

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120418\
 Data File : BG038327.D
 Acq On : 4 Dec 2018 13:22
 Operator : JU/SJ
 Sample : SSTDICCC040
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICCC040

Manual Integrations
 APPROVED

mohammad
 12/6/2018 10:32:47 AM

Quant Time: Dec 04 14:14:16 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG120418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 04 13:58:38 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.07	152	42136	20.00	ng	0.00
21) Naphthalene-d8	10.89	136	167330	20.00	ng	0.00
39) Acenaphthene-d10	14.71	164	104314	20.00	ng	0.00
64) Phenanthrene-d10	17.46	188	240648	20.00	ng	0.00
76) Chrysene-d12	21.74	240	230515	20.00	ng	0.00
87) Perylene-d12	25.05	264	249368	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.62	112	188388	77.19	ng	0.00
7) Phenol-d6	7.22	99	279246	80.16	ng	0.00
23) Nitrobenzene-d5	9.25	82	282149	90.26	ng	0.00
42) 2,4,6-Tribromophenol	16.20	330	107110	90.03	ng	0.00
45) 2-Fluorobiphenyl	13.33	172	583738	81.52	ng	0.00
79) Terphenyl-d14	20.06	244	832038	84.93	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.51	88	40483	35.119	ng	96
3) Pyridine	3.91	79	130823	39.785	ng	92
4) n-Nitrosodimethylamine	3.83	42	56051	46.985	ng	# 99
6) Aniline	7.40	93	178750	39.757	ng	99
8) 2-Chlorophenol	7.63	128	110882	39.157	ng	96
9) Benzaldehyde	7.21	77	83101	32.986	ng	97
10) Phenol	7.25	94	147114	39.869	ng	97
11) bis(2-Chloroethyl)ether	7.49	93	122642	40.712	ng	100
12) 1,3-Dichlorobenzene	7.96	146	131138	40.199	ng	97
13) 1,4-Dichlorobenzene	8.11	146	132304	40.148	ng	98
14) 1,2-Dichlorobenzene	8.43	146	121574	38.642	ng	100
15) Benzyl Alcohol	8.31	79	109383	43.394	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.59	45	250563	44.794	ng	98
17) 2-Methylphenol	8.51	107	97449	39.063	ng	97
18) Hexachloroethane	9.15	117	48858	40.738	ng	98
19) n-Nitroso-di-n-propylamine	8.88	70	93439	37.069	ng	97
20) 3+4-Methylphenols	8.83	107	136581	38.621	ng	95
22) Acetophenone	8.91	105	168205	40.402	ng	# 95
24) Nitrobenzene	9.29	77	142864	44.678	ng	97
25) Isophorone	9.81	82	334116	59.385	ng	97
26) 2-Nitrophenol	10.00	139	77905	48.802	ng	97
27) 2,4-Dimethylphenol	10.05	122	108875	45.267	ng	96
28) bis(2-Chloroethoxy)methane	10.29	93	150839	41.926	ng	98
29) 2,4-Dichlorophenol	10.53	162	110100	40.696	ng	98
30) 1,2,4-Trichlorobenzene	10.75	180	123588	40.766	ng	98
31) Naphthalene	10.94	128	326210	39.636	ng	99
32) Benzoic acid	10.20	122	68192m	55.238	ng	
33) 4-Chloroaniline	11.06	127	149513	39.054	ng	100
34) Hexachlorobutadiene	11.20	225	79958	42.854	ng	98
35) Caprolactam	11.85	113	43959	40.595	ng	95
36) 4-Chloro-3-methylphenol	12.16	107	122540	41.460	ng	100
37) 2-Methylnaphthalene	12.54	142	234889	39.733	ng	98
38) 1-Methylnaphthalene	12.76	142	226452	39.795	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.90	216	146592	42.055	ng	99

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41) Hexachlorocyclopentadiene	12.87	237	85778	45.956	nq	96
43) 2,4,6-Trichlorophenol	13.14	196	98562	41.495	nq	100
44) 2,4,5-Trichlorophenol	13.22	196	101338	40.548	nq	98
46) 1,1'-Biphenyl	13.55	154	349178	39.844	nq	99
47) 2-Chloronaphthalene	13.59	162	263183	39.356	nq	99
48) 2-Nitroaniline	13.81	65	91248	43.355	nq	97
49) Acenaphthylene	14.43	152	406993	40.472	nq	99
50) Dimethylphthalate	14.16	163	340211	41.139	nq	99
51) 2,6-Dinitrotoluene	14.29	165	78934	42.173	nq	98
52) Acenaphthene	14.77	154	240607	39.743	nq	99
53) 3-Nitroaniline	14.63	138	84911	41.113	nq	# 96
54) 2,4-Dinitrophenol	14.83	184	43431	55.072	nq	# 84
55) Dibenzofuran	15.11	168	403913	40.348	nq	98
56) 4-Nitrophenol	14.92	139	66172	41.542	nq	92
57) 2,4-Dinitrotoluene	15.07	165	107977	42.401	nq	96
58) Fluorene	15.76	166	302961	40.561	nq	97
59) 2,3,4,6-Tetrachlorophenol	15.33	232	89835	42.957	nq	98
60) Diethylphthalate	15.52	149	365834	41.649	nq	99
61) 4-Chlorophenyl-phenylether	15.74	204	180607	41.324	nq	98
62) 4-Nitroaniline	15.79	138	83627	40.760	nq	90
63) Azobenzene	16.04	77	327037	41.642	nq	96
65) 4,6-Dinitro-2-methylphenol	15.84	198	68385	44.928	nq	96
66) n-Nitrosodiphenylamine	15.96	169	295247	40.554	nq	100
67) 4-Bromophenyl-phenylether	16.64	248	114056	43.181	nq	97
68) Hexachlorobenzene	16.75	284	116621	42.775	nq	95
69) Atrazine	16.90	200	110535	41.187	nq	100
70) Pentachlorophenol	17.10	266	81728	44.138	nq	91
71) Phenanthrene	17.50	178	498817	40.486	nq	98
72) Anthracene	17.59	178	495061	40.762	nq	100
73) Carbazole	17.87	167	498596	40.918	nq	99
74) Di-n-butylphthalate	18.39	149	602417	44.042	nq	98
75) Fluoranthene	19.50	202	616264	43.840	nq	99
77) Benzidine	19.69	184	317529	39.256	nq	99
78) Pyrene	19.87	202	616691	40.918	nq	99
80) Butylbenzylphthalate	20.74	149	260435	39.514	nq	97
81) Benzo(a)anthracene	21.72	228	573350	41.159	nq	99
82) 3,3'-Dichlorobenzidine	21.64	252	230209	42.424	nq	100
83) Chrysene	21.79	228	545545	40.932	nq	99
84) Bis(2-ethylhexyl)phthalate	21.59	149	382691	40.715	nq	99
85) Di-n-octyl phthalate	22.82	149	670733	44.109	nq	99
86) Indeno(1,2,3-cd)pyrene	28.85	276	638129	43.108	nq	# 95
88) Benzo(b)fluoranthene	23.99	252	551639	40.199	nq	99
89) Benzo(k)fluoranthene	24.06	252	547219	40.412	nq	100
90) Benzo(a)pyrene	24.89	252	632475	48.227	nq	99
91) Dibenzo(a,h)anthracene	28.91	278	536099	42.485	nq	99
92) Benzo(g,h,i)perylene	30.05	276	494391	39.404	nq	# 94

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

