

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG113018\
 Data File : BG038252.D
 Acq On : 30 Nov 2018 11:27
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 12/3/2018 3:32:55 PM

Quant Time: Nov 30 13:14:21 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG111318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 13 16:39:28 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.08	152	46114	20.00	ng	0.00
21) Naphthalene-d8	10.90	136	210118	20.00	ng	0.00
39) Acenaphthene-d10	14.72	164	128489	20.00	ng	0.00
64) Phenanthrene-d10	17.47	188	286499	20.00	ng	0.00
76) Chrysene-d12	21.75	240	266527	20.00	ng	0.00
87) Perylene-d12	25.07	264	281995	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.63	112	223365	83.63	ng	0.00
7) Phenol-d6	7.23	99	327380	85.87	ng	0.00
23) Nitrobenzene-d5	9.26	82	336306	85.68	ng	0.00
42) 2,4,6-Tribromophenol	16.21	330	119525	81.57	ng	0.00
45) 2-Fluorobiphenyl	13.34	172	729552	82.72	ng	0.00
79) Terphenyl-d14	20.07	244	951031	83.96	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.52	88	45509	36.074	ng	97
3) Pyridine	3.92	79	136020	37.797	ng	90
4) n-Nitrosodimethylamine	3.85	42	63970	48.997	ng	# 89
6) Aniline	7.41	93	208637	42.401	ng	98
8) 2-Chlorophenol	7.64	128	133103	42.949	ng	98
9) Benzaldehyde	7.22	77	104628	37.949	ng	97
10) Phenol	7.26	94	173303	42.915	ng	94
11) bis(2-Chloroethyl)ether	7.50	93	139837	42.416	ng	100
12) 1,3-Dichlorobenzene	7.97	146	146698	41.090	ng	96
13) 1,4-Dichlorobenzene	8.12	146	147553	40.913	ng	99
14) 1,2-Dichlorobenzene	8.44	146	139095	40.397	ng	95
15) Benzyl Alcohol	8.32	79	130430	47.280	ng	93
16) 2,2'-oxybis(1-Chloropropan	8.60	45	299771	48.968	ng	98
17) 2-Methylphenol	8.52	107	119807	43.882	ng	97
18) Hexachloroethane	9.17	117	56238	42.847	ng	97
19) n-Nitroso-di-n-propylamine	8.89	70	117806	42.704	ng	95
20) 3+4-Methylphenols	8.85	107	170266	43.993	ng	97
22) Acetophenone	8.92	105	211443	40.445	ng	# 96
24) Nitrobenzene	9.30	77	173763	43.275	ng	98
25) Isophorone	9.82	82	291640	41.280	ng	99
26) 2-Nitrophenol	10.01	139	85504	42.655	ng	95
27) 2,4-Dimethylphenol	10.06	122	126922	42.024	ng	98
28) bis(2-Chloroethoxy)methane	10.30	93	187679	41.543	ng	97
29) 2,4-Dichlorophenol	10.54	162	142654	41.992	ng	98
30) 1,2,4-Trichlorobenzene	10.76	180	157145	41.280	ng	97
31) Naphthalene	10.96	128	421249	40.761	ng	99
32) Benzoic acid	10.18	122	52213m	32.283	ng	
33) 4-Chloroaniline	11.07	127	195057	40.575	ng	98
34) Hexachlorobutadiene	11.21	225	97056	41.425	ng	97
35) Caprolactam	11.87	113	53464	39.318	ng	# 87
36) 4-Chloro-3-methylphenol	12.17	107	158232	42.634	ng	96
37) 2-Methylnaphthalene	12.55	142	300201	40.440	ng	97
38) 1-Methylnaphthalene	12.77	142	289939	40.576	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.91	216	180416	42.020	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.88	237	81543	35.468	nq	97
43) 2,4,6-Trichlorophenol	13.15	196	120549	41.203	nq	96
44) 2,4,5-Trichlorophenol	13.23	196	121735	39.544	nq	98
46) 1,1'-Biphenyl	13.55	154	445899	41.308	nq	100
47) 2-Chloronaphthalene	13.60	162	335964	40.787	nq	97
48) 2-Nitroaniline	13.81	65	117130	45.181	nq	98
49) Acenaphthylene	14.45	152	509196	41.108	nq	99
50) Dimethylphthalate	14.17	163	415754	40.815	nq	100
51) 2,6-Dinitrotoluene	14.30	165	96528	41.870	nq	97
52) Acenaphthene	14.79	154	304365	40.815	nq	99
53) 3-Nitroaniline	14.64	138	105888	41.624	nq	# 97
54) 2,4-Dinitrophenol	14.84	184	10773	16.937	nq	# 82
55) Dibenzofuran	15.11	168	499976	40.547	nq	97
56) 4-Nitrophenol	14.93	139	67008	34.152	nq	95
57) 2,4-Dinitrotoluene	15.09	165	132339	42.190	nq	98
58) Fluorene	15.77	166	373292	40.574	nq	100
59) 2,3,4,6-Tetrachlorophenol	15.34	232	102323	39.723	nq	99
60) Diethylphthalate	15.52	149	438588	40.537	nq	99
61) 4-Chlorophenyl-phenylether	15.75	204	221324	41.113	nq	98
62) 4-Nitroaniline	15.80	138	105333	41.680	nq	93
63) Azobenzene	16.04	77	406968	42.070	nq	96
65) 4,6-Dinitro-2-methylphenol	15.84	198	37106	20.477	nq	94
66) n-Nitrosodiphenylamine	15.97	169	362713	41.848	nq	99
67) 4-Bromophenyl-phenylether	16.65	248	135567	43.111	nq	97
68) Hexachlorobenzene	16.76	284	140899	43.409	nq	96
69) Atrazine	16.91	200	133107	41.660	nq	97
70) Pentachlorophenol	17.11	266	82054	37.222	nq	94
71) Phenanthrene	17.51	178	606200	41.327	nq	99
72) Anthracene	17.60	178	598459	41.390	nq	99
73) Carbazole	17.88	167	603256	41.584	nq	99
74) Di-n-butylphthalate	18.41	149	699845	42.976	nq	99
75) Fluoranthene	19.52	202	694839	41.519	nq	99
77) Benzidine	19.70	184	372712	39.853	nq	99
78) Pyrene	19.88	202	713707	40.957	nq	99
80) Butylbenzylphthalate	20.75	149	317885	41.714	nq	99
81) Benzo(a)anthracene	21.73	228	667612	41.450	nq	99
82) 3,3'-Dichlorobenzidine	21.65	252	256900	40.946	nq	100
83) Chrysene	21.80	228	636685	41.315	nq	99
84) Bis(2-ethylhexyl)phthalate	21.60	149	461844	42.497	nq	99
85) Di-n-octyl phthalate	22.84	149	764993	43.511	nq	98
86) Indeno(1,2,3-cd)pyrene	28.88	276	643017	37.569	nq	# 92
88) Benzo(b)fluoranthene	24.01	252	654093	42.150	nq	99
89) Benzo(k)fluoranthene	24.08	252	625433	40.844	nq	99
90) Benzo(a)pyrene	24.92	252	627898	42.338	nq	99
91) Dibenzo(a,h)anthracene	28.95	278	527442	36.963	nq	98
92) Benzo(g,h,i)perylene	30.09	276	507545	35.772	nq	99

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

