

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110520\
 Data File : BG047225.D
 Acq On : 5 Nov 2020 18:21
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Nov 06 02:17:59 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG110420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 04 16:19:40 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.004	152	49834	20.00	ng	# 0.00
21) Naphthalene-d8	10.818	136	189492	20.00	ng	0.00
39) Acenaphthene-d10	14.643	164	135238	20.00	ng	0.00
64) Phenanthrene-d10	17.387	188	333166	20.00	ng	# 0.00
76) Chrysene-d12	21.646	240	345804	20.00	ng	# 0.00
86) Perylene-d12	24.831	264	362498	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.559	112	246790	80.27	ng	0.00
7) Phenol-d6	7.163	99	361039	81.12	ng	0.00
23) Nitrobenzene-d5	9.173	82	362183	82.03	ng	0.00
42) 2,4,6-Tribromophenol	16.129	330	197461	87.19	ng	0.00
45) 2-Fluorobiphenyl	13.262	172	855458	81.13	ng	0.00
79) Terphenyl-d14	19.995	244	1650349	83.93	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.432	88	56249	36.94	ng	# 94
3) Pyridine	3.832	79	157969	38.79	ng	94
4) n-Nitrosodimethylamine	3.750	42	82987	40.64	ng	# 74
6) Aniline	7.328	93	210785	40.12	ng	94
8) 2-Chlorophenol	7.569	128	129975	39.82	ng	93
9) Benzaldehyde	7.140	77	122397	40.88	ng	96
10) Phenol	7.187	94	169674	40.19	ng	76
11) bis(2-Chloroethyl)ether	7.422	93	133989	40.01	ng	95
12) 1,3-Dichlorobenzene	7.898	146	165901	39.05	ng	# 93
13) 1,4-Dichlorobenzene	8.045	146	168513	39.50	ng	97
14) 1,2-Dichlorobenzene	8.362	146	157673	38.97	ng	95
15) Benzyl Alcohol	8.239	79	143656	42.00	ng	89
16) 2,2'-oxybis(1-Chloropr...	8.538	45	204031	39.48	ng	97
17) 2-Methylphenol	8.444	107	113120	39.59	ng	# 87
18) Hexachloroethane	9.091	117	62005	39.70	ng	91
19) n-Nitroso-di-n-propyla...	8.808	70	117517	40.18	ng	91
20) 3+4-Methylphenols	8.773	107	161285	42.14	ng	89
22) Acetophenone	8.826	105	216663	39.65	ng	92
24) Nitrobenzene	9.214	77	177302	39.46	ng	# 86
25) Isophorone	9.731	82	307145	40.60	ng	# 95
26) 2-Nitrophenol	9.925	139	79206	40.88	ng	# 74
27) 2,4-Dimethylphenol	9.978	122	106289	40.85	ng	88
28) bis(2-Chloroethoxy)met...	10.219	93	185326	40.47	ng	96
29) 2,4-Dichlorophenol	10.460	162	151144	42.04	ng	97
30) 1,2,4-Trichlorobenzene	10.677	180	180188	40.66	ng	97
31) Naphthalene	10.865	128	432955	39.90	ng	98
32) Benzoic acid	10.113	122	82584	37.17	ng	# 80
33) 4-Chloroaniline	10.971	127	185174	40.57	ng	# 91
34) Hexachlorobutadiene	11.153	225	134714	40.17	ng	97
35) Caprolactam	11.740	113	51586	43.91	ng	96
36) 4-Chloro-3-methylphenol	12.093	107	157094	42.87	ng	91
37) 2-Methylnaphthalene	12.469	142	334048	40.97	ng	# 94
38) 1-Methylnaphthalene	12.692	142	317781	41.01	ng	95
40) 1,2,4,5-Tetrachloroben...	12.839	216	219186	41.05	ng	98
41) Hexachlorocyclopentadiene	12.816	237	114877	41.51	ng	96
43) 2,4,6-Trichlorophenol	13.074	196	146325	42.15	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.145	196	150371	42.66	ng	# 95
46) 1,1'-Biphenyl	13.474	154	449248	40.58	ng	95
47) 2-Chloronaphthalene	13.521	162	374185	41.43	ng	98
48) 2-Nitroaniline	13.720	65	126110	42.13	ng	# 81
49) Acenaphthylene	14.367	152	539647	40.85	ng	98
50) Dimethylphthalate	14.096	163	495697	41.50	ng	100
51) 2,6-Dinitrotoluene	14.214	165	107398	43.36	ng	# 69
52) Acenaphthene	14.707	154	353820	41.90	ng	97
53) 3-Nitroaniline	14.543	138	106999	43.39	ng	# 76
54) 2,4-Dinitrophenol	14.749	184	67803	42.66	ng	# 72
55) Dibenzofuran	15.042	168	564896	40.68	ng	96
56) 4-Nitrophenol	14.843	139	83836	43.89	ng	# 59
57) 2,4-Dinitrotoluene	15.001	165	152787	42.24	ng	# 81
58) Fluorene	15.689	166	466383	41.79	ng	97
59) 2,3,4,6-Tetrachlorophenol	15.266	232	156468	42.28	ng	# 96
60) Diethylphthalate	15.454	149	497364	42.16	ng	96
61) 4-Chlorophenyl-phenyle...	15.683	204	276070	41.77	ng	99
62) 4-Nitroaniline	15.706	138	115695	43.69	ng	# 60
63) Azobenzene	15.971	77	436163	41.25	ng	89
65) 4,6-Dinitro-2-methylph...	15.765	198	96043	41.46	ng	84
66) n-Nitrosodiphenylamine	15.894	169	412603	40.26	ng	99
67) 4-Bromophenyl-phenylether	16.576	248	195773	42.07	ng	98
68) Hexachlorobenzene	16.693	284	206207	40.79	ng	95
69) Atrazine	16.840	200	181518	41.41	ng	98
70) Pentachlorophenol	17.034	266	120304	43.67	ng	94
71) Phenanthrene	17.428	178	794728	41.10	ng	98
72) Anthracene	17.522	178	766391	40.80	ng	97
73) Carbazole	17.786	167	684051	41.25	ng	97
74) Di-n-butylphthalate	18.344	149	849428	41.89	ng	# 97
75) Fluoranthene	19.437	202	1010727	41.04	ng	97
77) Benzidine	19.613	184	482369	41.59	ng	97
78) Pyrene	19.796	202	1013896	42.10	ng	97
80) Butylbenzylphthalate	20.677	149	396529	43.65	ng	92
81) Benzo(a)anthracene	21.629	228	1020876	41.65	ng	98
82) 3,3'-Dichlorobenzidine	21.541	252	363326	42.84	ng	# 99
83) Chrysene	21.693	228	984387	41.72	ng	99
84) Bis(2-ethylhexyl)phtha...	21.529	149	550054	43.29	ng	95
85) Di-n-octyl phthalate	22.727	149	912358	42.72	ng	# 96
87) Indeno(1,2,3-cd)pyrene	28.474	276	1150361	41.81	ng	# 93
88) Benzo(b)fluoranthene	23.826	252	1054570	42.61	ng	# 97
89) Benzo(k)fluoranthene	23.891	252	984639	40.73	ng	# 96
90) Benzo(a)pyrene	24.684	252	963554	41.72	ng	# 98
91) Dibenzo(a,h)anthracene	28.538	278	938957	41.93	ng	# 96
92) Benzo(g,h,i)perylene	29.614	276	920481	41.45	ng	# 94

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

