

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052322\
 Data File : BG053660.D
 Acq On : 23 May 2022 12:01
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Reviewed By :Christian
 Giraldo

Quant Time: May 24 05:08:48 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG050522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 24 05:05:32 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.058	152	32181	20.000	ng	0.00
21) Naphthalene-d8	10.872	136	123885	20.000	ng	0.00
39) Acenaphthene-d10	14.690	164	93024	20.000	ng	0.00
64) Phenanthrene-d10	17.439	188	216891	20.000	ng	0.00
76) Chrysene-d12	21.716	240	187606	20.000	ng	0.00
86) Perylene-d12	24.982	264	187219	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.626	112	149070	78.615	ng	0.00
7) Phenol-d6	7.229	99	227060	78.901	ng	0.00
23) Nitrobenzene-d5	9.227	82	242636	83.373	ng	0.00
42) 2,4,6-Tribromophenol	16.182	330	111145	81.010	ng	0.00
45) 2-Fluorobiphenyl	13.315	172	533005	83.728	ng	0.00
79) Terphenyl-d14	20.041	244	868925	86.263	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.493	88	33546	33.718	ng	95
3) Pyridine	3.899	79	91546	37.678	ng	93
4) n-Nitrosodimethylamine	3.810	42	44616	37.696	ng	# 98
6) Aniline	7.382	93	126467	39.678	ng	99
8) 2-Chlorophenol	7.629	128	78751	39.727	ng	96
9) Benzaldehyde	7.194	77	62783m	34.341	ng	
10) Phenol	7.253	94	111880	39.751	ng	95
11) bis(2-Chloroethyl)ether	7.470	93	82976	39.529	ng	98
12) 1,3-Dichlorobenzene	7.952	146	92109	39.388	ng	95
13) 1,4-Dichlorobenzene	8.093	146	91636	38.832	ng	96
14) 1,2-Dichlorobenzene	8.416	146	90109	40.663	ng	96
15) Benzyl Alcohol	8.299	79	93121	38.705	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.586	45	133314	39.022	ng	99
17) 2-Methylphenol	8.510	107	75902	39.900	ng	96
18) Hexachloroethane	9.144	117	34404	38.391	ng	95
19) n-Nitroso-di-n-propyla...	8.862	70	80473	38.974	ng	98
20) 3+4-Methylphenols	8.839	107	103961	38.695	ng	94
22) Acetophenone	8.880	105	140844	41.540	ng	# 99
24) Nitrobenzene	9.268	77	114835	40.768	ng	98
25) Isophorone	9.791	82	202579	41.361	ng	100
26) 2-Nitrophenol	9.984	139	48531	42.540	ng	95
27) 2,4-Dimethylphenol	10.043	122	69612	40.545	ng	94
28) bis(2-Chloroethoxy)met...	10.266	93	112620	39.980	ng	98
29) 2,4-Dichlorophenol	10.531	162	87634	42.292	ng	91
30) 1,2,4-Trichlorobenzene	10.736	180	100198	42.559	ng	98
31) Naphthalene	10.924	128	263435	41.804	ng	100
32) Benzoic acid	10.225	122	34834m	35.205	ng	
33) 4-Chloroaniline	11.030	127	110409	41.836	ng	98
34) Hexachlorobutadiene	11.206	225	76972	44.090	ng	95
35) Caprolactam	11.806	113	29298	38.908	ng	92
36) 4-Chloro-3-methylphenol	12.158	107	98493	39.737	ng	98
37) 2-Methylnaphthalene	12.522	142	190617	40.532	ng	99
38) 1-Methylnaphthalene	12.740	142	185583	40.446	ng	98
40) 1,2,4,5-Tetrachloroben...	12.892	216	125000	42.630	ng	100
41) Hexachlorocyclopentadiene	12.869	237	30525	31.571	ng	98
43) 2,4,6-Trichlorophenol	13.133	196	81491	42.576	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
44) 2,4,5-Trichlorophenol	13.215	196	86259	42.344	ng	98	05/24/2022
46) 1,1'-Biphenyl	13.521	154	264942	41.783	ng	98	Supervised By :Jagrut Upadhyay
47) 2-Chloronaphthalene	13.574	162	205630	41.392	ng	99	
48) 2-Nitroaniline	13.773	65	77990	42.131	ng	97	
49) Acenaphthylene	14.414	152	321769	40.589	ng	99	
50) Dimethylphthalate	14.144	163	276096	39.058	ng	99	
51) 2,6-Dinitrotoluene	14.261	165	57663	40.039	ng	95	05/24/2022
52) Acenaphthene	14.755	154	190825	40.392	ng	99	
53) 3-Nitroaniline	14.596	138	62457	40.137	ng	96	
54) 2,4-Dinitrophenol	14.819	184	30865	37.372	ng	92	
55) Dibenzofuran	15.089	168	338096	39.842	ng	98	
56) 4-Nitrophenol	14.925	139	41164	37.037	ng	90	
57) 2,4-Dinitrotoluene	15.054	165	84353	40.408	ng	92	
58) Fluorene	15.736	166	269526	39.579	ng	97	
59) 2,3,4,6-Tetrachlorophenol	15.324	232	79514	38.471	ng	98	
60) Diethylphthalate	15.495	149	285050	39.037	ng	99	
61) 4-Chlorophenyl-phenyle...	15.724	204	150536	39.615	ng	98	
62) 4-Nitroaniline	15.759	138	64249	39.808	ng	99	
63) Azobenzene	16.017	77	295684	39.544	ng	98	
65) 4,6-Dinitro-2-methylph...	15.830	198	57447	44.450	ng	94	
66) n-Nitrosodiphenylamine	15.941	169	239118	41.991	ng	99	
67) 4-Bromophenyl-phenylether	16.617	248	107297	42.359	ng	98	
68) Hexachlorobenzene	16.752	284	113976	41.085	ng	97	
69) Atrazine	16.881	200	92439	38.560	ng	97	
70) Pentachlorophenol	17.098	266	50493	36.484	ng	96	
71) Phenanthrene	17.480	178	445936	40.435	ng	97	
72) Anthracene	17.568	178	441062	40.688	ng	99	
73) Carbazole	17.839	167	425822	40.082	ng	100	
74) Di-n-butylphthalate	18.385	149	479044	40.093	ng	99	
75) Fluoranthene	19.489	202	540541	37.853	ng	98	
77) Benzidine	19.660	184	202009	49.958	ng	98	
78) Pyrene	19.853	202	541799	44.612	ng	99	
80) Butylbenzylphthalate	20.723	149	193613	43.475	ng	96	
81) Benzo(a)anthracene	21.698	228	484044	40.452	ng	99	
82) 3,3'-Dichlorobenzidine	21.604	252	131205	43.239	ng	96	
83) Chrysene	21.763	228	462899	41.420	ng	98	
84) Bis(2-ethylhexyl)phtha...	21.581	149	258748	42.095	ng	98	
85) Di-n-octyl phthalate	22.802	149	430803	41.447	ng	99	
87) Indeno(1,2,3-cd)pyrene	28.730	276	539848	41.401	ng	# 95	
88) Benzo(b)fluoranthene	23.942	252	457116	41.249	ng	99	
89) Benzo(k)fluoranthene	24.013	252	457822	42.042	ng	96	
90) Benzo(a)pyrene	24.829	252	389430	41.287	ng	99	
91) Dibenzo(a,h)anthracene	28.783	278	446821	41.586	ng	98	
92) Benzo(g,h,i)perylene	29.905	276	439220	41.240	ng	95	

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

Manual Integrations

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