

Data Path : Z:\HPCHEM1\BNA G\DATA\BG080116\
 Data File : BG023386.D
 Acq On : 1 Aug 2016 16:36
 Operator : UM/SJ
 Sample : SSTDICC050
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC050

Manual Integrations
 APPROVED

sohil
 8/2/2016 5:10:36 PM

Quant Time: Aug 01 17:48:47 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG080116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 01 17:40:36 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.13	152	164042	20.00	ng	0.00
21) Naphthalene-d8	10.96	136	772641	20.00	ng	0.00
38) Acenaphthene-d10	14.79	164	541453	20.00	ng	0.00
63) Phenanthrene-d10	17.55	188	1269888	20.00	ng	0.00
75) Chrysene-d12	21.87	240	1057173	20.00	ng	0.00
86) Perylene-d12	25.25	264	1095242	20.00	ng	0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.67	112	976945	101.67	ng	0.00
7) Phenol-d6	7.28	99	1506866	102.14	ng	0.00
23) Nitrobenzene-d5	9.31	82	1569077	99.31	ng	0.01
41) 2,4,6-Tribromophenol	16.29	330	786243	140.30	ng	0.00
44) 2-Fluorobiphenyl	13.41	172	2974057	102.80	ng	0.01
78) Terphenyl-d14	20.16	244	3267146	100.46	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.53	88	200860	48.74	ng	# 91
3) Pyridine	3.93	79	706164	48.96	ng	# 92
4) n-Nitrosodimethylamine	3.85	42	257329	39.49	ng	# 64
6) Aniline	7.45	93	1109494	56.31	ng	96
8) 2-Chlorophenol	7.69	128	556587	51.91	ng	98
9) Benzaldehyde	7.26	77	406945	38.40	ng	93
10) Phenol	7.31	94	820751	52.73	ng	# 77
11) bis(2-Chloroethyl)ether	7.54	93	641314	48.47	ng	92
12) 1,3-Dichlorobenzene	8.02	146	630864	52.22	ng	# 92
13) 1,4-Dichlorobenzene	8.16	146	653875	53.05	ng	95
14) 1,2-Dichlorobenzene	8.49	146	628773	52.88	ng	94
15) Benzyl Alcohol	8.37	79	599582	60.39	ng	92
16) 2,2'-oxybis(1-Chloropropan	8.66	45	946633	44.96	ng	90
17) 2-Methylphenol	8.57	107	543810	51.11	ng	# 86
18) Hexachloroethane	9.22	117	247925	48.28	ng	95
19) n-Nitroso-di-n-propylamine	8.94	70	657894	49.58	ng	98
20) 3+4-Methylphenols	8.90	107	779005	53.85	ng	89
22) Acetophenone	8.96	105	1053473	52.22	ng	# 89
24) Nitrobenzene	9.35	77	862501	47.01	ng	# 89
25) Isophorone	9.87	82	1650427	50.57	ng	# 93
26) 2-Nitrophenol	10.06	139	389340	56.18	ng	# 75
27) 2,4-Dimethylphenol	10.12	122	640617	54.84	ng	84
28) bis(2-Chloroethoxy)methane	10.35	93	974237	53.21	ng	97
29) 2,4-Dichlorophenol	10.61	162	652560	57.70	ng	99
30) 1,2,4-Trichlorobenzene	10.82	180	672208	53.79	ng	94
31) Naphthalene	11.01	128	1927078	53.68	ng	97
32) Benzoic acid	10.30	122	590582	71.26	ng	# 86
33) 4-Chloroaniline	11.12	127	960865	57.93	ng	94
34) Hexachlorobutadiene	11.29	225	446029	54.46	ng	95
35) Caprolactam	11.91	113	273783	56.93	ng	93
36) 4-Chloro-3-methylphenol	12.25	107	838301	56.42	ng	92
