

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102720\
 Data File : BG047097.D
 Acq On : 27 Oct 2020 16:24
 Operator : CG/JU
 Sample : L4214-08
 Misc : LOQ-WATER 5ppm
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LOQ-WATER-02-QT4-2020

Manual Integrations
 APPROVED

mohammad
 10/28/2020 8:06:13 AM

Quant Time: Oct 27 17:00:36 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG102720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 27 12:36:22 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.056	152	39654	20.00	ng	# 0.00
21) Naphthalene-d8	10.864	136	148237	20.00	ng	# 0.00
39) Acenaphthene-d10	14.684	164	109260	20.00	ng	0.00
64) Phenanthrene-d10	17.433	188	278515	20.00	ng	# 0.00
76) Chrysene-d12	21.699	240	339868	20.00	ng	# 0.00
86) Perylene-d12	24.924	264	338175	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.612	112	332714	148.74	ng	0.00
7) Phenol-d6	7.210	99	458881	141.85	ng	0.00
23) Nitrobenzene-d5	9.219	82	328892	106.25	ng	0.00
42) 2,4,6-Tribromophenol	16.170	330	266091	151.94	ng	0.00
45) 2-Fluorobiphenyl	13.309	172	825767	106.85	ng	0.00
79) Terphenyl-d14	20.036	244	1819596	104.10	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.502	88	5541	5.48	ng	91
3) Pyridine	3.902	79	13868	5.06	ng	# 91
4) n-Nitrosodimethylamine	3.814	42	7262	5.00	ng	# 63
6) Aniline	7.374	93	19923	5.21	ng	97
8) 2-Chlorophenol	7.615	128	12435	5.23	ng	92
9) Benzaldehyde	7.192	77	13074	7.21	ng	93
10) Phenol	7.233	94	14718	4.82	ng	# 72
11) bis(2-Chloroethyl)ether	7.468	93	11893	4.95	ng	95
12) 1,3-Dichlorobenzene	7.944	146	15573	5.14	ng	# 89
13) 1,4-Dichlorobenzene	8.091	146	15647m	5.11	ng	
14) 1,2-Dichlorobenzene	8.409	146	14899	5.14	ng	91
15) Benzyl Alcohol	8.291	79	11249	4.46	ng	# 78
16) 2,2'-oxybis(1-Chloropr...	8.591	45	17395	4.57	ng	88
17) 2-Methylphenol	8.491	107	10275	4.89	ng	# 83
18) Hexachloroethane	9.137	117	5857	5.11	ng	# 68
19) n-Nitroso-di-n-propyla...	8.855	70	11103	5.04	ng	# 91
20) 3+4-Methylphenols	8.820	107	12541	4.43	ng	91
22) Acetophenone	8.879	105	21723	5.44	ng	# 93
24) Nitrobenzene	9.260	77	18479	5.75	ng	# 91
25) Isophorone	9.783	82	29983	5.42	ng	# 91
26) 2-Nitrophenol	9.983	139	6470	4.64	ng	# 83
27) 2,4-Dimethylphenol	10.024	122	10967	5.76	ng	# 93
28) bis(2-Chloroethoxy)met...	10.259	93	16958	4.96	ng	# 91
29) 2,4-Dichlorophenol	10.512	162	11741	4.48	ng	95
30) 1,2,4-Trichlorobenzene	10.729	180	16086	4.98	ng	95
31) Naphthalene	10.911	128	41683	5.38	ng	96
33) 4-Chloroaniline	11.023	127	15788	4.68	ng	# 93
34) Hexachlorobutadiene	11.199	225	12366	5.10	ng	93
35) Caprolactam	11.763	113	3889	4.45	ng	# 82
36) 4-Chloro-3-methylphenol	12.139	107	12358	4.54	ng	96
37) 2-Methylnaphthalene	12.521	142	30911	5.13	ng	95
38) 1-Methylnaphthalene	12.733	142	27966	4.90	ng	# 95
40) 1,2,4,5-Tetrachloroben...	12.886	216	20498	5.14	ng	97
41) Hexachlorocyclopentadiene	12.862	237	10890	4.99	ng	94
43) 2,4,6-Trichlorophenol	13.121	196	12221	4.70	ng	95
44) 2,4,5-Trichlorophenol	13.197	196	12711	4.81	ng	91

