

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102122\
 Data File : BG055233.D
 Acq On : 21 Oct 2022 13:36
 Operator : CG/JU
 Sample : SP6027
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SP6027

Quant Time: Oct 22 01:28:11 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG101922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 20 09:21:42 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By : Christian Giraldo 10/24/2022
 Supervised By : Jagrut Upadhyay 10/24/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.287	152	34858	20.000	ng	0.00
21) Naphthalene-d8	11.119	136	156768	20.000	ng	0.00
39) Acenaphthene-d10	14.903	164	106127	20.000	ng	0.00
64) Phenanthrene-d10	17.635	188	237349	20.000	ng	0.00
76) Chrysene-d12	21.918	240	206924	20.000	ng	0.00
86) Perylene-d12	25.332	264	217933	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	0.000	112	0	0.000	ng	
7) Phenol-d6	0.000	99	0d	0.000	ng	
23) Nitrobenzene-d5	0.000	82	0d	0.000	ng	
42) 2,4,6-Tribromophenol	0.000	330	0	0.000	ng	
45) 2-Fluorobiphenyl	0.000	172	0d	0.000	ng	
79) Terphenyl-d14	0.000	244	0d	0.000	ng	
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.616	88	43680	46.359	ng	98
3) Pyridine	4.033	79	117559	39.623	ng	99
4) n-Nitrosodimethylamine	3.945	42	60334	43.457	ng	96
6) Aniline	7.600	93	139706	36.501	ng	100
8) 2-Chlorophenol	7.847	128	127254	54.324	ng	97
9) Benzaldehyde	7.412	77	114679	55.880	ng	98
10) Phenol	7.453	94	175387	52.584	ng	99
11) bis(2-Chloroethyl)ether	7.688	93	121872	47.457	ng	97
12) 1,3-Dichlorobenzene	8.176	146	127012	49.984	ng	98
13) 1,4-Dichlorobenzene	8.323	146	129442	49.868	ng	99
14) 1,2-Dichlorobenzene	8.646	146	124864	50.485	ng	97
15) Benzyl Alcohol	8.522	79	122861	48.456	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.816	45	202377	43.727	ng	96
17) 2-Methylphenol	8.722	107	117377	50.840	ng	99
18) Hexachloroethane	9.386	117	47069	50.098	ng	91
19) n-Nitroso-di-n-propyla...	9.092	70	109297	43.164	ng	95
20) 3+4-Methylphenols	9.051	107	162290	50.012	ng	99
22) Acetophenone	9.116	105	190172	44.847	ng	# 98
24) Nitrobenzene	9.504	77	151799	43.718	ng	95
25) Isophorone	10.026	82	291958	43.946	ng	97
26) 2-Nitrophenol	10.220	139	70710	51.768	ng	99
27) 2,4-Dimethylphenol	10.267	122	135238	62.174	ng	97
28) bis(2-Chloroethoxy)met...	10.502	93	171362	53.677	ng	99
29) 2,4-Dichlorophenol	10.761	162	124897	53.780	ng	98
30) 1,2,4-Trichlorobenzene	10.978	180	114830	48.293	ng	99
31) Naphthalene	11.172	128	387839	46.885	ng	99
32) Benzoic acid	10.408	122	66596	42.773	ng	94
33) 4-Chloroaniline	11.272	127	110491	32.437	ng	97
34) Hexachlorobutadiene	11.448	225	66346	46.208	ng	100
35) Caprolactam	12.036	113	46414	45.478	ng	92
36) 4-Chloro-3-methylphenol	12.371	107	147644	48.893	ng	99
37) 2-Methylnaphthalene	12.753	142	277030	44.028	ng	99
38) 1-Methylnaphthalene	12.970	142	261775	42.143	ng	100
40) 1,2,4,5-Tetrachloroben...	13.111	216	126105	50.552	ng	98
41) Hexachlorocyclopentadiene	13.088	237	143056	125.926	ng	96
43) 2,4,6-Trichlorophenol	13.346	196	96240	50.690	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.417	196	106126	50.950	ng	97
46) 1,1'-Biphenyl	13.740	154	371878	49.486	ng	99
47) 2-Chloronaphthalene	13.787	162	281148	48.596	ng	98
48) 2-Nitroaniline	13.981	65	107808	45.918	ng	98
49) Acenaphthylene	14.627	152	468139	47.300	ng	100
50) Dimethylphthalate	14.345	163	378591	45.352	ng	100
51) 2,6-Dinitrotoluene	14.468	165	82302	49.135	ng	88
52) Acenaphthene	14.968	154	327540m	51.978	ng	
53) 3-Nitroaniline	14.797	138	61070	29.270	ng	96
54) 2,4-Dinitrophenol	15.003	184	92367	109.385	ng	97
55) Dibenzofuran	15.297	168	439691	48.442	ng	98
56) 4-Nitrophenol	15.097	139	148144	100.980	ng	98
57) 2,4-Dinitrotoluene	15.250	165	116684	49.820	ng	95
58) Fluorene	15.943	166	355806	46.547	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.514	232	88244	48.649	ng	97
60) Diethylphthalate	15.690	149	405199	46.398	ng	99
61) 4-Chlorophenyl-phenylether	15.925	204	176702	49.805	ng	96
62) 4-Nitroaniline	15.955	138	97819	46.218	ng	99
63) Azobenzene	16.213	77	405130	45.900	ng	98
65) 4,6-Dinitro-2-methylph...	16.008	198	63141	49.802	ng	99
66) n-Nitrosodiphenylamine	16.137	169	321560	48.543	ng	99
67) 4-Bromophenyl-phenylether	16.813	248	112712	48.721	ng	98
68) Hexachlorobenzene	16.942	284	123705	48.858	ng	99
69) Atrazine	17.071	200	122170	55.957	ng	99
70) Pentachlorophenol	17.283	266	132531	96.990	ng	96
71) Phenanthrene	17.676	178	582650	46.004	ng	100
72) Anthracene	17.770	178	591058	47.756	ng	100
73) Carbazole	18.029	167	565967	49.658	ng	99
74) Di-n-butylphthalate	18.564	149	715958	48.161	ng	100
75) Fluoranthene	19.668	202	653001	46.545	ng	98
77) Benzidine	19.833	184	409002	77.728	ng	100
78) Pyrene	20.026	202	662098	42.464	ng	100
80) Butylbenzylphthalate	20.884	149	304582	41.793	ng	100
81) Benzo(a)anthracene	21.901	228	608039	45.302	ng	100
82) 3,3'-Dichlorobenzidine	21.801	252	111119	24.111	ng	97
83) Chrysene	21.965	228	571487	45.721	ng	99
84) Bis(2-ethylhexyl)phtha...	21.766	149	441071	46.255	ng	99
85) Di-n-octyl phthalate	23.041	149	733281	51.579	ng	97
87) Indeno(1,2,3-cd)pyrene	29.269	276	729417	48.122	ng	# 88
88) Benzo(b)fluoranthene	24.245	252	637377	49.250	ng	# 97
89) Benzo(k)fluoranthene	24.316	252	606736	46.292	ng	# 97
90) Benzo(a)pyrene	25.173	252	555173	49.980	ng	# 96
91) Dibenzo(a,h)anthracene	29.339	278	614290	49.468	ng	# 94
92) Benzo(g,h,i)perylene	30.508	276	596934	49.622	ng	# 90

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

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