37) 2-Methylnaphthalene	12.62	142	1475507m	55.43	ng	
39) 1,2,4,5-Tetrachlorobenzene	12.98	216	975025	55.62	ng	99
40) Hexachlorocyclopentadiene	12.96	237	507127	61.05	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.22	196	598763	58.15	nq	97
43) 2,4,5-Trichlorophenol	13.30	196	610910	55.93	nq	# 95
45) 1,1'-Biphenyl	13.62	154	2002398	53.03	nq	91
46) 2-Chloronaphthalene	13.67	162	1565652	51.74	nq	93
47) 2-Nitroaniline	13.87	65	686622	53.64	nq	# 86
48) Acenaphthylene	14.51	152	2415694	52.14	nq	# 93
49) Dimethylphthalate	14.24	163	1975727	49.45	nq	# 95
50) 2,6-Dinitrotoluene	14.37	165	490837	52.46	nq	# 78
51) Acenaphthene	14.85	154	1567242	52.15	nq	97
52) 3-Nitroaniline	14.70	138	555116	58.90	nq	# 83
53) 2,4-Dinitrophenol	14.91	184	336298	61.38	nq	# 84
54) Dibenzofuran	15.19	168	2194405	51.61	nq	98
55) 4-Nitrophenol	15.01	139	438835	53.68	nq	# 75
56) 2,4-Dinitrotoluene	15.15	165	702242	53.30	nq	88
57) Fluorene	15.84	166	1914444	50.99	nq	98
58) 2,3,4,6-Tetrachlorophenol	15.42	232	538822	58.27	nq	93
59) Diethylphthalate	15.60	149	1985787	48.52	nq	# 92
60) 4-Chlorophenyl-phenylether	15.83	204	1028899	50.62	nq	98
61) 4-Nitroaniline	15.87	138	582102	58.91	nq	# 64
62) Azobenzene	16.12	77	2042627	48.14	nq	89
64) 4,6-Dinitro-2-methylphenol	15.92	198	467622	57.33	nq	95
65) n-Nitrosodiphenylamine	16.05	169	1819261	53.66	nq	# 91
66) 4-Bromophenyl-phenylether	16.73	248	753152	58.66	nq	95
67) Hexachlorobenzene	16.85	284	811025	62.82	nq	99
68) Atrazine	17.00	200	694813	53.43	nq	95
69) Pentachlorophenol	17.20	266	483131	60.63	nq	98
70) Phenanthrene	17.60	178	2881069	51.87	nq	# 89
71) Anthracene	17.69	178	2908365	51.82	nq	# 89
72) Carbazole	17.96	167	2629206	51.15	nq	93
73) Di-n-butylphthalate	18.50	149	2882187	50.25	nq	# 96
74) Fluoranthene	19.61	202	2986766	47.12	nq	87
76) Benzidine	19.78	184	1439128	48.40	nq	96
77) Pyrene	19.97	202	2984274	49.74	nq	# 91
79) Butylbenzylphthalate	20.85	149	1268272	48.05	nq	95
80) Benzo(a)anthracene	21.84	228	2729518	49.32	nq	95
81) 3,3'-Dichlorobenzidine	21.76	252	1215405	54.16	nq	# 94
82) Chrysene	21.91	228	2629247	49.17	nq	93
83) Bis(2-ethylhexyl)phthalate	21.73	149	1739398	47.10	nq	96
84) Di-n-octyl phthalate	22.99	149	2949817	48.83	nq	99
85) Indeno(1,2,3-cd)pyrene	29.13	276	3485545	55.88	nq	# 100
87) Benzo(b)fluoranthene	24.17	252	3019092	51.89	nq	# 94
88) Benzo(k)fluoranthene	24.24	252	2887959	49.65	nq	# 94
89) Benzo(a)pyrene	25.09	252	2883285	50.57	nq	# 94
90) Dibenzo(a,h)anthracene	29.19	278	2987204	55.39	nq	# 89
91) Benzo(g,h,i)perylene	30.33	276	3015982	53.12	nq	# 90

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

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