

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071422\
 Data File : BG054284.D
 Acq On : 14 Jul 2022 17:38
 Operator : CG/JU
 Sample : N3711-07MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PSC124055MS

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 07/15/2022
 Supervised By :Jagrut Upadhyay 07/15/2022

Quant Time: Jul 15 01:58:24 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 12 15:49:20 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.041	152	15279	20.000	ng	0.00
21) Naphthalene-d8	10.850	136	58245	20.000	ng	# 0.00
39) Acenaphthene-d10	14.669	164	46668	20.000	ng	0.00
64) Phenanthrene-d10	17.413	188	127354	20.000	ng	# 0.00
76) Chrysene-d12	21.684	240	157023	20.000	ng	# 0.00
86) Perylene-d12	24.916	264	172756	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.609	112	71215	93.643	ng	0.00
7) Phenol-d6	7.207	99	91969	87.681	ng	0.00
23) Nitrobenzene-d5	9.199	82	57799	53.851	ng	0.00
42) 2,4,6-Tribromophenol	16.155	330	91637	101.368	ng	0.00
45) 2-Fluorobiphenyl	13.288	172	174815	55.041	ng	0.00
79) Terphenyl-d14	20.016	244	430533	56.383	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.482	88	12034	36.183	ng	96
3) Pyridine	3.882	79	31072	37.407	ng	96
4) n-Nitrosodimethylamine	3.793	42	22634	53.936	ng	# 71
6) Aniline	7.360	93	53741	43.586	ng	96
8) 2-Chlorophenol	7.607	128	45565	53.456	ng	92
9) Benzaldehyde	7.172	77	23163m	36.774	ng	
10) Phenol	7.236	94	52691	49.790	ng	89
11) bis(2-Chloroethyl)ether	7.454	93	35128	45.011	ng	91
12) 1,3-Dichlorobenzene	7.930	146	51765	50.015	ng	91
13) 1,4-Dichlorobenzene	8.077	146	54045	50.957	ng	98
14) 1,2-Dichlorobenzene	8.394	146	50973	50.052	ng	99
15) Benzyl Alcohol	8.276	79	42706	50.505	ng	96
16) 2,2'-oxybis(1-Chloropr...	8.570	45	42031	38.785	ng	89
17) 2-Methylphenol	8.482	107	36986	48.916	ng	97
18) Hexachloroethane	9.128	117	17632	48.862	ng	# 60
19) n-Nitroso-di-n-propyla...	8.840	70	31162	43.517	ng	# 91
20) 3+4-Methylphenols	8.811	107	49413	46.576	ng	98
22) Acetophenone	8.858	105	66756	47.425	ng	# 96
24) Nitrobenzene	9.246	77	48849	46.295	ng	# 84
25) Isophorone	9.769	82	88838	45.755	ng	# 89
26) 2-Nitrophenol	9.957	139	25864	49.754	ng	92
27) 2,4-Dimethylphenol	10.016	122	41751	57.418	ng	95
28) bis(2-Chloroethoxy)met...	10.245	93	48829	49.404	ng	97
29) 2,4-Dichlorophenol	10.497	162	56718	53.818	ng	90
30) 1,2,4-Trichlorobenzene	10.715	180	66232	50.936	ng	# 97
31) Naphthalene	10.897	128	135897	46.398	ng	99
32) Benzoic acid	10.192	122	11953m	36.554	ng	
33) 4-Chloroaniline	11.003	127	43255	35.512	ng	94
34) Hexachlorobutadiene	11.185	225	59590	55.231	ng	95
35) Caprolactam	11.772	113	13559m	45.176	ng	
36) 4-Chloro-3-methylphenol	12.131	107	49884	49.199	ng	86
37) 2-Methylnaphthalene	12.501	142	106507	47.450	ng	99
38) 1-Methylnaphthalene	12.718	142	101043	46.366	ng	98
40) 1,2,4,5-Tetrachloroben...	12.871	216	100723	53.171	ng	# 95
41) Hexachlorocyclopentadiene	12.848	237	82028	100.450	ng	95
43) 2,4,6-Trichlorophenol	13.112	196	56355	51.800	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.188	196	62730	52.713	ng	# 91
46) 1,1'-Biphenyl	13.506	154	157174	50.083	ng	96
47) 2-Chloronaphthalene	13.547	162	127957	48.773	ng	96
48) 2-Nitroaniline	13.746	65	33454	45.055	ng	89
49) Acenaphthylene	14.393	152	190064	47.755	ng	99
50) Dimethylphthalate	14.117	163	174849	47.719	ng	99
51) 2,6-Dinitrotoluene	14.240	165	39073	49.061	ng	# 76
52) Acenaphthene	14.734	154	116172	48.258	ng	99
53) 3-Nitroaniline	14.569	138	26339	37.310	ng	87
54) 2,4-Dinitrophenol	14.786	184	41101	92.574	ng	# 78
55) Dibenzofuran	15.063	168	209221	49.481	ng	96
56) 4-Nitrophenol	14.892	139	47936	105.695	ng	94
57) 2,4-Dinitrotoluene	15.033	165	56574	49.058	ng	# 82
58) Fluorene	15.715	166	171740	48.893	ng	96
59) 2,3,4,6-Tetrachlorophenol	15.298	232	60721	50.943	ng	# 86
60) Diethylphthalate	15.480	149	170191	47.622	ng	94
61) 4-Chlorophenyl-phenyle...	15.703	204	118405	53.068	ng	# 86
62) 4-Nitroaniline	15.732	138	33296	45.157	ng	92
63) Azobenzene	15.997	77	117371	43.579	ng	81
65) 4,6-Dinitro-2-methylph...	15.797	198	35674	47.616	ng	81
66) n-Nitrosodiphenylamine	15.915	169	156191	48.623	ng	95
67) 4-Bromophenyl-phenylether	16.596	248	88729	52.236	ng	# 90
68) Hexachlorobenzene	16.725	284	103812	53.544	ng	# 89
69) Atrazine	16.860	200	80482	56.728	ng	93
70) Pentachlorophenol	17.072	266	80690	91.074	ng	98
71) Phenanthrene	17.454	178	304866	47.855	ng	97
72) Anthracene	17.548	178	307626	49.308	ng	99
73) Carbazole	17.812	167	257384	48.716	ng	98
74) Di-n-butylphthalate	18.365	149	288573	46.126	ng	98
75) Fluoranthene	19.463	202	429180	49.538	ng	95
77) Benzidine	19.640	184	247772	89.889	ng	98
78) Pyrene	19.828	202	427367	47.291	ng	99
80) Butylbenzylphthalate	20.703	149	128714	43.958	ng	# 84
81) Benzo(a)anthracene	21.667	228	478460	48.169	ng	98
82) 3,3'-Dichlorobenzidine	21.573	252	124538	41.022	ng	# 97
83) Chrysene	21.731	228	442090	47.058	ng	97
84) Bis(2-ethylhexyl)phtha...	21.567	149	184366	44.378	ng	# 92
85) Di-n-octyl phthalate	22.771	149	317216	44.352	ng	# 92
87) Indeno(1,2,3-cd)pyrene	28.635	276	615172	47.334	ng	# 98
88) Benzo(b)fluoranthene	23.893	252	532546	51.985	ng	# 97
89) Benzo(k)fluoranthene	23.958	252	486129	47.833	ng	# 97
90) Benzo(a)pyrene	24.769	252	459331	52.403	ng	# 98
91) Dibenzo(a,h)anthracene	28.682	278	530770	49.852	ng	# 95
92) Benzo(g,h,i)perylene	29.786	276	530021	49.911	ng	# 92

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

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