

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG021523\
 Data File : BG056638.D
 Acq On : 15 Feb 2023 17:49
 Operator : CG/JU
 Sample : SSTDICC050
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC050

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 02/16/2023
 Supervised By : Jagrut Upadhyay 02/16/2023

Quant Time: Feb 16 01:27:22 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG021523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 16 01:15:40 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.441	152	17671	20.000	ng	0.00
21) Naphthalene-d8	11.279	136	71684	20.000	ng	0.00
39) Acenaphthene-d10	15.045	164	52732	20.000	ng	0.00
64) Phenanthrene-d10	17.777	188	133381	20.000	ng	# 0.00
76) Chrysene-d12	22.102	240	125254	20.000	ng	0.00
86) Perylene-d12	25.645	264	141789	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.944	112	110460	101.707	ng	0.00
7) Phenol-d6	7.560	99	163544	101.942	ng	0.00
23) Nitrobenzene-d5	9.616	82	146572	104.691	ng	0.00
42) 2,4,6-Tribromophenol	16.520	330	74168	102.731	ng	0.00
45) 2-Fluorobiphenyl	13.676	172	399147	101.358	ng	0.00
79) Terphenyl-d14	20.351	244	769103	100.463	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.759	88	24094	50.602	ng	97
3) Pyridine	4.182	79	76378	50.713	ng	96
4) n-Nitrosodimethylamine	4.088	42	35387	50.407	ng	93
6) Aniline	7.754	93	110623	49.616	ng	97
8) 2-Chlorophenol	7.995	128	54597	50.808	ng	92
9) Benzaldehyde	7.566	77	38986	45.952	ng	95
10) Phenol	7.589	94	83983	51.176	ng	99
11) bis(2-Chloroethyl)ether	7.842	93	60935	49.488	ng	96
12) 1,3-Dichlorobenzene	8.330	146	62059	48.473	ng	94
13) 1,4-Dichlorobenzene	8.477	146	63894	49.555	ng	94
14) 1,2-Dichlorobenzene	8.806	146	61501	49.624	ng	97
15) Benzyl Alcohol	8.670	79	74122	51.600	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.964	45	81071m	49.023	ng	
17) 2-Methylphenol	8.864	107	56680	49.148	ng	98
18) Hexachloroethane	9.546	117	21577	52.829	ng	98
19) n-Nitroso-di-n-propyla...	9.246	70	60927	50.729	ng	97
20) 3+4-Methylphenols	9.193	107	79802	50.373	ng	93
22) Acetophenone	9.270	105	107075	48.886	ng	98
24) Nitrobenzene	9.658	77	79583	48.757	ng	95
25) Isophorone	10.180	82	167841	49.259	ng	97
26) 2-Nitrophenol	10.374	139	29286	50.802	ng	97
27) 2,4-Dimethylphenol	10.415	122	63407	48.133	ng	98
28) bis(2-Chloroethoxy)met...	10.656	93	85123	49.021	ng	98
29) 2,4-Dichlorophenol	10.909	162	66501	50.934	ng	97
30) 1,2,4-Trichlorobenzene	11.138	180	76500	49.187	ng	96
31) Naphthalene	11.332	128	195164	49.124	ng	99
32) Benzoic acid	10.545	122	44067m	51.802	ng	
33) 4-Chloroaniline	11.426	127	89360	49.231	ng	97
34) Hexachlorobutadiene	11.608	225	54677	49.194	ng	97
35) Caprolactam	12.190	113	20864	49.684	ng	# 81
36) 4-Chloro-3-methylphenol	12.501	107	72723	50.509	ng	95
37) 2-Methylnaphthalene	12.901	142	150052	47.835	ng	99
38) 1-Methylnaphthalene	13.118	142	141253	48.257	ng	100
40) 1,2,4,5-Tetrachloroben...	13.259	216	100174	51.102	ng	98
41) Hexachlorocyclopentadiene	13.242	237	57615	54.189	ng	98
43) 2,4,6-Trichlorophenol	13.482	196	64114	53.906	ng	95

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.553	196	75780	53.162	ng	98
46) 1,1'-Biphenyl	13.888	154	200721	50.249	ng	99
47) 2-Chloronaphthalene	13.935	162	158826	50.245	ng	99
48) 2-Nitroaniline	14.123	65	43748	47.563	ng	92
49) Acenaphthylene	14.775	152	258652	50.359	ng	98
50) Dimethylphthalate	14.487	163	221635	50.843	ng	98
51) 2,6-Dinitrotoluene	14.605	165	41866	49.251	ng	91
52) Acenaphthene	15.110	154	158194	49.932	ng	97
53) 3-Nitroaniline	14.934	138	45107	53.307	ng	# 94
54) 2,4-Dinitrophenol	15.128	184	25701	51.034	ng	# 59
55) Dibenzofuran	15.439	168	266605	49.810	ng	99
56) 4-Nitrophenol	15.210	139	34353	52.988	ng	96
57) 2,4-Dinitrotoluene	15.380	165	55922	52.601	ng	92
58) Fluorene	16.085	166	208947	49.373	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.651	232	65946	52.717	ng	# 92
60) Diethylphthalate	15.833	149	222886	50.688	ng	99
61) 4-Chlorophenyl-phenyle...	16.068	204	130292	50.395	ng	99
62) 4-Nitroaniline	16.091	138	46331	52.537	ng	91
63) Azobenzene	16.361	77	217452	49.412	ng	99
65) 4,6-Dinitro-2-methylph...	16.144	198	35900	52.121	ng	94
66) n-Nitrosodiphenylamine	16.279	169	193062	49.755	ng	99
67) 4-Bromophenyl-phenylether	16.961	248	83087	48.518	ng	96
68) Hexachlorobenzene	17.084	284	86321	49.476	ng	97
69) Atrazine	17.208	200	75676	50.790	ng	97
70) Pentachlorophenol	17.419	266	61875	53.551	ng	95
71) Phenanthrene	17.824	178	361058	49.756	ng	99
72) Anthracene	17.913	178	367436	49.485	ng	98
73) Carbazole	18.171	167	314322	49.106	ng	98
74) Di-n-butylphthalate	18.706	149	373746	52.177	ng	100
75) Fluoranthene	19.804	202	473219	49.423	ng	100
77) Benzidine	19.969	184	156229	52.150	ng	97
78) Pyrene	20.169	202	470284	49.716	ng	99
80) Butylbenzylphthalate	21.038	149	155477	54.734	ng	98
81) Benzo(a)anthracene	22.078	228	463911	49.804	ng	98
82) 3,3'-Dichlorobenzidine	21.973	252	154652	51.173	ng	99
83) Chrysene	22.155	228	435381	49.998	ng	98
84) Bis(2-ethylhexyl)phtha...	21.955	149	229937	54.019	ng	100
85) Di-n-octyl phthalate	23.283	149	381514	53.926	ng	99
87) Indeno(1,2,3-cd)pyrene	29.752	276	561234	50.541	ng	# 98
88) Benzo(b)fluoranthene	24.511	252	468655	50.593	ng	99
89) Benzo(k)fluoranthene	24.587	252	453639	49.800	ng	97
90) Benzo(a)pyrene	25.480	252	447718	50.283	ng	98
91) Dibenzo(a,h)anthracene	29.828	278	462611	50.300	ng	97
92) Benzo(g,h,i)perylene	31.032	276	455659	50.704	ng	97

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

Manual Integrations

Reviewed By :Christian Giraldo 02/16/2023
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Abundance

TIC: BG056638.D\data.ms

Time-->