

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG012015\
 Data File : BG015887.D
 Acq On : 20 Jan 2015 12:51
 Operator : TP/IZ
 Sample : SSTDICC2.5
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC2.5

Quant Time: Jan 20 15:27:14 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG012015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 20 15:14:03 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	19907	20.00	ng	0.00
21) Naphthalene-d8	10.47	136	74009	20.00	ng	0.00
38) Acenaphthene-d10	14.33	164	61839	20.00	ng	0.00
63) Phenanthrene-d10	17.08	188	186239	20.00	ng	0.00
75) Chrysene-d12	21.28	240	285079	20.00	ng	0.00
86) Perylene-d12	23.54	264	276321	20.00	ng	0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.28	112	27512	12.33	ng	0.00
7) Phenol-d6	6.86	99	36262	10.23	ng	0.00
23) Nitrobenzene-d5	8.84	82	50215	10.49	ng	0.00
41) 2,4,6-Tribromophenol	15.83	330	20450	10.78	ng	0.00
44) 2-Fluorobiphenyl	12.95	172	90826	13.43	ng	0.00
78) Terphenyl-d14	19.72	244	261023	13.80	ng	0.00

Target Compounds

						Qvalue
6) Aniline	7.02	93	26243	9.01	ng	# 87
8) 2-Chlorophenol	7.26	128	13683	11.09	ng	# 69
9) Benzaldehyde	6.84	77	16904	10.15	ng	# 80
10) Phenol	6.89	94	24220	10.74	ng	# 91
11) bis(2-Chloroethyl)ether	7.11	93	20434	12.24	ng	# 76
12) 1,3-Dichlorobenzene	7.58	146	14106	9.56	ng	# 77
13) 1,4-Dichlorobenzene	7.72	146	15655m	10.21	ng	# 79
14) 1,2-Dichlorobenzene	8.03	146	12048	8.34	ng	# 90
15) Benzyl Alcohol	7.93	79	21176	8.62	ng	# 97
17) 2-Methylphenol	8.13	107	14148	9.76	ng	# 85
19) n-Nitroso-di-n-propylamine	8.49	70	14895	7.00	ng	# 87
20) 3+4-Methylphenols	8.46	107	15348	7.60	ng	# 74
22) Acetophenone	8.50	105	25552	10.51	ng	# 88
24) Nitrobenzene	8.89	77	27298	9.33	ng	# 96
25) Isophorone	9.40	82	43028	8.61	ng	# 67
26) 2-Nitrophenol	9.60	139	5962	9.25	ng	# 55
27) 2,4-Dimethylphenol	9.66	122	13044	11.38	ng	# 78
28) bis(2-Chloroethoxy)methane	9.89	93	22315	10.85	ng	# 91
29) 2,4-Dichlorophenol	10.13	162	13928	10.64	ng	# 69
30) 1,2,4-Trichlorobenzene	10.34	180	19309	12.44	ng	# 95
31) Naphthalene	10.52	128	41336	11.66	ng	# 88
33) 4-Chloroaniline	10.64	127	18860	11.54	ng	# 74
36) 4-Chloro-3-methylphenol	11.78	107	18201	9.73	ng	# 78
37) 2-Methylnaphthalene	12.14	142	32406	11.46	ng	# 87
39) 1,2,4,5-Tetrachlorobenzene	12.52	216	28706	13.44	ng	# 87
42) 2,4,6-Trichlorophenol	12.76	196	15046	11.18	ng	# 96
45) 1,1'-Biphenyl	13.16	154	48887	13.03	ng	# 75
46) 2-Chloronaphthalene	13.20	162	39706	12.68	ng	# 90
47) 2-Nitroaniline	13.42	65	16454	10.02	ng	# 92
48) Acenaphthylene	14.06	152	52959	11.17	ng	# 98
49) Dimethylphthalate	13.79	163	51123	11.32	ng	# 74
50) 2,6-Dinitrotoluene	13.92	165	11670	11.71	ng	# 87
51) Acenaphthene	14.40	154	33549	11.36	ng	# 81
52) 3-Nitroaniline	14.24	138	10208	11.79	ng	#

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
54) Dibenzofuran	14.73	168	57960	12.16	ng	# 88
56) 2,4-Dinitrotoluene	14.71	165	15065	10.28	ng	# 79
57) Fluorene	15.38	166	52524	11.97	ng	89
58) 2,3,4,6-Tetrachlorophenol	14.97	232	13962	8.80	ng	# 94
59) Diethylphthalate	15.16	149	51139	10.87	ng	# 88
60) 4-Chlorophenyl-phenylether	15.38	204	32857	10.96	ng	92
61) 4-Nitroaniline	15.41	138	8173	8.39	ng	# 79
62) Azobenzene	15.67	77	92859	12.78	ng	87
65) n-Nitrosodiphenylamine	15.60	169	48469	9.90	ng	# 92
66) 4-Bromophenyl-phenylether	16.28	248	25110	9.69	ng	# 62
67) Hexachlorobenzene	16.39	284	29624	10.54	ng	# 89
68) Atrazine	16.55	200	21072	11.99	ng	100
70) Phenanthrene	17.13	178	95264	11.54	ng	# 85
71) Anthracene	17.22	178	94391	11.21	ng	# 94
72) Carbazole	17.49	167	69265	9.40	ng	# 91
73) Di-n-butylphthalate	18.06	149	81184	8.90	ng	# 95
74) Fluoranthene	19.14	202	130141	11.30	ng	86
76) Benzidine	19.34	184	37242	8.22	ng	# 93
77) Pyrene	19.51	202	141594	11.45	ng	# 84
79) Butylbenzylphthalate	20.41	149	39116	8.01	ng	89
80) Benzo(a)anthracene	21.26	228	155269	11.38	ng	# 86
81) 3,3'-Dichlorobenzidine	21.20	252	54324	9.35	ng	95
82) Chrysene	21.31	228	159322	12.57	ng	# 86
83) Bis(2-ethylhexyl)phthalate	21.19	149	58417	8.87	ng	97
84) Di-n-octyl phthalate	22.08	149	90877	9.15	ng	# 100
85) Indeno(1,2,3-cd)pyrene	25.85	276	139034	8.31	ng	# 94
87) Benzo(b)fluoranthene	22.86	252	149446	10.06	ng	100
88) Benzo(k)fluoranthene	22.89	252	157067m	11.47	ng	
89) Benzo(a)pyrene	23.44	252	150541	10.82	ng	# 90
90) Dibenzo(a,h)anthracene	25.87	278	114194	7.98	ng	# 95
91) Benzo(g,h,i)perylene	26.56	276	127887	9.10	ng	96

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

