

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG071818\
 Data File : BG035737.D
 Acq On : 19 Jul 2018 4:11
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 7/19/2018 3:45:28 PM

Quant Time: Jul 19 07:11:28 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 18 17:45:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.04	152	46532	20.00	ng	0.00
21) Naphthalene-d8	10.84	136	189070	20.00	ng	0.00
38) Acenaphthene-d10	14.66	164	127744	20.00	ng	0.00
63) Phenanthrene-d10	17.41	188	337921	20.00	ng	0.00
75) Chrysene-d12	21.67	240	379019	20.00	ng	0.00
86) Perylene-d12	24.88	264	367459	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.62	112	221089	78.88	ng	0.00
7) Phenol-d6	7.21	99	330333	79.00	ng	0.00
23) Nitrobenzene-d5	9.20	82	341888	75.54	ng	0.00
41) 2,4,6-Tribromophenol	16.15	330	139801	83.03	ng	0.00
44) 2-Fluorobiphenyl	13.29	172	765175	80.25	ng	0.00
78) Terphenyl-d14	20.01	244	1304746	75.88	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.54	88	51704	36.245	ng	# 73
3) Pyridine	3.93	79	142854	38.360	ng	# 76
4) n-Nitrosodimethylamine	3.85	42	56439	35.296	ng	# 77
6) Aniline	7.36	93	203624	37.569	ng	97
8) 2-Chlorophenol	7.61	128	114917	40.314	ng	96
9) Benzaldehyde	7.18	77	94311	36.757	ng	97
10) Phenol	7.23	94	168627	39.914	ng	95
11) bis(2-Chloroethyl)ether	7.46	93	130184	39.348	ng	89
12) 1,3-Dichlorobenzene	7.93	146	136215	39.571	ng	94
13) 1,4-Dichlorobenzene	8.07	146	138421	39.577	ng	93
14) 1,2-Dichlorobenzene	8.39	146	132400	39.405	ng	99
15) Benzyl Alcohol	8.28	79	143004	37.659	ng	95
16) 2,2'-oxybis(1-Chloropropan	8.56	45	124166m	44.764	ng	
17) 2-Methylphenol	8.48	107	110437	38.448	ng	94
18) Hexachloroethane	9.12	117	52343	39.141	ng	97
19) n-Nitroso-di-n-propylamine	8.84	70	124554	35.372	ng	# 83
20) 3+4-Methylphenols	8.82	107	159198	39.952	ng	95
22) Acetophenone	8.86	105	222058	37.768	ng	# 91
24) Nitrobenzene	9.24	77	173391	38.094	ng	97
25) Isophorone	9.76	82	306933	36.532	ng	99
26) 2-Nitrophenol	9.95	139	70995	43.918	ng	92
27) 2,4-Dimethylphenol	10.01	122	107259	39.198	ng	87
28) bis(2-Chloroethoxy)methane	10.24	93	177177	38.271	ng	99
29) 2,4-Dichlorophenol	10.50	162	127271	40.745	ng	93
30) 1,2,4-Trichlorobenzene	10.71	180	156286	40.803	ng	92
31) Naphthalene	10.89	128	377996	38.380	ng	97
32) Benzoic acid	10.15	122	36814	19.985	ng	# 81
33) 4-Chloroaniline	11.00	127	164924	38.421	ng	# 95
34) Hexachlorobutadiene	11.18	225	118982	39.837	ng	93
35) Caprolactam	11.78	113	47101	35.138	ng	# 76
36) 4-Chloro-3-methylphenol	12.12	107	156823	37.456	ng	98
37) 2-Methylnaphthalene	12.49	142	286527	39.050	ng	# 92
39) 1,2,4,5-Tetrachlorobenzene	12.86	216	194797	40.402	ng	98
40) Hexachlorocyclopentadiene	12.84	237	102558	40.559	ng	89

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.10	196	121453	43.074	nq	94
43) 2,4,5-Trichlorophenol	13.18	196	129787	41.146	nq	97
45) 1,1'-Biphenyl	13.50	154	401758	40.566	nq	97
46) 2-Chloronaphthalene	13.54	162	311872	40.483	nq	99
47) 2-Nitroaniline	13.74	65	108883	39.979	nq	97
48) Acenaphthylene	14.39	152	510586	39.566	nq	97
49) Dimethylphthalate	14.12	163	430568	38.470	nq	99
50) 2,6-Dinitrotoluene	14.24	165	93956	41.224	nq	92
51) Acenaphthene	14.72	154	314929	39.177	nq	98
52) 3-Nitroaniline	14.57	138	92521	39.718	nq #	96
54) Dibenzofuran	15.06	168	501606	39.235	nq	92
55) 4-Nitrophenol	14.89	139	58688	35.376	nq #	82
56) 2,4-Dinitrotoluene	15.02	165	139478	42.657	nq #	85
57) Fluorene	15.71	166	422975	39.474	nq	99
58) 2,3,4,6-Tetrachlorophenol	15.29	232	121569	40.197	nq #	92
59) Diethylphthalate	15.48	149	443308	38.520	nq	98
60) 4-Chlorophenyl-phenylether	15.70	204	248468	39.452	nq	97
61) 4-Nitroaniline	15.74	138	97233	39.903	nq #	83
62) Azobenzene	15.99	77	422630	37.230	nq	95
64) 4,6-Dinitro-2-methylphenol	15.79	198	29703	15.790	nq #	75
65) n-Nitrosodiphenylamine	15.92	169	396277	41.200	nq	94
66) 4-Bromophenyl-phenylether	16.59	248	159036	40.566	nq	95
67) Hexachlorobenzene	16.72	284	168254	41.675	nq	92
68) Atrazine	16.86	200	156245	39.454	nq	95
69) Pentachlorophenol	17.06	266	100336	40.802	nq	95
70) Phenanthrene	17.45	178	673491	39.942	nq	99
71) Anthracene	17.54	178	670485	39.934	nq	97
72) Carbazole	17.81	167	632983	39.698	nq	98
73) Di-n-butylphthalate	18.36	149	751120	39.863	nq #	98
74) Fluoranthene	19.45	202	890315	39.199	nq	98
76) Benzidine	19.64	184	336861	33.806	nq	95
77) Pyrene	19.82	202	902869	37.673	nq	98
79) Butylbenzylphthalate	20.70	149	344005	40.139	nq	96
80) Benzo(a)anthracene	21.65	228	904082	38.277	nq	100
81) 3,3'-Dichlorobenzidine	21.57	252	331473	40.249	nq	96
82) Chrysene	21.72	228	849746	38.823	nq	98
83) Bis(2-ethylhexyl)phthalate	21.55	149	478985	41.110	nq	100
84) Di-n-octyl phthalate	22.76	149	813691	41.709	nq #	93
85) Indeno(1,2,3-cd)pyrene	28.55	276	960973	39.656	nq #	94
87) Benzo(b)fluoranthene	23.87	252	853042	38.195	nq #	96
88) Benzo(k)fluoranthene	23.93	252	873255	39.994	nq #	97
89) Benzo(a)pyrene	24.73	252	812678	39.113	nq #	96
90) Dibenzo(a,h)anthracene	28.60	278	811786	39.796	nq	97
91) Benzo(g,h,i)perylene	29.69	276	792714	39.713	nq #	95

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

