

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060320\
 Data File : BG045595.D
 Acq On : 4 Jun 2020 2:24
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 6/4/2020 12:11:38 PM

Quant Time: Jun 04 04:17:00 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG051520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 01 15:07:00 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.95	152	70441	20.00	ng	0.00
21) Naphthalene-d8	10.76	136	293474	20.00	ng	0.00
39) Acenaphthene-d10	14.60	164	213231	20.00	ng	0.00
64) Phenanthrene-d10	17.35	188	492444	20.00	ng	0.00
76) Chrysene-d12	21.61	240	442697	20.00	ng	0.00
87) Perylene-d12	24.75	264	480305	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.54	112	311753	86.49	ng	0.00
7) Phenol-d6	7.15	99	465395	90.72	ng	0.00
23) Nitrobenzene-d5	9.12	82	458757	85.24	ng	0.00
42) 2,4,6-Tribromophenol	16.09	330	219760	80.59	ng	0.00
45) 2-Fluorobiphenyl	13.22	172	1039046	82.65	ng	0.00
79) Terphenyl-d14	19.96	244	1623571	85.54	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.41	88	69376	38.814	ng	97
3) Pyridine	3.80	79	205973	41.377	ng	99
4) n-Nitrosodimethylamine	3.72	42	71578	43.941	ng	88
6) Aniline	7.28	93	290371	43.359	ng	98
8) 2-Chlorophenol	7.52	128	173533	43.902	ng	98
9) Benzaldehyde	7.08	77	135839	40.641	ng	97
10) Phenol	7.18	94	237308	44.882	ng	96
11) bis(2-Chloroethyl)ether	7.37	93	178856	43.368	ng	98
12) 1,3-Dichlorobenzene	7.84	146	205426	43.247	ng	97
13) 1,4-Dichlorobenzene	7.99	146	208608	44.502	ng	96
14) 1,2-Dichlorobenzene	8.31	146	198806	44.095	ng	97
15) Benzyl Alcohol	8.20	79	193481	45.654	ng	95
16) 2,2'-oxybis(1-Chloropropan	8.48	45	172450	44.052	ng	99
17) 2-Methylphenol	8.42	107	178220	45.033	ng	99
18) Hexachloroethane	9.04	117	80793	45.001	ng	85
19) n-Nitroso-di-n-propylamine	8.76	70	158961	44.808	ng	99
20) 3+4-Methylphenols	8.75	107	241031	44.725	ng	93
22) Acetophenone	8.77	105	292521	42.055	ng	97
24) Nitrobenzene	9.16	77	246691	42.784	ng	98
25) Isophorone	9.68	82	449538	42.415	ng	100
26) 2-Nitrophenol	9.87	139	108041	43.802	ng	97
27) 2,4-Dimethylphenol	9.95	122	156726	42.841	ng	96
28) bis(2-Chloroethoxy)methane	10.17	93	237936	42.807	ng	95
29) 2,4-Dichlorophenol	10.43	162	189215	43.799	ng	97
30) 1,2,4-Trichlorobenzene	10.63	180	224683	44.014	ng	97
31) Naphthalene	10.81	128	583592	42.674	ng	99
32) Benzoic acid	10.13	122	100971	42.131	ng	92
33) 4-Chloroaniline	10.92	127	246112	43.053	ng	97
34) Hexachlorobutadiene	11.10	225	159732	44.267	ng	97
35) Caprolactam	11.69	113	65785	41.487	ng	86
36) 4-Chloro-3-methylphenol	12.07	107	214495	44.062	ng	95
37) 2-Methylnaphthalene	12.42	142	443296	43.490	ng	99
38) 1-Methylnaphthalene	12.64	142	413271	42.921	ng	98
40) 1,2,4,5-Tetrachlorobenzene	12.79	216	249014	41.810	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.78	237	109038	31.719	nq	97
43) 2,4,6-Trichlorophenol	13.04	196	173663	41.975	nq	98
44) 2,4,5-Trichlorophenol	13.12	196	192024	40.615	nq	98
46) 1,1'-Biphenyl	13.43	154	539391	41.168	nq	98
47) 2-Chloronaphthalene	13.47	162	436474	41.024	nq	99
48) 2-Nitroaniline	13.67	65	143928	41.125	nq	94
49) Acenaphthylene	14.32	152	703284	41.153	nq	99
50) Dimethylphthalate	14.06	163	587874	40.190	nq	100
51) 2,6-Dinitrotoluene	14.17	165	133979	40.591	nq	96
52) Acenaphthene	14.66	154	458883m	40.114	nq	
53) 3-Nitroaniline	14.51	138	135381	40.348	nq	99
54) 2,4-Dinitrophenol	14.71	184	63212	32.699	nq	96
55) Dibenzofuran	15.00	168	696044	40.973	nq	99
56) 4-Nitrophenol	14.85	139	92359	36.358	nq	91
57) 2,4-Dinitrotoluene	14.96	165	192003	40.166	nq	98
58) Fluorene	15.65	166	575936	41.267	nq	97
59) 2,3,4,6-Tetrachlorophenol	15.24	232	165740	39.846	nq	98
60) Diethylphthalate	15.42	149	605541	39.773	nq	98
61) 4-Chlorophenyl-phenylether	15.64	204	338050	42.329	nq	99
62) 4-Nitroaniline	15.67	138	140212	40.028	nq	96
63) Azobenzene	15.94	77	583601	41.109	nq	98
65) 4,6-Dinitro-2-methylphenol	15.73	198	109645	38.231	nq	95
66) n-Nitrosodiphenylamine	15.85	169	500772	42.354	nq	97
67) 4-Bromophenyl-phenylether	16.53	248	233080	43.710	nq	94
68) Hexachlorobenzene	16.66	284	238447	42.754	nq	97
69) Atrazine	16.80	200	189467	38.905	nq	98
70) Pentachlorophenol	17.01	266	172983	59.734	nq	97
71) Phenanthrene	17.39	178	901415	41.930	nq	100
72) Anthracene	17.48	178	905022	41.921	nq	98
73) Carbazole	17.75	167	864028	41.058	nq	100
74) Di-n-butylphthalate	18.31	149	1016820	40.257	nq	99
75) Fluoranthene	19.40	202	1084831	39.673	nq	100
77) Benzidine	19.58	184	463308	37.263	nq	99
78) Pyrene	19.76	202	1075093	42.602	nq	100
80) Butylbenzylphthalate	20.65	149	450942	41.399	nq	98
81) Benzo(a)anthracene	21.59	228	1047929	42.493	nq	99
82) 3,3'-Dichlorobenzidine	21.50	252	395210	40.854	nq	100
83) Chrysene	21.65	228	999817	42.164	nq	100
84) Bis(2-ethylhexyl)phthalate	21.50	149	625506	41.775	nq	99
85) Di-n-octyl phthalate	22.69	149	1051141	40.874	nq	100
86) Indeno(1,2,3-cd)pyrene	28.33	276	1283689	41.398	nq	# 96
88) Benzo(b)fluoranthene	23.76	252	1105942	42.934	nq	99
89) Benzo(k)fluoranthene	23.82	252	1048593	43.329	nq	98
90) Benzo(a)pyrene	24.61	252	1031239	42.986	nq	98
91) Dibenzo(a,h)anthracene	28.40	278	1039191	43.106	nq	97
92) Benzo(g,h,i)perylene	29.45	276	1018578	42.245	nq	97

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

