

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG011123\
 Data File : BG056253.D
 Acq On : 11 Jan 2023 16:40
 Operator : CG/JU
 Sample : 01064-03MS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 S4-03MS

Quant Time: Jan 12 04:21:09 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 12 03:38:40 2023
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By : Christian Giraldo 01/12/2023
 Supervised By : Jagrut Upadhyay 01/12/2023

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.318	152	43850	20.000	ng	0.00
21) Naphthalene-d8	11.150	136	156975	20.000	ng	# 0.00
39) Acenaphthene-d10	14.928	164	124758	20.000	ng	0.00
64) Phenanthrene-d10	17.666	188	338358	20.000	ng	0.00
76) Chrysene-d12	21.961	240	335768	20.000	ng	0.00
86) Perylene-d12	25.398	264	373661	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.833	112	358181	145.845	ng	0.00
7) Phenol-d6	7.454	99	501452	132.956	ng	0.00
23) Nitrobenzene-d5	9.493	82	259394	84.526	ng	0.00
42) 2,4,6-Tribromophenol	16.409	330	226977	143.653	ng	0.00
45) 2-Fluorobiphenyl	13.559	172	765639	84.746	ng	0.00
79) Terphenyl-d14	20.239	244	1638771	87.415	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.647	88	50557	45.070	ng	# 93
3) Pyridine	4.064	79	134624	39.403	ng	94
4) n-Nitrosodimethylamine	3.970	42	32043	46.105	ng	99
6) Aniline	7.631	93	84602	17.105	ng	99
8) 2-Chlorophenol	7.877	128	129209	52.871	ng	98
9) Benzaldehyde	7.437	77	65753	29.321	ng	95
10) Phenol	7.478	94	193129	50.493	ng	99
11) bis(2-Chloroethyl)ether	7.719	93	119925	46.267	ng	97
12) 1,3-Dichlorobenzene	8.206	146	150879	50.727	ng	96
13) 1,4-Dichlorobenzene	8.353	146	149223	49.961	ng	98
14) 1,2-Dichlorobenzene	8.676	146	143408	49.761	ng	98
15) Benzyl Alcohol	8.547	79	123301	44.920	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.829	45	23396	51.078	ng	92
17) 2-Methylphenol	8.747	107	129518	46.148	ng	98
18) Hexachloroethane	9.417	117	45024	49.621	ng	94
19) n-Nitroso-di-n-propyla...	9.117	70	92642	40.394	ng	99
20) 3+4-Methylphenols	9.076	107	176115	44.632	ng	98
22) Acetophenone	9.141	105	204185	46.003	ng	# 99
24) Nitrobenzene	9.534	77	142003	46.693	ng	99
25) Isophorone	10.051	82	276589	42.836	ng	97
26) 2-Nitrophenol	10.251	139	77862	50.039	ng	97
27) 2,4-Dimethylphenol	10.292	122	140130	49.375	ng	99
28) bis(2-Chloroethoxy)met...	10.527	93	159380	47.492	ng	99
29) 2,4-Dichlorophenol	10.786	162	162571	51.813	ng	99
30) 1,2,4-Trichlorobenzene	11.003	180	175338	49.243	ng	96
31) Naphthalene	11.197	128	386017	46.726	ng	98
32) Benzoic acid	10.416	122	86684	41.165	ng	94
33) 4-Chloroaniline	11.297	127	40117	10.171	ng	93
34) Hexachlorobutadiene	11.473	225	123410	48.190	ng	96
35) Caprolactam	12.067	113	48660m	45.806	ng	
36) 4-Chloro-3-methylphenol	12.390	107	154012	48.373	ng	95
37) 2-Methylnaphthalene	12.783	142	302075	43.305	ng	100
38) 1-Methylnaphthalene	12.995	142	278441	42.985	ng	96
40) 1,2,4,5-Tetrachloroben...	13.142	216	233712	49.982	ng	99
41) Hexachlorocyclopentadiene	13.118	237	296911	111.376	ng	96
43) 2,4,6-Trichlorophenol	13.371	196	158751	51.344	ng	99

