787	ng	96
68) Hexachlorobenzene	16.651	284	161220	48.379	ng	99
69) Atrazine	16.809	200	106120	45.804	ng	97
70) Pentachlorophenol	16.997	266	98148	52.067	ng	97
71) Phenanthrene	17.397	178	559929	47.057	ng	99
72) Anthracene	17.485	178	567723	47.299	ng	100
73) Carbazole	17.755	167	520380	46.760	ng	99
74) Di-n-butylphthalate	18.325	149	634279	46.795	ng	99
75) Fluoranthene	19.406	202	707272	46.698	ng	99
77) Benzidine	19.588	184	248752	57.269	ng	98
78) Pyrene	19.770	202	729289	47.251	ng	99
80) Butylbenzylphthalate	20.669	149	265373	48.556	ng	96
81) Benzo(a)anthracene	21.609	228	713514	47.287	ng	99
82) 3,3'-Dichlorobenzidine	21.527	252	235008	47.806	ng	98
83) Chrysene	21.674	228	685632	47.912	ng	98
84) Bis(2-ethylhexyl)phtha...	21.533	149	391109	48.243	ng	99
85) Di-n-octyl phthalate	22.749	149	683599	49.118	ng	99
87) Indeno(1,2,3-cd)pyrene	28.484	276	938100	48.401	ng	99
88) Benzo(b)fluoranthene	23.813	252	782460	48.325	ng	100
89) Benzo(k)fluoranthene	23.883	252	753114	47.546	ng	98
90) Benzo(a)pyrene	24.676	252	694604	48.262	ng	98
91) Dibenzo(a,h)anthracene	28.560	278	790631	48.923	ng	97
92) Benzo(g,h,i)perylene	29.629	276	779923	48.330	ng	99

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

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