

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060617\
 Data File : BG027253.D
 Acq On : 6 Jun 2017 17:32
 Operator : SJ/MA
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 6/8/2017 8:24:50 AM

Quant Time: Jun 06 19:06:38 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 05 16:00:35 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.57	152	66170	20.00	ng	-0.01
21) Naphthalene-d8	11.39	136	317523	20.00	ng	-0.01
38) Acenaphthene-d10	15.15	164	193437	20.00	ng	-0.01
63) Phenanthrene-d10	17.90	188	405448	20.00	ng	-0.02
75) Chrysene-d12	22.29	240	434221	20.00	ng	-0.02
86) Perylene-d12	25.98	264	403250	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	6.15	112	365328	84.25	ng	-0.01
7) Phenol-d6	7.70	99	548442	86.16	ng	0.00
23) Nitrobenzene-d5	9.73	82	478636	94.80	ng	-0.01
41) 2,4,6-Tribromophenol	16.63	330	218690	85.85	ng	-0.01
44) 2-Fluorobiphenyl	13.78	172	1140421	82.48	ng	-0.01
78) Terphenyl-d14	20.48	244	1359135	79.85	ng	-0.02
Target Compounds						
2) 1,4-Dioxane	4.16	88	70554	39.240	ng	Qvalue # 79
3) Pyridine	4.54	79	229382	41.385	ng	# 73
4) n-Nitrosodimethylamine	4.45	42	87712	42.740	ng	# 55
6) Aniline	7.89	93	336337	41.960	ng	# 92
8) 2-Chlorophenol	8.13	128	190720	41.653	ng	90
9) Benzaldehyde	7.71	77	186090	42.284	ng	85
10) Phenol	7.73	94	272944	41.931	ng	76
11) bis(2-Chloroethyl)ether	7.97	93	220294	41.253	ng	85
12) 1,3-Dichlorobenzene	8.46	146	205684	39.827	ng	# 94
13) 1,4-Dichlorobenzene	8.60	146	212622	40.116	ng	96
14) 1,2-Dichlorobenzene	8.93	146	204497	39.558	ng	96
15) Benzyl Alcohol	8.79	79	196762	43.873	ng	# 85
16) 2,2'-oxybis(1-Chloropropan	9.08	45	320161	41.901	ng	91
17) 2-Methylphenol	8.98	107	191929	41.712	ng	# 87
18) Hexachloroethane	9.66	117	75588	42.337	ng	86
19) n-Nitroso-di-n-propylamine	9.36	70	188065	42.234	ng	# 85
20) 3+4-Methylphenols	9.31	107	271568	42.607	ng	89
22) Acetophenone	9.38	105	338082	41.635	ng	# 83
24) Nitrobenzene	9.77	77	259744	45.157	ng	# 85
25) Isophorone	10.29	82	499850	42.103	ng	# 96
26) 2-Nitrophenol	10.48	139	109519	46.456	ng	# 76
27) 2,4-Dimethylphenol	10.52	122	216918	42.495	ng	93
28) bis(2-Chloroethoxy)methane	10.76	93	316093	41.119	ng	98
29) 2,4-Dichlorophenol	11.02	162	202052	43.359	ng	99
30) 1,2,4-Trichlorobenzene	11.24	180	212808	41.242	ng	99
31) Naphthalene	11.44	128	696740	40.150	ng	99
32) Benzoic acid	10.64	122	154527	44.474	ng	# 88
33) 4-Chloroaniline	11.53	127	297260	41.107	ng	# 87
34) Hexachlorobutadiene	11.71	225	128378	40.482	ng	98
35) Caprolactam	12.32	113	72138	45.137	ng	# 75
36) 4-Chloro-3-methylphenol	12.60	107	245213	41.984	ng	91
37) 2-Methylnaphthalene	13.00	142	504954m	39.890	ng	
39) 1,2,4,5-Tetrachlorobenzene	13.36	216	252172	41.847	ng	98
40) Hexachlorocyclopentadiene	13.34	237	131876	41.117	ng	94

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.58	196	165108	45.173	ng	96
43) 2,4,5-Trichlorophenol	13.66	196	183703	44.915	ng	94
45) 1,1'-Biphenyl	13.98	154	654957	41.482	ng	97
46) 2-Chloronaphthalene	14.04	162	492461	41.601	ng	100
47) 2-Nitroaniline	14.23	65	158264	47.153	ng	# 80
48) Acenaphthylene	14.88	152	826194	41.239	ng	98
49) Dimethylphthalate	14.59	163	613909	40.008	ng	97
50) 2,6-Dinitrotoluene	14.71	165	135903	44.111	ng	# 78
51) Acenaphthene	15.21	154	541880	41.583	ng	98
52) 3-Nitroaniline	15.04	138	137116	42.653	ng	87
53) 2,4-Dinitrophenol	15.24	184	62139	50.331	ng	# 42
54) Dibenzofuran	15.54	168	732348	40.216	ng	91
55) 4-Nitrophenol	15.32	139	106678	40.111	ng	# 56
56) 2,4-Dinitrotoluene	15.49	165	175056	44.491	ng	# 87
57) Fluorene	16.19	166	645461	39.906	ng	99
58) 2,3,4,6-Tetrachlorophenol	15.76	232	155610	42.620	ng	100
59) Diethylphthalate	15.94	149	608361	39.747	ng	98
60) 4-Chlorophenyl-phenylether	16.18	204	324452	39.822	ng	93
61) 4-Nitroaniline	16.21	138	133779	41.361	ng	# 66
62) Azobenzene	16.47	77	586457	41.341	ng	89
64) 4,6-Dinitro-2-methylphenol	16.25	198	94125	51.065	ng	88
65) n-Nitrosodiphenylamine	16.39	169	539932	42.477	ng	99
66) 4-Bromophenyl-phenylether	17.07	248	203350	42.376	ng	100
67) Hexachlorobenzene	17.20	284	215643	41.779	ng	90
68) Atrazine	17.32	200	181543	42.709	ng	98
69) Pentachlorophenol	17.53	266	126043	41.660	ng	97
70) Phenanthrene	17.94	178	925053	40.552	ng	96
71) Anthracene	18.03	178	937172	41.008	ng	99
72) Carbazole	18.30	167	854690	39.652	ng	96
73) Di-n-butylphthalate	18.83	149	896110	42.874	ng	# 93
74) Fluoranthene	19.93	202	983280	40.349	ng	94
76) Benzidine	20.09	184	468135	35.789	ng	97
77) Pyrene	20.29	202	1003569	38.252	ng	96
79) Butylbenzylphthalate	21.18	149	398807	43.020	ng	91
80) Benzo(a)anthracene	22.27	228	1020130	41.196	ng	95
81) 3,3'-Dichlorobenzidine	22.16	252	401739	41.702	ng	98
82) Chrysene	22.34	228	966789	40.665	ng	96
83) Bis(2-ethylhexyl)phthalate	22.13	149	581940	41.539	ng	99
84) Di-n-octyl phthalate	23.52	149	983343	42.380	ng	# 88
85) Indeno(1,2,3-cd)pyrene	30.25	276	1188292	41.271	ng	# 93
87) Benzo(b)fluoranthene	24.80	252	1051945	42.062	ng	# 97
88) Benzo(k)fluoranthene	24.88	252	1017308	41.463	ng	# 94
89) Benzo(a)pyrene	25.81	252	979180	41.591	ng	# 94
90) Dibenzo(a,h)anthracene	30.34	278	1044109	42.366	ng	# 94
91) Benzo(g,h,i)perylene	31.59	276	990719	41.509	ng	# 91

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

