

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100820\
 Data File : BG046817.D
 Acq On : 8 Oct 2020 16:07
 Operator : CG/JU
 Sample : L4293-04MS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 B-6-MW-2MS

Manual Integrations
 APPROVED

mohammad
 10/9/2020 2:46:34 PM

Quant Time: Oct 08 16:46:09 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG092220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 07 11:38:17 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.108	152	63102	20.00	ng	0.00
21) Naphthalene-d8	10.922	136	242952	20.00	ng	0.00
39) Acenaphthene-d10	14.735	164	166471	20.00	ng	0.00
64) Phenanthrene-d10	17.473	188	407026	20.00	ng	# 0.00
76) Chrysene-d12	21.751	240	466323	20.00	ng	# 0.00
86) Perylene-d12	25.023	264	509379	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.670	112	225526	57.26	ng	0.00
7) Phenol-d6	7.256	99	180302	32.45	ng	0.00
23) Nitrobenzene-d5	9.271	82	486079	106.44	ng	0.00
42) 2,4,6-Tribromophenol	16.216	330	341312	155.49	ng	0.00
45) 2-Fluorobiphenyl	13.361	172	1072004	91.17	ng	0.00
79) Terphenyl-d14	20.076	244	1921933	82.60	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.566	88	28226	15.36	ng	90
3) Pyridine	3.960	79	83896	16.14	ng	96
4) n-Nitrosodimethylamine	3.878	42	45790	16.83	ng	# 75
6) Aniline	7.426	93	155659	23.18	ng	# 92
8) 2-Chlorophenol	7.673	128	156391	39.84	ng	94
9) Benzaldehyde	7.244	77	54113	13.53	ng	91
10) Phenol	7.285	94	76640	13.86	ng	90
11) bis(2-Chloroethyl)ether	7.520	93	188083	44.00	ng	91
12) 1,3-Dichlorobenzene	7.996	146	207510	40.94	ng	# 90
13) 1,4-Dichlorobenzene	8.143	146	205652	40.94	ng	97
14) 1,2-Dichlorobenzene	8.466	146	202814	43.29	ng	96
15) Benzyl Alcohol	8.337	79	129854	28.43	ng	# 86
16) 2,2'-oxybis(1-Chloropr...	8.631	45	275570	36.74	ng	98
17) 2-Methylphenol	8.543	107	109014	30.70	ng	# 88
18) Hexachloroethane	9.201	117	71094	37.52	ng	90
19) n-Nitroso-di-n-propyla...	8.913	70	165486	43.44	ng	97
20) 3+4-Methylphenols	8.866	107	131510	27.17	ng	94
22) Acetophenone	8.931	105	314220	44.53	ng	# 93
24) Nitrobenzene	9.312	77	253098	50.77	ng	# 88
25) Isophorone	9.835	82	444245	44.56	ng	95
26) 2-Nitrophenol	10.023	139	109234	60.62	ng	# 75
27) 2,4-Dimethylphenol	10.076	122	156636	43.48	ng	87
28) bis(2-Chloroethoxy)met...	10.317	93	252987	41.14	ng	98
29) 2,4-Dichlorophenol	10.558	162	200519	46.28	ng	98
30) 1,2,4-Trichlorobenzene	10.781	180	218286	42.20	ng	99
31) Naphthalene	10.969	128	576185	43.72	ng	100
32) Benzoic acid	10.153	122	32048	17.41	ng	# 79
33) 4-Chloroaniline	11.069	127	138165	23.67	ng	96
34) Hexachlorobutadiene	11.257	225	152449	40.20	ng	97
35) Caprolactam	11.845	113	10735m	7.31	ng	
36) 4-Chloro-3-methylphenol	12.180	107	191930	41.20	ng	90
37) 2-Methylnaphthalene	12.567	142	441429	45.18	ng	# 94
38) 1-Methylnaphthalene	12.785	142	411119m	45.34	ng	
40) 1,2,4,5-Tetrachloroben...	12.932	216	281779	47.09	ng	# 98
41) Hexachlorocyclopentadiene	12.914	237	330858	95.05	ng	99
43) 2,4,6-Trichlorophenol	13.161	196	190306	50.57	ng	95

