

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061016\
 Data File : BG022304.D
 Acq On : 11 Jun 2016 2:27
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

umangi
 6/13/2016 7:32:03 PM

Quant Time: Jun 13 15:56:22 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 06 16:12:45 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.10	152	47271	20.00	ng	-0.02
21) Naphthalene-d8	10.92	136	230292	20.00	ng	-0.01
38) Acenaphthene-d10	14.74	164	122002	20.00	ng	-0.01
63) Phenanthrene-d10	17.48	188	247948	20.00	ng	-0.01
75) Chrysene-d12	21.76	240	227257	20.00	ng	-0.01
86) Perylene-d12	25.06	264	211490	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.65	112	219918	89.95	ng	-0.01
7) Phenol-d6	7.27	99	325727	84.74	ng	0.00
23) Nitrobenzene-d5	9.28	82	297334	90.01	ng	-0.01
41) 2,4,6-Tribromophenol	16.23	330	94735	68.72	ng	0.00
44) 2-Fluorobiphenyl	13.36	172	684099	85.45	ng	0.00
78) Terphenyl-d14	20.08	244	768078	80.24	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.50	88	47090	40.17	ng	# 75
3) Pyridine	3.91	79	135720	42.84	ng	94
4) n-Nitrosodimethylamine	3.83	42	36030	50.86	ng	# 86
6) Aniline	7.43	93	209355	42.87	ng	# 82
8) 2-Chlorophenol	7.67	128	137884	40.80	ng	83
9) Benzaldehyde	7.24	77	90978	43.00	ng	# 79
10) Phenol	7.29	94	170534	43.03	ng	86
11) bis(2-Chloroethyl)ether	7.52	93	135346	42.83	ng	# 73
12) 1,3-Dichlorobenzene	7.99	146	140873	38.89	ng	# 93
13) 1,4-Dichlorobenzene	8.14	146	146229	40.52	ng	95
14) 1,2-Dichlorobenzene	8.46	146	143120	39.34	ng	97
15) Benzyl Alcohol	8.35	79	105413	42.44	ng	90
16) 2,2'-oxybis(1-Chloropropan	8.64	45	150056	45.77	ng	84
17) 2-Methylphenol	8.56	107	118162	39.95	ng	# 87
18) Hexachloroethane	9.19	117	54757	41.55	ng	# 84
19) n-Nitroso-di-n-propylamine	8.92	70	105036	40.37	ng	# 72
20) 3+4-Methylphenols	8.89	107	165437	39.00	ng	97
22) Acetophenone	8.93	105	210286	42.37	ng	# 84
24) Nitrobenzene	9.32	77	149348	44.96	ng	# 87
25) Isophorone	9.84	82	300602	38.90	ng	95
26) 2-Nitrophenol	10.04	139	80462	44.67	ng	# 77
27) 2,4-Dimethylphenol	10.09	122	135037	41.42	ng	93
28) bis(2-Chloroethoxy)methane	10.32	93	182747	41.27	ng	92
29) 2,4-Dichlorophenol	10.58	162	124886	41.35	ng	98
30) 1,2,4-Trichlorobenzene	10.78	180	135350	40.11	ng	96
31) Naphthalene	10.98	128	451550	39.28	ng	# 94
32) Benzoic acid	10.26	122	56659	49.70	ng	91
33) 4-Chloroaniline	11.09	127	191107	39.98	ng	# 91
34) Hexachlorobutadiene	11.25	225	71427	39.42	ng	97
35) Caprolactam	11.88	113	48787	29.45	ng	# 55
36) 4-Chloro-3-methylphenol	12.21	107	138665	38.73	ng	87
37) 2-Methylnaphthalene	12.58	142	320952	37.82	ng	96
39) 1,2,4,5-Tetrachlorobenzene	12.94	216	135733	43.22	ng	98
40) Hexachlorocyclopentadiene	12.91	237	56953	38.43	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.18	196	92350	43.83	ng	96
43) 2,4,5-Trichlorophenol	13.26	196	95134	40.87	ng	96
45) 1,1'-Biphenyl	13.57	154	401230	43.70	ng	94
46) 2-Chloronaphthalene	13.62	162	301099	42.78	ng	98
47) 2-Nitroaniline	13.83	65	76496	45.58	ng #	70
48) Acenaphthylene	14.46	152	498178	40.34	ng	98
49) Dimethylphthalate	14.19	163	367962	36.38	ng	99
50) 2,6-Dinitrotoluene	14.32	165	84775	38.55	ng #	85
51) Acenaphthene	14.80	154	296888	40.13	ng	96
52) 3-Nitroaniline	14.66	138	88788	36.40	ng #	85
53) 2,4-Dinitrophenol	14.87	184	33256	42.90	ng #	80
54) Dibenzofuran	15.14	168	448438	38.84	ng	100
55) 4-Nitrophenol	14.97	139	60230	35.78	ng #	79
56) 2,4-Dinitrotoluene	15.11	165	111753	37.89	ng #	88
57) Fluorene	15.78	166	374676	37.67	ng	97
58) 2,3,4,6-Tetrachlorophenol	15.37	232	76823m	36.80	ng	
59) Diethylphthalate	15.54	149	375798	34.78	ng	100
60) 4-Chlorophenyl-phenylether	15.77	204	171305	38.03	ng	94
61) 4-Nitroaniline	15.81	138	90425	34.96	ng #	78
62) Azobenzene	16.06	77	342081	41.80	ng	90
64) 4,6-Dinitro-2-methylphenol	15.87	198	52242	42.29	ng	86
65) n-Nitrosodiphenylamine	15.99	169	308902	43.25	ng	98
66) 4-Bromophenyl-phenylether	16.67	248	96803	41.67	ng	96
67) Hexachlorobenzene	16.79	284	104814	38.71	ng	98
68) Atrazine	16.93	200	96712	36.58	ng	98
69) Pentachlorophenol	17.14	266	43655	39.33	ng	97
70) Phenanthrene	17.53	178	533376	39.61	ng	99
71) Anthracene	17.62	178	532731	39.89	ng	97
72) Carbazole	17.89	167	490451	38.77	ng	97
73) Di-n-butylphthalate	18.42	149	627042	37.92	ng	98
74) Fluoranthene	19.53	202	573395	37.04	ng	95
76) Benzidine	19.70	184	281064	47.29	ng	100
77) Pyrene	19.89	202	569283	40.29	ng	99
79) Butylbenzylphthalate	20.75	149	295847	45.15	ng	96
80) Benzo(a)anthracene	21.74	228	521840	40.95	ng	100
81) 3,3'-Dichlorobenzidine	21.65	252	152281	42.57	ng #	98
82) Chrysene	21.81	228	501621	40.45	ng	99
83) Bis(2-ethylhexyl)phthalate	21.61	149	423650	43.97	ng	97
84) Di-n-octyl phthalate	22.84	149	714255	44.65	ng #	88
85) Indeno(1,2,3-cd)pyrene	28.83	276	549513	39.50	ng #	85
87) Benzo(b)fluoranthene	24.00	252	508454	41.22	ng #	96
88) Benzo(k)fluoranthene	24.07	252	488863	40.26	ng #	96
89) Benzo(a)pyrene	24.90	252	491488	41.67	ng #	96
90) Dibenzo(a,h)anthracene	28.88	278	451325	40.63	ng #	94
91) Benzo(g,h,i)perylene	30.00	276	456564	39.51	ng #	92

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

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