

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG123019\
 Data File : BG043865.D
 Acq On : 30 Dec 2019 15:51
 Operator : JU
 Sample : SSTDICC100
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC100

Manual Integrations
 APPROVED

mohammad

12/31/2019 1:48:05 PM

Quant Time: Dec 30 17:19:56 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG123019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 30 14:25:12 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.97	152	179159	20.00	ng	0.00
21) Naphthalene-d8	10.77	136	712923	20.00	ng	0.00
39) Acenaphthene-d10	14.60	164	444403	20.00	ng	0.00
64) Phenanthrene-d10	17.34	188	826524	20.00	ng	0.00
76) Chrysene-d12	21.60	240	664541	20.00	ng	0.00
87) Perylene-d12	24.72	264	781073	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.54	112	1657919	169.19	ng	0.00
7) Phenol-d6	7.14	99	2273272	165.56	ng	0.01
23) Nitrobenzene-d5	9.13	82	2622905	203.95	ng	0.00
42) 2,4,6-Tribromophenol	16.09	330	849198	182.22	ng	0.00
45) 2-Fluorobiphenyl	13.23	172	4028067	162.48	ng	0.00
79) Terphenyl-d14	19.96	244	4004653	147.92	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.44	88	385385	90.528	ng	99
3) Pyridine	3.83	79	1089553	86.829	ng	99
4) n-Nitrosodimethylamine	3.75	42	511895	91.224	ng	99
6) Aniline	7.30	93	1572963	87.393	ng	96
8) 2-Chlorophenol	7.53	128	1003072	87.285	ng	96
9) Benzaldehyde	7.10	77	806815	100.383	ng	97
10) Phenol	7.16	94	1222370	86.541	ng	94
11) bis(2-Chloroethyl)ether	7.39	93	1066582	87.632	ng	96
12) 1,3-Dichlorobenzene	7.86	146	1168422	86.844	ng	96
13) 1,4-Dichlorobenzene	8.00	146	1160189	87.462	ng	95
14) 1,2-Dichlorobenzene	8.33	146	1105493	86.705	ng	98
15) Benzyl Alcohol	8.21	79	982471	90.943	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.50	45	1669200	88.493	ng	98
17) 2-Methylphenol	8.41	107	922741	90.414	ng	96
18) Hexachloroethane	9.05	117	467176	90.145	ng	97
19) n-Nitroso-di-n-propylamine	8.79	70	908390	89.023	ng	94
20) 3+4-Methylphenols	8.75	107	1264534	88.847	ng	97
22) Acetophenone	8.79	105	1571114	88.554	ng	# 93
24) Nitrobenzene	9.17	77	1222347	92.673	ng	97
25) Isophorone	9.70	82	2271636	90.877	ng	95
26) 2-Nitrophenol	9.88	139	640045	102.500	ng	96
27) 2,4-Dimethylphenol	9.95	122	890237	94.746	ng	100
28) bis(2-Chloroethoxy)methane	10.18	93	1481831	88.870	ng	96
29) 2,4-Dichlorophenol	10.42	162	1049126	93.726	ng	98
30) 1,2,4-Trichlorobenzene	10.63	180	1154453	91.814	ng	97
31) Naphthalene	10.82	128	2872473	83.273	ng	# 90
32) Benzoic acid	10.15	122	815901m	112.503	ng	
33) 4-Chloroaniline	10.92	127	1489935	90.684	ng	94
34) Hexachlorobutadiene	11.12	225	723503	94.283	ng	100
35) Caprolactam	11.73	113	396291m	95.231	ng	
36) 4-Chloro-3-methylphenol	12.06	107	1101327	92.042	ng	99
37) 2-Methylnaphthalene	12.43	142	2232015	85.951	ng	97
38) 1-Methylnaphthalene	12.64	142	2124814	86.325	ng	# 98
40) 1,2,4,5-Tetrachlorobenzene	12.80	216	1205026	92.607	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.78	237	692025	101.945	nq	99
43) 2,4,6-Trichlorophenol	13.03	196	859142	95.717	nq	98
44) 2,4,5-Trichlorophenol	13.11	196	877627	95.007	nq	98
46) 1,1'-Biphenyl	13.44	154	2621224	82.126	nq	84
47) 2-Chloronaphthalene	13.48	162	2215125	86.137	nq #	88
48) 2-Nitroaniline	13.67	65	794257	96.040	nq	97
49) Acenaphthylene	14.32	152	3012454	81.548	nq #	86
50) Dimethylphthalate	14.07	163	2563815	81.115	nq #	93
51) 2,6-Dinitrotoluene	14.17	165	668311	93.675	nq	98
52) Acenaphthene	14.66	154	2323553	89.648	nq	94
53) 3-Nitroaniline	14.50	138	734644	90.601	nq	98
54) 2,4-Dinitrophenol	14.70	184	368500	114.902	nq	97
55) Dibenzofuran	15.00	168	2818989	78.849	nq #	83
56) 4-Nitrophenol	14.81	139	576469	92.687	nq	100
57) 2,4-Dinitrotoluene	14.96	165	884613	91.116	nq	96
58) Fluorene	15.65	166	2283219	80.610	nq #	98
59) 2,3,4,6-Tetrachlorophenol	15.22	232	732025	92.124	nq	99
60) Diethylphthalate	15.43	149	2527328	79.936	nq #	90
61) 4-Chlorophenyl-phenylether	15.64	204	1329998	86.495	nq	95
62) 4-Nitroaniline	15.67	138	695532	86.933	nq	98
63) Azobenzene	15.93	77	2387254	81.498	nq	88
65) 4,6-Dinitro-2-methylphenol	15.72	198	445585	111.060	nq	95
66) n-Nitrosodiphenylamine	15.86	169	2136040	87.516	nq	91
67) 4-Bromophenyl-phenylether	16.53	248	885070	95.268	nq	98
68) Hexachlorobenzene	16.66	284	940066	93.775	nq	97
69) Atrazine	16.80	200	613416	77.368	nq	98
70) Pentachlorophenol	16.99	266	561567	101.566	nq	99
71) Phenanthrene	17.39	178	3110320	78.410	nq #	82
72) Anthracene	17.48	178	3070727	78.399	nq #	83
73) Carbazole	17.74	167	2859080	78.926	nq #	88
77) Benzidine	19.56	184	1314143	76.758	nq #	91
80) Butylbenzylphthalate	20.65	149	1824026	88.425	nq #	93
81) Benzo(a)anthracene	21.58	228	3392662	82.290	nq #	83
82) 3,3'-Dichlorobenzidine	21.49	252	1398680	85.437	nq	94
83) Chrysene	21.64	228	3229305	82.508	nq #	82
84) Bis(2-ethylhexyl)phthalate	21.50	149	2360670	83.093	nq #	87
85) Di-n-octyl phthalate	22.70	149	3937732	82.206	nq #	92
86) Indeno(1,2,3-cd)pyrene	28.30	276	5132787	96.465	nq #	92
88) Benzo(b)fluoranthene	23.74	252	4035171	87.313	nq #	91
89) Benzo(k)fluoranthene	23.81	252	3665060	83.660	nq #	90
90) Benzo(a)pyrene	24.59	252	3843548	88.245	nq	93
91) Dibenzo(a,h)anthracene	28.36	278	4000404	91.473	nq	96
92) Benzo(g,h,i)perylene	29.41	276	4091845	94.147	nq	99

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

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