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) 1,1'-Biphenyl	13.520	154	43486	5.26	ng	96
47) 2-Chloronaphthalene	13.561	162	33373	4.90	ng	98
48) 2-Nitroaniline	13.761	65	8903	4.01	ng	84
49) Acenaphthylene	14.407	152	49688	5.03	ng	98
50) Dimethylphthalate	14.131	163	45777	5.09	ng	99
51) 2,6-Dinitrotoluene	14.255	165	9092	4.84	ng	# 77
52) Acenaphthene	14.748	154	33790	5.24	ng	94
53) 3-Nitroaniline	14.578	138	8149	4.37	ng	# 97
54) 2,4-Dinitrophenol	14.795	184	3591	2.90	ng	# 87
55) Dibenzofuran	15.083	168	53849	5.20	ng	96
56) 4-Nitrophenol	14.883	139	5629	3.86	ng	# 62
57) 2,4-Dinitrotoluene	15.042	165	13288	4.93	ng	# 84
58) Fluorene	15.729	166	43648	5.22	ng	96
59) 2,3,4,6-Tetrachlorophenol	15.306	232	13645	4.86	ng	# 95
60) Diethylphthalate	15.494	149	47591	5.31	ng	# 93
61) 4-Chlorophenyl-phenyle...	15.723	204	25604	5.13	ng	92
62) 4-Nitroaniline	15.741	138	8779	4.38	ng	# 78
63) Azobenzene	16.011	77	43631	5.47	ng	86
65) 4,6-Dinitro-2-methylph...	15.800	198	8612	5.03	ng	85
66) n-Nitrosodiphenylamine	15.935	169	42091	5.27	ng	95
67) 4-Bromophenyl-phenylether	16.617	248	18963	5.17	ng	88
68) Hexachlorobenzene	16.734	284	18190	4.71	ng	95
69) Atrazine	16.875	200	17537	5.33	ng	95
70) Pentachlorophenol	17.087	266	8403	3.73	ng	# 86
71) Phenanthrene	17.474	178	77391	5.20	ng	99
72) Anthracene	17.562	178	80486	5.52	ng	98
73) Carbazole	17.833	167	68063	5.25	ng	94
74) Di-n-butylphthalate	18.385	149	89134	5.56	ng	96
75) Fluoranthene	19.478	202	106204	5.55	ng	96
77) Benzidine	19.654	184	42826	5.51	ng	# 93
78) Pyrene	19.842	202	112505	5.19	ng	98
80) Butylbenzylphthalate	20.718	149	37804	4.55	ng	# 84
81) Benzo(a)anthracene	21.675	228	116145	5.22	ng	97
82) 3,3'-Dichlorobenzidine	21.593	252	41216	5.10	ng	# 99
83) Chrysene	21.740	228	114199m	5.39	ng	
84) Bis(2-ethylhexyl)phtha...	21.575	149	55168	4.71	ng	# 96
85) Di-n-octyl phthalate	22.792	149	91577	4.66	ng	# 96
87) Indeno(1,2,3-cd)pyrene	28.602	276	120225	4.87	ng	# 65
88) Benzo(b)fluoranthene	23.896	252	111531	5.05	ng	# 95
89) Benzo(k)fluoranthene	23.967	252	113142	5.27	ng	# 92
90) Benzo(a)pyrene	24.766	252	98028	4.75	ng	# 96
91) Dibenzo(a,h)anthracene	28.673	278	99527	4.96	ng	# 90
92) Benzo(g,h,i)perylene	29.754	276	100091	5.02	ng	# 93

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

