

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG010515\
 Data File : BG015838.D
 Acq On : 5 Jan 2015 12:05
 Operator : TP/IZ
 Sample : SSTDICC2.5
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDICC2.5

Manual Integrations
 APPROVED

TEJASKUMAR
 1/6/2015 6:24:45 PM

Quant Time: Jan 06 05:09:51 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG010515.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 06 03:51:36 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	24976	20.00	ng	-0.01
21) Naphthalene-d8	10.51	136	88062	20.00	ng	0.00
38) Acenaphthene-d10	14.37	164	64898	20.00	ng	0.00
63) Phenanthrene-d10	17.11	188	168268	20.00	ng	0.00
75) Chrysene-d12	21.29	240	255814	20.00	ng	0.00
86) Perylene-d12	23.56	264	246952	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.31	112	6538	2.25	ng	-0.02
7) Phenol-d6	6.90	99	9740	2.41	ng	-0.01
23) Nitrobenzene-d5	8.88	82	9833	2.19	ng	0.00
41) 2,4,6-Tribromophenol	15.86	330	5174	2.10	ng	0.00
44) 2-Fluorobiphenyl	12.98	172	22492	2.48	ng	0.00
78) Terphenyl-d14	19.74	244	56712	2.48	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.25	88	1942	2.92	ng	# 78
3) Pyridine	3.64	79	4233	2.32	ng	# 73
4) n-Nitrosodimethylamine	3.57	42	1470	1.87	ng	# 47
6) Aniline	7.05	93	5911	2.16	ng	# 98
8) 2-Chlorophenol	7.28	128	3766	2.52	ng	# 74
9) Benzaldehyde	6.87	77	3536	2.60	ng	# 94
10) Phenol	6.92	94	5675	2.58	ng	# 70
11) bis(2-Chloroethyl)ether	7.15	93	4085	2.52	ng	# 84
12) 1,3-Dichlorobenzene	7.60	146	4218m	2.35	ng	# 81
13) 1,4-Dichlorobenzene	7.76	146	4223	2.31	ng	# 89
14) 1,2-Dichlorobenzene	8.07	146	4427	2.55	ng	# 90
15) Benzyl Alcohol	7.96	79	4062	2.06	ng	# 72
16) 2,2'-oxybis(1-Chloropropan	8.25	45	2402	2.32	ng	# 87
17) 2-Methylphenol	8.17	107	3594	2.39	ng	# 60
18) Hexachloroethane	8.79	117	1949	2.31	ng	# 74
19) n-Nitroso-di-n-propylamine	8.52	70	3702	2.35	ng	# 84
20) 3+4-Methylphenols	8.49	107	4365	2.18	ng	# 87
22) Acetophenone	8.54	105	5977	2.32	ng	# 77
24) Nitrobenzene	8.91	77	5194	2.30	ng	# 96
25) Isophorone	9.44	82	8667	2.18	ng	# 82
26) 2-Nitrophenol	9.62	139	2012	2.32	ng	# 97
27) 2,4-Dimethylphenol	9.69	122	3305	2.28	ng	# 86
28) bis(2-Chloroethoxy)methane	9.93	93	4450	2.27	ng	# 85
29) 2,4-Dichlorophenol	10.16	162	3071	2.10	ng	# 89
30) 1,2,4-Trichlorobenzene	10.38	180	4263	2.22	ng	# 90
31) Naphthalene	10.56	128	11145	2.39	ng	# 92
33) 4-Chloroaniline	10.68	127	4599	2.42	ng	# 81
34) Hexachlorobutadiene	10.85	225	4494	2.52	ng	# 32
35) Caprolactam	11.43	113	1351	2.10	ng	# 70
36) 4-Chloro-3-methylphenol	11.80	107	4001	2.10	ng	# 99
37) 2-Methylnaphthalene	12.17	142	8095	2.32	ng	# 95
39) 1,2,4,5-Tetrachlorobenzene	12.55	216	6301	2.61	ng	# 91
40) Hexachlorocyclopentadiene	12.53	237	2246	1.40	ng	# 91
42) 2,4,6-Trichlorophenol	12.79	196	2890	1.95	ng	# 91

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,5-Trichlorophenol	12.87	196	3248	2.00	ng	# 89
45) 1,1'-Biphenyl	13.20	154	11768	2.54	ng	86
46) 2-Chloronaphthalene	13.24	162	9457	2.56	ng	# 90
47) 2-Nitroaniline	13.44	65	2551	2.11	ng	# 76
48) Acenaphthylene	14.08	152	14949	2.37	ng	# 93
49) Dimethylphthalate	13.82	163	11865	2.48	ng	97
50) 2,6-Dinitrotoluene	13.94	165	1977	1.91	ng	# 77
51) Acenaphthene	14.43	154	8571	2.31	ng	# 80
52) 3-Nitroaniline	14.28	138	2333	2.38	ng	# 76
54) Dibenzofuran	14.77	168	14877	2.47	ng	97
55) 4-Nitrophenol	14.59	139	1671	2.07	ng	# 33
56) 2,4-Dinitrotoluene	14.73	165	2783	1.87	ng	# 90
57) Fluorene	15.41	166	11362	2.32	ng	# 90
58) 2,3,4,6-Tetrachlorophenol	14.99	232	3913	2.14	ng	# 97
59) Diethylphthalate	15.19	149	13010	2.46	ng	# 86
60) 4-Chlorophenyl-phenylether	15.41	204	7118	2.22	ng	94
61) 4-Nitroaniline	15.43	138	2631	2.28	ng	# 60
62) Azobenzene	15.70	77	12892	2.26	ng	96
64) 4,6-Dinitro-2-methylphenol	15.51	198	1077	0.93	ng	# 60
65) n-Nitrosodiphenylamine	15.63	169	10953	2.26	ng	92
66) 4-Bromophenyl-phenylether	16.30	248	4805	2.08	ng	95
67) Hexachlorobenzene	16.42	284	6525	2.50	ng	# 88
68) Atrazine	16.57	200	4997	2.31	ng	82
69) Pentachlorophenol	16.77	266	1709	1.11	ng	# 50
70) Phenanthrene	17.16	178	20883	2.33	ng	# 90
71) Anthracene	17.24	178	19844	2.23	ng	93
72) Carbazole	17.51	167	18996	2.23	ng	94
73) Di-n-butylphthalate	18.08	149	21629	2.15	ng	# 92
74) Fluoranthene	19.17	202	31211	2.30	ng	95
76) Benzidine	19.35	184	13906	2.35	ng	98
77) Pyrene	19.54	202	32688	2.39	ng	95
79) Butylbenzylphthalate	20.43	149	12491	2.44	ng	99
80) Benzo(a)anthracene	21.28	228	35328	2.33	ng	98
81) 3,3'-Dichlorobenzidine	21.22	252	12786	2.29	ng	# 93
82) Chrysene	21.33	228	34500	2.47	ng	93
83) Bis(2-ethylhexyl)phthalate	21.21	149	17766	2.38	ng	94
84) Di-n-octyl phthalate	22.09	149	27433	2.26	ng	# 96
85) Indeno(1,2,3-cd)pyrene	25.87	276	37614	2.28	ng	# 62
87) Benzo(b)fluoranthene	22.87	252	35523m	2.30	ng	
88) Benzo(k)fluoranthene	22.92	252	33965m	2.29	ng	
89) Benzo(a)pyrene	23.45	252	31529	2.27	ng	# 96
90) Dibenzo(a,h)anthracene	25.88	278	30978	2.27	ng	93
91) Benzo(g,h,i)perylene	26.57	276	31709	2.34	ng	# 94

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

