

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060616\
 Data File : BG022167.D
 Acq On : 6 Jun 2016 13:41
 Operator : UM/SJ
 Sample : SSTDICC25
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC25

Manual Integrations
 APPROVED

sohil
 6/7/2016 7:27:27 PM

Quant Time: Jun 06 14:22:41 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 06 13:59:41 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.12	152	28111	20.00	ng	0.00
21) Naphthalene-d8	10.94	136	149328	20.00	ng	0.00
38) Acenaphthene-d10	14.75	164	97924	20.00	ng	0.00
63) Phenanthrene-d10	17.49	188	237363	20.00	ng	0.00
75) Chrysene-d12	21.77	240	242711	20.00	ng	0.00
86) Perylene-d12	25.06	264	232288	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.67	112	73956	43.81	ng	0.00
7) Phenol-d6	7.28	99	119361	47.34	ng	0.00
23) Nitrobenzene-d5	9.29	82	112742	38.81	ng	0.00
41) 2,4,6-Tribromophenol	16.23	330	57846	54.81	ng	0.00
44) 2-Fluorobiphenyl	13.37	172	330916	49.51	ng	0.00
78) Terphenyl-d14	20.08	244	512341	52.67	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.52	88	17400	23.16	ng	# 84
3) Pyridine	3.93	79	47121	20.90	ng	95
4) n-Nitrosodimethylamine	3.84	42	10584	9.47	ng	89
6) Aniline	7.44	93	75310	22.58	ng	# 82
8) 2-Chlorophenol	7.68	128	51556	25.48	ng	# 77
9) Benzaldehyde	7.26	77	35893	24.62	ng	# 71
10) Phenol	7.31	94	62319	22.98	ng	81
11) bis(2-Chloroethyl)ether	7.53	93	51395	23.68	ng	# 70
12) 1,3-Dichlorobenzene	8.01	146	53500	23.67	ng	# 93
13) 1,4-Dichlorobenzene	8.15	146	54879	23.85	ng	94
14) 1,2-Dichlorobenzene	8.48	146	53958	24.44	ng	98
15) Benzyl Alcohol	8.36	79	35712	18.53	ng	87
16) 2,2'-oxybis(1-Chloropropan	8.65	45	49724	10.69	ng	80
17) 2-Methylphenol	8.57	107	45394	24.21	ng	# 87
18) Hexachloroethane	9.21	117	20406	24.37	ng	# 79
19) n-Nitroso-di-n-propylamine	8.92	70	40792	20.82	ng	# 68
20) 3+4-Methylphenols	8.89	107	62872	24.50	ng	96
22) Acetophenone	8.95	105	85083	20.46	ng	# 81
24) Nitrobenzene	9.34	77	55541	17.85	ng	# 84
25) Isophorone	9.85	82	128001	20.13	ng	# 92
26) 2-Nitrophenol	10.05	139	29726	25.03	ng	# 72
27) 2,4-Dimethylphenol	10.11	122	57433	21.29	ng	88
28) bis(2-Chloroethoxy)methane	10.33	93	72158	21.36	ng	91
29) 2,4-Dichlorophenol	10.59	162	52618	22.88	ng	91
30) 1,2,4-Trichlorobenzene	10.80	180	56103	22.68	ng	94
31) Naphthalene	10.99	128	189069	23.66	ng	97
32) Benzoic acid	10.21	122	10852	7.56	ng	85
33) 4-Chloroaniline	11.10	127	79245	22.91	ng	# 89
34) Hexachlorobutadiene	11.26	225	29373	18.41	ng	91
35) Caprolactam	11.86	113	28115	26.71	ng	# 45
36) 4-Chloro-3-methylphenol	12.21	107	60558	20.94	ng	86
37) 2-Methylnaphthalene	12.58	142	141636m	25.82	ng	
39) 1,2,4,5-Tetrachlorobenzene	12.95	216	64027	20.92	ng	98
40) Hexachlorocyclopentadiene	12.93	237	27860	16.39	ng	90

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.19	196	43375	22.23	ng	98
43) 2,4,5-Trichlorophenol	13.27	196	48699	23.13	ng	92
45) 1,1'-Biphenyl	13.58	154	190019	23.03	ng	96
46) 2-Chloronaphthalene	13.63	162	145429	23.84	ng	94
47) 2-Nitroaniline	13.84	65	33935	16.22	ng	# 56
48) Acenaphthylene	14.48	152	255173	25.30	ng	99
49) Dimethylphthalate	14.19	163	207938	24.56	ng	100
50) 2,6-Dinitrotoluene	14.32	165	45670	28.29	ng	# 74
51) Acenaphthene	14.81	154	148364	23.01	ng	96
52) 3-Nitroaniline	14.66	138	50017	27.94	ng	# 77
53) 2,4-Dinitrophenol	14.88	184	11102	21.33	ng	# 72
54) Dibenzofuran	15.15	168	238352	27.08	ng	94
55) 4-Nitrophenol	14.98	139	33731	22.00	ng	# 64
56) 2,4-Dinitrotoluene	15.11	165	64126	29.39	ng	# 81
57) Fluorene	15.79	166	205027	26.08	ng	98
58) 2,3,4,6-Tetrachlorophenol	15.38	232	41285	23.63	ng	93
59) Diethylphthalate	15.55	149	224135	25.36	ng	99
60) 4-Chlorophenyl-phenylether	15.78	204	93631	24.18	ng	89
61) 4-Nitroaniline	15.82	138	51980	26.38	ng	# 78
62) Azobenzene	16.07	77	169118	22.33	ng	87
64) 4,6-Dinitro-2-methylphenol	15.88	198	26399	28.81	ng	# 81
65) n-Nitrosodiphenylamine	15.99	169	174259	24.48	ng	99
66) 4-Bromophenyl-phenylether	16.67	248	55774	21.93	ng	93
67) Hexachlorobenzene	16.80	284	65065	24.24	ng	94
68) Atrazine	16.93	200	65386	23.26	ng	99
69) Pentachlorophenol	17.14	266	23487	15.32	ng	95
70) Phenanthrene	17.53	178	325935	24.91	ng	98
71) Anthracene	17.63	178	332542	25.04	ng	99
72) Carbazole	17.90	167	309333	24.96	ng	98
73) Di-n-butylphthalate	18.43	149	406818	25.80	ng	# 99
74) Fluoranthene	19.54	202	377379	24.16	ng	97
76) Benzidine	19.72	184	176266	23.33	ng	97
77) Pyrene	19.90	202	376494	26.90	ng	99
79) Butylbenzylphthalate	20.76	149	173318	26.68	ng	90
80) Benzo(a)anthracene	21.74	228	338846	24.26	ng	100
81) 3,3'-Dichlorobenzidine	21.66	252	94004	8.50	ng	99
82) Chrysene	21.82	228	328983	24.52	ng	99
83) Bis(2-ethylhexyl)phthalate	21.62	149	258579	27.21	ng	98
84) Di-n-octyl phthalate	22.86	149	436904	27.45	ng	# 85
85) Indeno(1,2,3-cd)pyrene	28.84	276	371806	23.31	ng	100
87) Benzo(b)fluoranthene	24.01	252	331488	24.59	ng	# 97
88) Benzo(k)fluoranthene	24.08	252	337134	25.58	ng	# 96
89) Benzo(a)pyrene	24.91	252	324781	24.87	ng	98
90) Dibenzo(a,h)anthracene	28.90	278	304094	23.23	ng	97
91) Benzo(g,h,i)perylene	30.02	276	314962	24.51	ng	# 94

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

