

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG051415\
 Data File : BG017170.D
 Acq On : 15 May 2015 10:08
 Operator : TP/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

apatel
 5/15/2015 11:31:43 AM

Quant Time: May 15 11:19:00 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG051315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 13 16:12:13 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	43269	20.00	ng	0.00
21) Naphthalene-d8	10.50	136	204117	20.00	ng	0.00
38) Acenaphthene-d10	14.36	164	139799	20.00	ng	0.00
63) Phenanthrene-d10	17.11	188	311054	20.00	ng	0.00
75) Chrysene-d12	21.30	240	318491	20.00	ng	0.00
86) Perylene-d12	23.56	264	307004	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.30	112	132594	54.37	ng	0.00
7) Phenol-d6	6.91	99	195808	57.21	ng	0.00
23) Nitrobenzene-d5	8.87	82	228886	52.35	ng	0.00
41) 2,4,6-Tribromophenol	15.86	330	80829	47.83	ng	0.00
44) 2-Fluorobiphenyl	12.98	172	458502	48.13	ng	0.00
78) Terphenyl-d14	19.74	244	767465	50.22	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.18	88	24841	25.82	ng	97
3) Pyridine	3.57	79	78888	27.10	ng	93
4) n-Nitrosodimethylamine	3.49	42	53811	24.35	ng	# 95
6) Aniline	7.04	93	125327	26.47	ng	97
8) 2-Chlorophenol	7.28	128	75921	26.02	ng	92
9) Benzaldehyde	6.85	77	64704	26.52	ng	95
10) Phenol	6.94	94	103332	28.47	ng	96
11) bis(2-Chloroethyl)ether	7.14	93	75226	27.64	ng	99
12) 1,3-Dichlorobenzene	7.59	146	78201	24.00	ng	96
13) 1,4-Dichlorobenzene	7.74	146	82877	25.20	ng	94
14) 1,2-Dichlorobenzene	8.06	146	77050	24.53	ng	96
15) Benzyl Alcohol	7.95	79	83887	25.23	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.25	45	124202	28.15	ng	97
17) 2-Methylphenol	8.17	107	70286	26.71	ng	97
18) Hexachloroethane	8.78	117	35019	25.49	ng	# 87
19) n-Nitroso-di-n-propylamine	8.51	70	79209	26.56	ng	# 95
20) 3+4-Methylphenols	8.50	107	99375	27.17	ng	94
22) Acetophenone	8.53	105	120729	24.37	ng	97
24) Nitrobenzene	8.91	77	109851	25.68	ng	98
25) Isophorone	9.43	82	204615	25.39	ng	98
26) 2-Nitrophenol	9.62	139	48678	25.46	ng	89
27) 2,4-Dimethylphenol	9.69	122	80002	25.42	ng	94
28) bis(2-Chloroethoxy)methane	9.92	93	101459	25.86	ng	95
29) 2,4-Dichlorophenol	10.17	162	78664	26.49	ng	99
30) 1,2,4-Trichlorobenzene	10.37	180	79432	23.58	ng	97
31) Naphthalene	10.55	128	255752	25.60	ng	99
32) Benzoic acid	9.83	122	51778	24.28	ng	# 76
33) 4-Chloroaniline	10.67	127	124591	26.37	ng	98
34) Hexachlorobutadiene	10.84	225	48156	22.58	ng	95
35) Caprolactam	11.43	113	37138	24.88	ng	95
36) 4-Chloro-3-methylphenol	11.82	107	99577	26.54	ng	99
37) 2-Methylnaphthalene	12.17	142	195845	24.72	ng	98
39) 1,2,4,5-Tetrachlorobenzene	12.54	216	93603	23.67	ng	99
40) Hexachlorocyclopentadiene	12.52	237	51266	21.25	ng	94

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	12.79	196	68468	24.40	ng	97
43) 2,4,5-Trichlorophenol	12.88	196	74544	25.79	ng	94
45) 1,1'-Biphenyl	13.19	154	246597	24.49	ng	98
46) 2-Chloronaphthalene	13.23	162	196873	24.69	ng	98
47) 2-Nitroaniline	13.45	65	82093	26.47	ng	91
48) Acenaphthylene	14.08	152	342853	25.11	ng	96
49) Dimethylphthalate	13.82	163	261482	23.92	ng	98
50) 2,6-Dinitrotoluene	13.95	165	62181	25.67	ng	96
51) Acenaphthene	14.43	154	197638	24.50	ng	99
52) 3-Nitroaniline	14.27	138	70154	26.10	ng	96
53) 2,4-Dinitrophenol	14.48	184	28162	20.02	ng	92
54) Dibenzofuran	14.76	168	293448	24.47	ng	100
55) 4-Nitrophenol	14.63	139	52197	24.12	ng	90
56) 2,4-Dinitrotoluene	14.73	165	85260	25.23	ng	93
57) Fluorene	15.41	166	259872	24.49	ng	# 97
58) 2,3,4,6-Tetrachlorophenol	15.00	232	67121	25.30	ng	98
59) Diethylphthalate	15.19	149	277668	23.65	ng	98
60) 4-Chlorophenyl-phenylether	15.41	204	132632	24.33	ng	94
61) 4-Nitroaniline	15.44	138	79135	26.39	ng	93
62) Azobenzene	15.70	77	290881	25.93	ng	94
64) 4,6-Dinitro-2-methylphenol	15.50	198	47291	22.18	ng	98
65) n-Nitrosodiphenylamine	15.63	169	236905	24.81	ng	99
66) 4-Bromophenyl-phenylether	16.30	248	81523	24.03	ng	96
67) Hexachlorobenzene	16.41	284	86676	24.21	ng	100
68) Atrazine	16.58	200	88033	25.20	ng	97
69) Pentachlorophenol	16.77	266	49519	22.16	ng	96
70) Phenanthrene	17.15	178	411688	25.67	ng	98
71) Anthracene	17.24	178	412516	25.38	ng	99
72) Carbazole	17.52	167	387811	25.96	ng	99
73) Di-n-butylphthalate	18.08	149	517328	25.88	ng	98
74) Fluoranthene	19.17	202	487347	25.30	ng	97
76) Benzidine	19.36	184	223752	21.52	ng	96
77) Pyrene	19.53	202	488702	25.01	ng	97
79) Butylbenzylphthalate	20.44	149	227022	25.55	ng	97
80) Benzo(a)anthracene	21.28	228	447620	24.53	ng	99
81) 3,3'-Dichlorobenzidine	21.22	252	163056	23.09	ng	95
82) Chrysene	21.33	228	423681	24.92	ng	99
83) Bis(2-ethylhexyl)phthalate	21.21	149	315797	25.01	ng	99
84) Di-n-octyl phthalate	22.09	149	530173	25.22	ng	99
85) Indeno(1,2,3-cd)pyrene	25.88	276	492725	24.22	ng	97
87) Benzo(b)fluoranthene	22.88	252	434710	24.29	ng	# 98
88) Benzo(k)fluoranthene	22.92	252	437875m	25.77	ng	
89) Benzo(a)pyrene	23.46	252	404162	24.59	ng	97
90) Dibenzo(a,h)anthracene	25.89	278	416153	24.67	ng	99
91) Benzo(g,h,i)perylene	26.59	276	390594	24.60	ng	# 95

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

