

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120418\
 Data File : BG038329.D
 Acq On : 4 Dec 2018 14:41
 Operator : JU/SJ
 Sample : SSTDICC060
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC060

Manual Integrations
 APPROVED

mohammad
 12/6/2018 10:32:52 AM

Quant Time: Dec 04 15:20:56 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	39725	20.00	ng	0.00
21) Naphthalene-d8	10.89	136	180428	20.00	ng	0.00
39) Acenaphthene-d10	14.71	164	104773	20.00	ng	0.00
64) Phenanthrene-d10	17.46	188	284124	20.00	ng	0.00
76) Chrysene-d12	21.75	240	212723	20.00	ng	0.00
87) Perylene-d12	25.05	264	233107	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.62	112	283173	123.08	ng	0.00
7) Phenol-d6	7.23	99	400086	121.82	ng	0.00
23) Nitrobenzene-d5	9.25	82	398853	118.33	ng	0.00
42) 2,4,6-Tribromophenol	16.21	330	160646	134.44	ng	0.00
45) 2-Fluorobiphenyl	13.33	172	859743	119.55	ng	0.00
79) Terphenyl-d14	20.06	244	1151321	127.35	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.51	88	63795	58.702	ng	98
3) Pyridine	3.91	79	211773	68.312	ng	94
4) n-Nitrosodimethylamine	3.83	42	94113	83.679	ng	# 94
6) Aniline	7.40	93	259095	61.124	ng	99
8) 2-Chlorophenol	7.63	128	164199	61.504	ng	99
9) Benzaldehyde	7.22	77	110874	46.682	ng	99
10) Phenol	7.25	94	212541	61.097	ng	94
11) bis(2-Chloroethyl)ether	7.49	93	177585	62.529	ng	99
12) 1,3-Dichlorobenzene	7.96	146	189166	61.506	ng	98
13) 1,4-Dichlorobenzene	8.11	146	183264	58.988	ng	98
14) 1,2-Dichlorobenzene	8.43	146	169716	57.217	ng	98
15) Benzyl Alcohol	8.32	79	157011	66.069	ng	94
16) 2,2'-oxybis(1-Chloropropan	8.60	45	360563	68.371	ng	97
17) 2-Methylphenol	8.51	107	136572	58.068	ng	93
18) Hexachloroethane	9.15	117	69469	61.440	ng	94
19) n-Nitroso-di-n-propylamine	8.88	70	139125	58.544	ng	93
20) 3+4-Methylphenols	8.84	107	196027	58.795	ng	97
22) Acetophenone	8.91	105	250141	55.721	ng	96
24) Nitrobenzene	9.30	77	203312	58.966	ng	97
25) Isophorone	9.81	82	344714	56.821	ng	# 96
26) 2-Nitrophenol	10.01	139	104025	60.433	ng	94
27) 2,4-Dimethylphenol	10.05	122	146252	56.392	ng	97
28) bis(2-Chloroethoxy)methane	10.29	93	215923	55.660	ng	97
29) 2,4-Dichlorophenol	10.54	162	177330	60.788	ng	98
30) 1,2,4-Trichlorobenzene	10.75	180	197288	60.352	ng	95
31) Naphthalene	10.95	128	513134	57.822	ng	100
32) Benzoic acid	10.23	122	107567m	80.808	ng	
33) 4-Chloroaniline	11.06	127	238124	57.685	ng	98
34) Hexachlorobutadiene	11.21	225	130166	64.699	ng	97
35) Caprolactam	11.87	113	69916	59.878	ng	# 88
36) 4-Chloro-3-methylphenol	12.17	107	204878	64.285	ng	99
37) 2-Methylnaphthalene	12.54	142	350671	55.012	ng	98
38) 1-Methylnaphthalene	12.76	142	339898	55.395	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.90	216	221073	63.145	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.87	237	132318	70.580	nq	99
43) 2,4,6-Trichlorophenol	13.14	196	149066	62.482	nq	96
44) 2,4,5-Trichlorophenol	13.22	196	157109	62.587	nq	97
46) 1,1'-Biphenyl	13.54	154	518457	58.902	nq	99
47) 2-Chloronaphthalene	13.60	162	398716	59.362	nq	99
48) 2-Nitroaniline	13.80	65	139110	65.806	nq	96
49) Acenaphthylene	14.44	152	600458	59.448	nq	99
50) Dimethylphthalate	14.17	163	499455	60.131	nq	99
51) 2,6-Dinitrotoluene	14.30	165	113840	60.556	nq	95
52) Acenaphthene	14.78	154	359983	59.200	nq	98
53) 3-Nitroaniline	14.63	138	133821	64.511	nq #	98
54) 2,4-Dinitrophenol	14.83	184	68316	86.247	nq	89
55) Dibenzofuran	15.11	168	589904	58.669	nq	97
56) 4-Nitrophenol	14.92	139	96960	60.604	nq	87
57) 2,4-Dinitrotoluene	15.08	165	158719	62.053	nq	98
58) Fluorene	15.76	166	437406	58.305	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.33	232	132886	63.264	nq	97
60) Diethylphthalate	15.52	149	518987	58.826	nq	99
61) 4-Chlorophenyl-phenylether	15.74	204	268476	61.160	nq	98
62) 4-Nitroaniline	15.80	138	123501	59.931	nq	88
63) Azobenzene	16.04	77	475828	60.322	nq	96
65) 4,6-Dinitro-2-methylphenol	15.84	198	104686	58.253	nq	95
66) n-Nitrosodiphenylamine	15.96	169	429726	49.994	nq	99
67) 4-Bromophenyl-phenylether	16.64	248	169713	54.421	nq	97
68) Hexachlorobenzene	16.75	284	174247	54.132	nq	99
69) Atrazine	16.90	200	152841	48.236	nq	97
70) Pentachlorophenol	17.10	266	121390	55.526	nq	96
71) Phenanthrene	17.50	178	791931	54.440	nq	100
72) Anthracene	17.60	178	742836	51.804	nq	99
73) Carbazole	17.87	167	703173	48.877	nq	99
74) Di-n-butylphthalate	18.40	149	850841	52.686	nq	98
75) Fluoranthene	19.51	202	895897	53.980	nq	99
77) Benzidine	19.69	184	430627	57.692	nq	97
78) Pyrene	19.87	202	866858	62.328	nq	100
80) Butylbenzylphthalate	20.74	149	371513	61.082	nq	98
81) Benzo(a)anthracene	21.72	228	783553	60.953	nq	99
82) 3,3'-Dichlorobenzidine	21.64	252	310371	61.981	nq	99
83) Chrysene	21.79	228	743605	60.458	nq	99
84) Bis(2-ethylhexyl)phthalate	21.59	149	520626	60.023	nq	100
85) Di-n-octyl phthalate	22.82	149	862570	61.470	nq	97
86) Indeno(1,2,3-cd)pyrene	28.86	276	887946	65.002	nq #	94
88) Benzo(b)fluoranthene	24.00	252	770572	60.070	nq	98
89) Benzo(k)fluoranthene	24.07	252	776322	61.330	nq	98
90) Benzo(a)pyrene	24.90	252	747876	61.004	nq	97
91) Dibenzo(a,h)anthracene	28.93	278	733171	62.155	nq	98
92) Benzo(g,h,i)perylene	30.07	276	725165	61.828	nq	97

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

