

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG121417\
 Data File : BG031579.D
 Acq On : 15 Dec 2017 4:37
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 12/15/2017 6:47:06 PM

Quant Time: Dec 15 08:14:45 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG121117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 11 17:25:55 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.10	152	72545	20.00	ng	0.00
21) Naphthalene-d8	10.91	136	312626	20.00	ng	0.00
38) Acenaphthene-d10	14.72	164	202737	20.00	ng	0.00
63) Phenanthrene-d10	17.47	188	505053	20.00	ng	0.00
75) Chrysene-d12	21.74	240	554936	20.00	ng	0.00
86) Perylene-d12	25.02	264	515514	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.65	112	339269	82.89	ng	-0.01
7) Phenol-d6	7.26	99	446668	78.62	ng	0.00
23) Nitrobenzene-d5	9.27	82	417665	82.35	ng	0.00
41) 2,4,6-Tribromophenol	16.22	330	215663	90.80	ng	0.00
44) 2-Fluorobiphenyl	13.35	172	1152505	83.44	ng	0.00
78) Terphenyl-d14	20.06	244	1926689	78.99	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.50	88	75464	38.644	ng	# 81
3) Pyridine	3.90	79	205300	39.929	ng	86
4) n-Nitrosodimethylamine	3.82	42	89314	36.879	ng	# 86
6) Aniline	7.41	93	286021	38.225	ng	95
8) 2-Chlorophenol	7.66	128	184517	40.418	ng	88
9) Benzaldehyde	7.23	77	125151	38.001	ng	96
10) Phenol	7.29	94	228781	39.197	ng	92
11) bis(2-Chloroethyl)ether	7.51	93	186701	39.192	ng	90
12) 1,3-Dichlorobenzene	7.98	146	216520	39.882	ng	95
13) 1,4-Dichlorobenzene	8.13	146	224232	39.741	ng	94
14) 1,2-Dichlorobenzene	8.45	146	215100	40.020	ng	97
15) Benzyl Alcohol	8.34	79	159978	35.994	ng	91
16) 2,2'-oxybis(1-Chloropropan	8.62	45	225401m	42.117	ng	
17) 2-Methylphenol	8.55	107	154603	38.858	ng	97
18) Hexachloroethane	9.18	117	75416	39.652	ng	94
19) n-Nitroso-di-n-propylamine	8.90	70	153182	34.254	ng	# 94
20) 3+4-Methylphenols	8.88	107	215735	36.399	ng	96
22) Acetophenone	8.92	105	305516	40.736	ng	# 92
24) Nitrobenzene	9.31	77	212032	40.796	ng	# 91
25) Isophorone	9.83	82	377365	39.369	ng	# 90
26) 2-Nitrophenol	10.02	139	114140	44.412	ng	# 87
27) 2,4-Dimethylphenol	10.08	122	163340	41.838	ng	93
28) bis(2-Chloroethoxy)methane	10.31	93	255984	41.365	ng	95
29) 2,4-Dichlorophenol	10.57	162	200351	43.557	ng	89
30) 1,2,4-Trichlorobenzene	10.78	180	242356	41.248	ng	97
31) Naphthalene	10.96	128	601290	39.150	ng	98
32) Benzoic acid	10.25	122	64581	33.258	ng	90
33) 4-Chloroaniline	11.08	127	266022	39.892	ng	94
34) Hexachlorobutadiene	11.24	225	159741	40.987	ng	95
35) Caprolactam	11.85	113	71284	39.467	ng	# 65
36) 4-Chloro-3-methylphenol	12.20	107	205044	39.053	ng	# 83
37) 2-Methylnaphthalene	12.56	142	472501	39.734	ng	94
39) 1,2,4,5-Tetrachlorobenzene	12.93	216	341415	43.089	ng	# 96
40) Hexachlorocyclopentadiene	12.90	237	61330	34.193	ng	90

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.17	196	181101	44.736	nq	99
43) 2,4,5-Trichlorophenol	13.26	196	191660	43.951	nq	# 92
45) 1,1'-Biphenyl	13.56	154	689310	41.396	nq	97
46) 2-Chloronaphthalene	13.61	162	505759	41.498	nq	95
47) 2-Nitroaniline	13.81	65	136754	39.931	nq	# 81
48) Acenaphthylene	14.45	152	800328	40.811	nq	98
49) Dimethylphthalate	14.18	163	693861	41.339	nq	99
50) 2,6-Dinitrotoluene	14.30	165	152251	42.438	nq	# 77
51) Acenaphthene	14.79	154	507218	40.905	nq	99
52) 3-Nitroaniline	14.64	138	152333	41.653	nq	# 76
53) 2,4-Dinitrophenol	14.86	184	30227	26.957	nq	# 48
54) Dibenzofuran	15.12	168	791325	39.992	nq	98
55) 4-Nitrophenol	14.97	139	95104	40.324	nq	# 77
56) 2,4-Dinitrotoluene	15.10	165	213400	41.879	nq	# 72
57) Fluorene	15.77	166	650721	41.043	nq	97
58) 2,3,4,6-Tetrachlorophenol	15.36	232	174340	42.534	nq	# 90
59) Diethylphthalate	15.53	149	691104	41.078	nq	97
60) 4-Chlorophenyl-phenylether	15.76	204	389728	40.177	nq	92
61) 4-Nitroaniline	15.80	138	162547	42.354	nq	89
62) Azobenzene	16.05	77	523422	37.557	nq	91
64) 4,6-Dinitro-2-methylphenol	15.87	198	111817	37.746	nq	# 74
65) n-Nitrosodiphenylamine	15.98	169	625235	42.294	nq	99
66) 4-Bromophenyl-phenylether	16.65	248	252958	43.072	nq	96
67) Hexachlorobenzene	16.78	284	257484	42.458	nq	95
68) Atrazine	16.92	200	240248	44.446	nq	98
69) Pentachlorophenol	17.13	266	93795	35.560	nq	98
70) Phenanthrene	17.51	178	1066494	40.433	nq	99
71) Anthracene	17.60	178	1076648	41.271	nq	98
72) Carbazole	17.87	167	1009558	42.764	nq	99
73) Di-n-butylphthalate	18.41	149	1163954	42.694	nq	99
74) Fluoranthene	19.52	202	1374917	41.651	nq	97
76) Benzidine	19.69	184	602797	46.676	nq	98
77) Pyrene	19.88	202	1394287	40.436	nq	96
79) Butylbenzylphthalate	20.74	149	548917	45.378	nq	# 83
80) Benzo(a)anthracene	21.72	228	1362285	40.671	nq	98
81) 3,3'-Dichlorobenzidine	21.63	252	513984	44.090	nq	99
82) Chrysene	21.79	228	1282945	39.954	nq	97
83) Bis(2-ethylhexyl)phthalate	21.60	149	753770	43.312	nq	99
84) Di-n-octyl phthalate	22.81	149	1213503	42.272	nq	# 93
85) Indeno(1,2,3-cd)pyrene	28.80	276	1272468	37.080	nq	98
87) Benzo(b)fluoranthene	23.98	252	1284489m	41.677	nq	
88) Benzo(k)fluoranthene	24.05	252	1269047	40.348	nq	98
89) Benzo(a)pyrene	24.87	252	1190384	40.715	nq	99
90) Dibenzo(a,h)anthracene	28.84	278	1074065	38.417	nq	97
91) Benzo(g,h,i)perylene	29.97	276	1037789	37.756	nq	97

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