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44) 2,4,5-Trichlorophenol	13.441	196	178308	48.751	ng	97	01/12/2023
46) 1,1'-Biphenyl	13.770	154	443869	49.679	ng	99	Supervised By : Jagrut Upadhyay
47) 2-Chloronaphthalene	13.817	162	355055	50.346	ng	97	
48) 2-Nitroaniline	14.011	65	74982	49.902	ng	98	
49) Acenaphthylene	14.658	152	551452	48.051	ng	99	
50) Dimethylphthalate	14.370	163	463497	46.022	ng	97	
51) 2,6-Dinitrotoluene	14.493	165	105468	49.146	ng	96	01/12/2023
52) Acenaphthene	14.993	154	405368	54.309	ng	99	
53) 3-Nitroaniline	14.822	138	33737	16.519	ng	# 94	
54) 2,4-Dinitrophenol	15.028	184	143802	93.800	ng	95	
55) Dibenzofuran	15.327	168	587651	47.576	ng	98	
56) 4-Nitrophenol	15.116	139	160100	101.925	ng	95	
57) 2,4-Dinitrotoluene	15.275	165	146827	48.923	ng	97	
58) Fluorene	15.968	166	470601	47.369	ng	99	
59) 2,3,4,6-Tetrachlorophenol	15.545	232	174963	50.947	ng	97	
60) Diethylphthalate	15.721	149	436029	46.166	ng	98	
61) 4-Chlorophenyl-phenylether	15.950	204	292516	47.181	ng	98	
62) 4-Nitroaniline	15.980	138	96498	46.125	ng	95	
63) Azobenzene	16.244	77	362278	49.706	ng	97	
65) 4,6-Dinitro-2-methylph...	16.038	198	105772	46.398	ng	94	
66) n-Nitrosodiphenylamine	16.162	169	440006	46.870	ng	95	
67) 4-Bromophenyl-phenylether	16.843	248	199909	46.397	ng	95	
68) Hexachlorobenzene	16.973	284	201772	50.194	ng	98	
69) Atrazine	17.096	200	194558	50.006	ng	97	
70) Pentachlorophenol	17.307	266	269424	87.250	ng	97	
71) Phenanthrene	17.707	178	837434	47.758	ng	99	
72) Anthracene	17.801	178	855415	47.291	ng	99	
73) Carbazole	18.060	167	751286	49.412	ng	99	
74) Di-n-butylphthalate	18.594	149	756090	49.885	ng	100	
75) Fluoranthene	19.699	202	1094655	47.859	ng	100	
77) Benzidine	19.863	184	207996	17.677	ng	99	
78) Pyrene	20.057	202	1105457	46.710	ng	99	
80) Butylbenzylphthalate	20.921	149	328755	48.420	ng	97	
81) Benzo(a)anthracene	21.937	228	1118702	47.436	ng	99	
82) 3,3'-Dichlorobenzidine	21.837	252	165038	19.440	ng	96	
83) Chrysene	22.008	228	1038855	47.027	ng	99	
84) Bis(2-ethylhexyl)phtha...	21.808	149	476563	48.719	ng	99	
85) Di-n-octyl phthalate	23.089	149	812108	49.178	ng	100	
87) Indeno(1,2,3-cd)pyrene	29.358	276	1362188	50.126	ng	# 99	
88) Benzo(b)fluoranthene	24.305	252	1211278	51.020	ng	99	
89) Benzo(k)fluoranthene	24.376	252	1144457	48.409	ng	100	
90) Benzo(a)pyrene	25.239	252	1049955	45.343	ng	# 99	
91) Dibenzo(a,h)anthracene	29.429	278	1135880	49.855	ng	99	
92) Benzo(g,h,i)perylene	30.615	276	1111379	50.302	ng	98	

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

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