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.237	196	194882	51.45	ng	91
46) 1,1'-Biphenyl	13.566	154	595768	46.55	ng	94
47) 2-Chloronaphthalene	13.613	162	462717	44.49	ng	97
48) 2-Nitroaniline	13.807	65	164700	59.65	ng	# 77
49) Acenaphthylene	14.453	152	710504	46.42	ng	99
50) Dimethylphthalate	14.183	163	638918	48.05	ng	99
51) 2,6-Dinitrotoluene	14.295	165	136934	63.37	ng	# 72
52) Acenaphthene	14.794	154	569981m	54.83	ng	
53) 3-Nitroaniline	14.624	138	79641	32.61	ng	# 73
54) 2,4-Dinitrophenol	14.829	184	167262	141.73	ng	# 74
55) Dibenzofuran	15.129	168	741555	47.25	ng	95
56) 4-Nitrophenol	14.923	139	79606	39.58	ng	# 71
57) 2,4-Dinitrotoluene	15.082	165	199797	69.82	ng	# 81
58) Fluorene	15.775	166	613119	48.78	ng	97
59) 2,3,4,6-Tetrachlorophenol	15.347	232	196479	51.90	ng	# 94
60) Diethylphthalate	15.540	149	613034	45.86	ng	97
61) 4-Chlorophenyl-phenyle...	15.770	204	340263	47.43	ng	99
62) 4-Nitroaniline	15.787	138	134094	49.46	ng	# 65
63) Azobenzene	16.057	77	579989	44.14	ng	90
65) 4,6-Dinitro-2-methylph...	15.846	198	127796	75.64	ng	89
66) n-Nitrosodiphenylamine	15.981	169	557181	45.57	ng	98
67) 4-Bromophenyl-phenylether	16.663	248	235717	45.83	ng	97
68) Hexachlorobenzene	16.780	284	246642	44.57	ng	92
69) Atrazine	16.927	200	183436	37.34	ng	96
70) Pentachlorophenol	17.121	266	348598	109.13	ng	87
71) Phenanthrene	17.520	178	1026366	46.44	ng	97
72) Anthracene	17.609	178	1058478	48.54	ng	98
73) Carbazole	17.873	167	970359	48.90	ng	98
74) Di-n-butylphthalate	18.431	149	1060932	43.50	ng	# 98
75) Fluoranthene	19.524	202	1371804	48.74	ng	97
77) Benzidine	19.694	184	508874	35.40	ng	99
78) Pyrene	19.882	202	1399126	46.68	ng	99
80) Butylbenzylphthalate	20.764	149	532628	45.98	ng	90
81) Benzo(a)anthracene	21.733	228	1438671	46.21	ng	99
82) 3,3'-Dichlorobenzidine	21.639	252	325196	26.91	ng	96
83) Chrysene	21.798	228	1397125	46.97	ng	98
84) Bis(2-ethylhexyl)phtha...	21.633	149	763388	44.86	ng	95
85) Di-n-octyl phthalate	22.867	149	1338815	45.76	ng	# 95
87) Indeno(1,2,3-cd)pyrene	28.760	276	1726921	45.82	ng	# 90
88) Benzo(b)fluoranthene	23.983	252	1523677	45.64	ng	# 97
89) Benzo(k)fluoranthene	24.060	252	1469292	45.78	ng	# 98
90) Benzo(a)pyrene	24.871	252	1384831	44.09	ng	# 98
91) Dibenzo(a,h)anthracene	28.831	278	1424058	46.20	ng	# 96
92) Benzo(g,h,i)perylene	29.929	276	1442854	47.86	ng	# 93

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

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