

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG051222\
 Data File : BG053531.D
 Acq On : 12 May 2022 15:28
 Operator : CG/JU
 Sample : PB144706BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB144706BS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 05/13/2022
 Supervised By : Jagrut Upadhyay 05/13/2022

Quant Time: May 13 06:36:17 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG050522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 11 06:00:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.088	152	23246	20.000	ng	0.00
21) Naphthalene-d8	10.896	136	102412	20.000	ng	0.00
39) Acenaphthene-d10	14.714	164	85549	20.000	ng	0.00
64) Phenanthrene-d10	17.458	188	211918	20.000	ng	0.00
76) Chrysene-d12	21.740	240	205976	20.000	ng	0.00
86) Perylene-d12	25.018	264	215351	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.650	112	189755	138.536	ng	0.00
7) Phenol-d6	7.248	99	280068	134.728	ng	0.00
23) Nitrobenzene-d5	9.251	82	194267	80.749	ng	0.00
42) 2,4,6-Tribromophenol	16.201	330	157567	124.880	ng	0.00
45) 2-Fluorobiphenyl	13.334	172	456616	77.996	ng	0.00
79) Terphenyl-d14	20.060	244	945253	85.471	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.524	88	25718	35.785	ng	93
3) Pyridine	3.929	79	53718	30.607	ng	98
4) n-Nitrosodimethylamine	3.841	42	34643	40.520	ng	# 99
6) Aniline	7.407	93	100934	43.839	ng	95
8) 2-Chlorophenol	7.653	128	68722	47.993	ng	95
9) Benzaldehyde	7.219	77	46967	35.565	ng	88
10) Phenol	7.277	94	97873	48.140	ng	97
11) bis(2-Chloroethyl)ether	7.495	93	66777	44.040	ng	97
12) 1,3-Dichlorobenzene	7.976	146	68834	40.748	ng	93
13) 1,4-Dichlorobenzene	8.123	146	70070	41.106	ng	100
14) 1,2-Dichlorobenzene	8.441	146	68160	42.581	ng	99
15) Benzyl Alcohol	8.323	79	80612m	46.384	ng	
16) 2,2'-oxybis(1-Chloropr...	8.611	45	105145	42.606	ng	97
17) 2-Methylphenol	8.529	107	67566	49.169	ng	95
18) Hexachloroethane	9.175	117	26374	40.743	ng	97
19) n-Nitroso-di-n-propyla...	8.887	70	65999	44.250	ng	98
20) 3+4-Methylphenols	8.858	107	91010	46.895	ng	97
22) Acetophenone	8.905	105	111390	39.741	ng	# 98
24) Nitrobenzene	9.292	77	98631	42.358	ng	98
25) Isophorone	9.809	82	184073	45.463	ng	98
26) 2-Nitrophenol	10.009	139	39662	42.055	ng	94
27) 2,4-Dimethylphenol	10.062	122	72415	51.022	ng	95
28) bis(2-Chloroethoxy)met...	10.291	93	96926	41.623	ng	98
29) 2,4-Dichlorophenol	10.549	162	78868	46.042	ng	96
30) 1,2,4-Trichlorobenzene	10.761	180	79618	40.909	ng	98
31) Naphthalene	10.949	128	215781	41.421	ng	98
32) Benzoic acid	10.238	122	31312m	37.540	ng	
33) 4-Chloroaniline	11.055	127	70444	32.289	ng	97
34) Hexachlorobutadiene	11.231	225	57543	39.872	ng	97
35) Caprolactam	11.818	113	28542m	45.851	ng	
36) 4-Chloro-3-methylphenol	12.177	107	92997	45.386	ng	98
37) 2-Methylnaphthalene	12.547	142	161524	41.548	ng	97
38) 1-Methylnaphthalene	12.764	142	156436	41.243	ng	99
40) 1,2,4,5-Tetrachloroben...	12.917	216	106927	39.652	ng	99
41) Hexachlorocyclopentadiene	12.893	237	100938	90.582	ng	97
43) 2,4,6-Trichlorophenol	13.152	196	72208	41.022	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.234	196	80193	42.806	ng	96
46) 1,1'-Biphenyl	13.545	154	228343	39.158	ng	99
47) 2-Chloronaphthalene	13.592	162	178631	39.099	ng	99
48) 2-Nitroaniline	13.792	65	75302	44.233	ng	98
49) Acenaphthylene	14.438	152	312678	42.889	ng	99
50) Dimethylphthalate	14.162	163	267448	41.140	ng	99
51) 2,6-Dinitrotoluene	14.280	165	57102	43.114	ng	97
52) Acenaphthene	14.773	154	175004	40.279	ng	99
53) 3-Nitroaniline	14.615	138	48975	34.223	ng	99
54) 2,4-Dinitrophenol	14.826	184	64666	77.158	ng	97
55) Dibenzofuran	15.108	168	307215	39.366	ng	98
56) 4-Nitrophenol	14.932	139	88278	86.367	ng	87
57) 2,4-Dinitrotoluene	15.073	165	83723	43.610	ng	# 85
58) Fluorene	15.760	166	257656	41.142	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.343	232	76487	40.240	ng	97
60) Diethylphthalate	15.519	149	282312	42.040	ng	98
61) 4-Chlorophenyl-phenyle...	15.742	204	144071	41.227	ng	99
62) 4-Nitroaniline	15.778	138	61029	41.117	ng	97
63) Azobenzene	16.036	77	279152	40.595	ng	98
65) 4,6-Dinitro-2-methylph...	15.842	198	50904	40.312	ng	96
66) n-Nitrosodiphenylamine	15.960	169	232212	41.735	ng	99
67) 4-Bromophenyl-phenylether	16.641	248	103715	41.906	ng	98
68) Hexachlorobenzene	16.770	284	113294	41.798	ng	98
69) Atrazine	16.906	200	105414	45.004	ng	95
70) Pentachlorophenol	17.117	266	110519	75.962	ng	93
71) Phenanthrene	17.499	178	451071	41.860	ng	98
72) Anthracene	17.593	178	452331	42.707	ng	99
73) Carbazole	17.857	167	412917	39.779	ng	99
74) Di-n-butylphthalate	18.403	149	484084	41.465	ng	100
75) Fluoranthene	19.508	202	577150	41.366	ng	99
77) Benzidine	19.678	184	283279	63.808	ng	99
78) Pyrene	19.872	202	567023	42.525	ng	99
80) Butylbenzylphthalate	20.741	149	208667	42.677	ng	98
81) Benzo(a)anthracene	21.717	228	552313	42.041	ng	99
82) 3,3'-Dichlorobenzidine	21.623	252	152765	45.854	ng	98
83) Chrysene	21.787	228	505170	41.171	ng	97
84) Bis(2-ethylhexyl)phtha...	21.605	149	291354	43.173	ng	99
85) Di-n-octyl phthalate	22.833	149	499669	43.785	ng	99
87) Indeno(1,2,3-cd)pyrene	28.778	276	621704	41.450	ng	# 97
88) Benzo(b)fluoranthene	23.978	252	577264	45.286	ng	99
89) Benzo(k)fluoranthene	24.043	252	547876	43.739	ng	100
90) Benzo(a)pyrene	24.865	252	557470	51.382	ng	98
91) Dibenzo(a,h)anthracene	28.831	278	537865	43.521	ng	99
92) Benzo(g,h,i)perylene	29.958	276	538659	43.970	ng	97

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

