

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG051322\
 Data File : BG053538.D
 Acq On : 13 May 2022 11:44
 Operator : CG/JU
 Sample : N2749-19MS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-7MS

Quant Time: May 14 03:12:47 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

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 Upadhyay
 05/16/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.084	152	27209	20.000	ng	0.00
21) Naphthalene-d8	10.898	136	109826	20.000	ng	0.00
39) Acenaphthene-d10	14.710	164	85125	20.000	ng	0.00
64) Phenanthrene-d10	17.454	188	201832	20.000	ng	0.00
76) Chrysene-d12	21.736	240	202390	20.000	ng	0.00
86) Perylene-d12	25.008	264	214762	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.652	112	179372	111.882	ng	0.00
7) Phenol-d6	7.250	99	269344	110.698	ng	0.00
23) Nitrobenzene-d5	9.247	82	184803	71.629	ng	0.00
42) 2,4,6-Tribromophenol	16.203	330	135677	108.067	ng	0.00
45) 2-Fluorobiphenyl	13.336	172	412829	70.867	ng	0.00
79) Terphenyl-d14	20.056	244	758041	69.758	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.526	88	32034	38.082	ng	# 85
3) Pyridine	3.925	79	80695m	39.281	ng	
4) n-Nitrosodimethylamine	3.843	42	49969	49.933	ng	98
6) Aniline	7.403	93	71907	26.683	ng	94
8) 2-Chlorophenol	7.649	128	88351	52.715	ng	99
9) Benzaldehyde	7.215	77	60179m	38.932	ng	
10) Phenol	7.279	94	126149	53.011	ng	95
11) bis(2-Chloroethyl)ether	7.497	93	84102	47.387	ng	95
12) 1,3-Dichlorobenzene	7.973	146	91417	46.235	ng	93
13) 1,4-Dichlorobenzene	8.119	146	94482	47.354	ng	98
14) 1,2-Dichlorobenzene	8.437	146	91344	48.753	ng	97
15) Benzyl Alcohol	8.319	79	101638m	49.964	ng	
16) 2,2'-oxybis(1-Chloropr...	8.607	45	133874	46.346	ng	97
17) 2-Methylphenol	8.531	107	83940	52.188	ng	98
18) Hexachloroethane	9.171	117	34856	46.003	ng	88
19) n-Nitroso-di-n-propyla...	8.883	70	81345	46.596	ng	99
20) 3+4-Methylphenols	8.860	107	115648	50.911	ng	96
22) Acetophenone	8.907	105	140489	46.739	ng	# 98
24) Nitrobenzene	9.288	77	125991	50.455	ng	100
25) Isophorone	9.811	82	223267	51.420	ng	98
26) 2-Nitrophenol	10.005	139	51035	50.461	ng	98
27) 2,4-Dimethylphenol	10.064	122	90292	59.323	ng	93
28) bis(2-Chloroethoxy)met...	10.287	93	119842	47.990	ng	98
29) 2,4-Dichlorophenol	10.546	162	99250	54.029	ng	92
30) 1,2,4-Trichlorobenzene	10.757	180	102151	48.943	ng	96
31) Naphthalene	10.945	128	275418	49.300	ng	100
32) Benzoic acid	10.234	122	55400m	56.421	ng	
33) 4-Chloroaniline	11.051	127	32736	13.992	ng	93
34) Hexachlorobutadiene	11.233	225	75006	48.464	ng	98
35) Caprolactam	11.820	113	33271m	49.840	ng	
36) 4-Chloro-3-methylphenol	12.179	107	112589	51.239	ng	97
37) 2-Methylnaphthalene	12.543	142	196572	47.149	ng	98
38) 1-Methylnaphthalene	12.760	142	189293m	46.536	ng	
40) 1,2,4,5-Tetrachloroben...	12.913	216	131367	48.958	ng	99
41) Hexachlorocyclopentadiene	12.895	237	117560	104.518	ng	98
43) 2,4,6-Trichlorophenol	13.154	196	92173	52.625	ng	95

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.230	196	97266	52.178	ng	99
46) 1,1'-Biphenyl	13.547	154	278713	48.034	ng	100
47) 2-Chloronaphthalene	13.594	162	218221	48.002	ng	97
48) 2-Nitroaniline	13.788	65	85595	50.530	ng	97
49) Acenaphthylene	14.434	152	374299	51.597	ng	99
50) Dimethylphthalate	14.158	163	304404	47.058	ng	99
51) 2,6-Dinitrotoluene	14.282	165	66237	50.260	ng	96
52) Acenaphthene	14.775	154	210939	48.792	ng	98
53) 3-Nitroaniline	14.616	138	28712	20.164	ng	99
54) 2,4-Dinitrophenol	14.828	184	83459	98.223	ng	98
55) Dibenzofuran	15.110	168	361824	46.595	ng	98
56) 4-Nitrophenol	14.934	139	104500	102.747	ng	90
57) 2,4-Dinitrotoluene	15.075	165	94785	49.618	ng	# 94
58) Fluorene	15.756	166	298929	47.970	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.339	232	92143	48.718	ng	99
60) Diethylphthalate	15.515	149	311782	46.660	ng	98
61) 4-Chlorophenyl-phenyle...	15.744	204	164781	47.388	ng	97
62) 4-Nitroaniline	15.780	138	65924	44.636	ng	97
63) Azobenzene	16.038	77	316634	46.275	ng	98
65) 4,6-Dinitro-2-methylph...	15.838	198	59285	49.295	ng	96
66) n-Nitrosodiphenylamine	15.956	169	263405	49.707	ng	99
67) 4-Bromophenyl-phenylether	16.637	248	116826	49.562	ng	96
68) Hexachlorobenzene	16.767	284	129347	50.105	ng	99
69) Atrazine	16.902	200	113409	50.837	ng	99
70) Pentachlorophenol	17.113	266	137439	97.764	ng	97
71) Phenanthrene	17.501	178	546685	53.269	ng	97
72) Anthracene	17.589	178	515112	51.065	ng	99
73) Carbazole	17.859	167	463903	46.924	ng	100
74) Di-n-butylphthalate	18.405	149	524572	47.179	ng	99
75) Fluoranthene	19.504	202	701127	52.762	ng	98
77) Benzidine	19.680	184	248475	56.960	ng	98
78) Pyrene	19.868	202	692935	52.889	ng	100
80) Butylbenzylphthalate	20.744	149	233920	48.689	ng	97
81) Benzo(a)anthracene	21.719	228	682457	52.867	ng	98
82) 3,3'-Dichlorobenzidine	21.625	252	88339	26.986	ng	99
83) Chrysene	21.783	228	631094	52.345	ng	97
84) Bis(2-ethylhexyl)phtha...	21.607	149	335628	50.614	ng	100
85) Di-n-octyl phthalate	22.829	149	574088	51.198	ng	100
87) Indeno(1,2,3-cd)pyrene	28.774	276	776679	51.925	ng	# 95
88) Benzo(b)fluoranthene	23.974	252	769149	60.505	ng	99
89) Benzo(k)fluoranthene	24.045	252	669283	53.578	ng	98
90) Benzo(a)pyrene	24.867	252	700470	64.739	ng	98
91) Dibenzo(a,h)anthracene	28.833	278	638460	51.802	ng	98
92) Benzo(g,h,i)perylene	29.955	276	687432	56.268	ng	98

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

