

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG051322\
 Data File : BG053551.D
 Acq On : 13 May 2022 20:24
 Operator : CG/JU
 Sample : N2762-06MSD
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB16MSD

Quant Time: May 14 03:14:44 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

Manual Integrations APPROVED

Reviewed By : Christian Giraldo 05/16/2022
 Supervised By : Jagrut Upadhyay 05/16/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.082	152	24117	20.000	ng	0.00
21) Naphthalene-d8	10.896	136	99032	20.000	ng	0.00
39) Acenaphthene-d10	14.708	164	79215	20.000	ng	0.00
64) Phenanthrene-d10	17.457	188	193852	20.000	ng	0.00
76) Chrysene-d12	21.740	240	185198	20.000	ng	0.00
86) Perylene-d12	25.012	264	195528	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.656	112	165862	116.719	ng	0.00
7) Phenol-d6	7.248	99	253071	117.344	ng	0.00
23) Nitrobenzene-d5	9.245	82	174093	74.833	ng	0.00
42) 2,4,6-Tribromophenol	16.200	330	131144	112.249	ng	0.00
45) 2-Fluorobiphenyl	13.334	172	375738	69.312	ng	0.00
79) Terphenyl-d14	20.060	244	683685	68.756	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.529	88	29180	39.136	ng	# 90
3) Pyridine	3.929	79	72294	39.703	ng	98
4) n-Nitrosodimethylamine	3.841	42	44567	50.245	ng	96
6) Aniline	7.406	93	75511	31.612	ng	96
8) 2-Chlorophenol	7.653	128	81254	54.696	ng	95
9) Benzaldehyde	7.218	77	54485m	39.767	ng	
10) Phenol	7.277	94	115077	54.558	ng	94
11) bis(2-Chloroethyl)ether	7.494	93	74315	47.241	ng	96
12) 1,3-Dichlorobenzene	7.976	146	85600	48.843	ng	93
13) 1,4-Dichlorobenzene	8.117	146	86906	49.141	ng	97
14) 1,2-Dichlorobenzene	8.440	146	83979	50.569	ng	96
15) Benzyl Alcohol	8.323	79	96044m	53.268	ng	
16) 2,2'-oxybis(1-Chloropr...	8.611	45	123981	48.424	ng	98
17) 2-Methylphenol	8.528	107	77508	54.367	ng	97
18) Hexachloroethane	9.174	117	32804	48.846	ng	96
19) n-Nitroso-di-n-propyla...	8.887	70	75889	49.044	ng	99
20) 3+4-Methylphenols	8.857	107	108674	53.974	ng	98
22) Acetophenone	8.904	105	131234	48.419	ng	# 98
24) Nitrobenzene	9.292	77	115825	51.439	ng	97
25) Isophorone	9.809	82	211701	54.071	ng	98
26) 2-Nitrophenol	10.003	139	48473	53.152	ng	96
27) 2,4-Dimethylphenol	10.062	122	82371	60.017	ng	98
28) bis(2-Chloroethoxy)met...	10.291	93	111557	49.541	ng	97
29) 2,4-Dichlorophenol	10.549	162	95203	57.475	ng	90
30) 1,2,4-Trichlorobenzene	10.761	180	94490	50.207	ng	99
31) Naphthalene	10.943	128	252380	50.100	ng	98
32) Benzoic acid	10.244	122	48123m	54.663	ng	
33) 4-Chloroaniline	11.054	127	32925	15.607	ng	95
34) Hexachlorobutadiene	11.230	225	69064	49.489	ng	97
35) Caprolactam	11.824	113	32229m	53.542	ng	
36) 4-Chloro-3-methylphenol	12.176	107	108086	54.551	ng	95
37) 2-Methylnaphthalene	12.546	142	189108	50.303	ng	96
38) 1-Methylnaphthalene	12.764	142	181165	49.392	ng	97
40) 1,2,4,5-Tetrachloroben...	12.916	216	126438	50.637	ng	99
41) Hexachlorocyclopentadiene	12.893	237	105016	100.685	ng	97
43) 2,4,6-Trichlorophenol	13.151	196	88881	54.532	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.234	196	97326	56.106	ng	98
46) 1,1'-Biphenyl	13.545	154	269037	49.825	ng	99
47) 2-Chloronaphthalene	13.592	162	209067	49.420	ng	97
48) 2-Nitroaniline	13.792	65	85205	54.052	ng	99
49) Acenaphthylene	14.432	152	362305	53.669	ng	99
50) Dimethylphthalate	14.162	163	294399	48.907	ng	99
51) 2,6-Dinitrotoluene	14.279	165	65064	53.054	ng	97
52) Acenaphthene	14.773	154	206650	51.366	ng	99
53) 3-Nitroaniline	14.614	138	32138	24.253	ng	99
54) 2,4-Dinitrophenol	14.832	184	78702	99.451	ng	92
55) Dibenzofuran	15.108	168	353299	48.891	ng	98
56) 4-Nitrophenol	14.937	139	101490	107.232	ng	86
57) 2,4-Dinitrotoluene	15.072	165	90715	51.031	ng	# 84
58) Fluorene	15.760	166	291639	50.292	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.343	232	90389	51.356	ng	99
60) Diethylphthalate	15.519	149	303966	48.884	ng	98
61) 4-Chlorophenyl-phenyle...	15.742	204	162681	50.274	ng	98
62) 4-Nitroaniline	15.783	138	62575	45.529	ng	96
63) Azobenzene	16.036	77	312461	49.073	ng	99
65) 4,6-Dinitro-2-methylph...	15.842	198	58694	50.812	ng	98
66) n-Nitrosodiphenylamine	15.959	169	261866	51.451	ng	99
67) 4-Bromophenyl-phenylether	16.641	248	116021	51.247	ng	98
68) Hexachlorobenzene	16.770	284	127331	51.354	ng	97
69) Atrazine	16.905	200	110990	51.801	ng	96
70) Pentachlorophenol	17.117	266	134512	99.532	ng	93
71) Phenanthrene	17.498	178	523652	53.125	ng	97
72) Anthracene	17.592	178	504073	52.027	ng	99
73) Carbazole	17.857	167	453528	47.763	ng	99
74) Di-n-butylphthalate	18.403	149	517838	48.491	ng	99
75) Fluoranthene	19.508	202	659540	51.676	ng	99
77) Benzidine	19.678	184	216733	54.296	ng	98
78) Pyrene	19.866	202	651870	54.373	ng	99
80) Butylbenzylphthalate	20.741	149	228547	51.987	ng	97
81) Benzo(a)anthracene	21.716	228	635720	53.819	ng	99
82) 3,3'-Dichlorobenzidine	21.622	252	119034	39.738	ng	95
83) Chrysene	21.787	228	581988	52.753	ng	97
84) Bis(2-ethylhexyl)phtha...	21.605	149	322737	53.188	ng	98
85) Di-n-octyl phthalate	22.832	149	546713	53.283	ng	99
87) Indeno(1,2,3-cd)pyrene	28.771	276	717041	52.654	ng	# 98
88) Benzo(b)fluoranthene	23.978	252	660843	57.099	ng	99
89) Benzo(k)fluoranthene	24.043	252	626016	55.044	ng	99
90) Benzo(a)pyrene	24.865	252	636168	64.580	ng	97
91) Dibenzo(a,h)anthracene	28.830	278	600932	53.553	ng	99
92) Benzo(g,h,i)perylene	29.952	276	615194	55.308	ng	96

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

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