

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042115\
 Data File : BG016585.D
 Acq On : 21 Apr 2015 16:19
 Operator : TP/IZ
 Sample : SSTDICC010
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

apatel
 4/22/2015 5:20:58 PM

Quant Time: Apr 22 01:58:34 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 22 01:14:44 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	49040	20.00	ng	0.00
21) Naphthalene-d8	10.42	136	229996	20.00	ng	0.00
38) Acenaphthene-d10	14.28	164	134457	20.00	ng	0.00
63) Phenanthrene-d10	17.03	188	289183	20.00	ng	0.00
75) Chrysene-d12	21.20	240	304066	20.00	ng	0.00
86) Perylene-d12	23.42	264	284899	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.23	112	61579	19.78	ng	0.00
7) Phenol-d6	6.83	99	84764	18.28	ng	0.00
23) Nitrobenzene-d5	8.79	82	75210	18.96	ng	0.00
41) 2,4,6-Tribromophenol	15.77	330	18051	19.46	ng	0.00
44) 2-Fluorobiphenyl	12.90	172	175165	21.96	ng	0.00
78) Terphenyl-d14	19.65	244	276561	21.17	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.11	88	14216	10.02	ng	96
3) Pyridine	3.51	79	38167	9.11	ng	97
4) n-Nitrosodimethylamine	3.42	42	11193	9.02	ng	89
6) Aniline	6.97	93	58659	8.91	ng	99
8) 2-Chlorophenol	7.21	128	36189	9.70	ng	100
9) Benzaldehyde	6.78	77	25873	9.92	ng	98
10) Phenol	6.86	94	44545	9.00	ng	100
11) bis(2-Chloroethyl)ether	7.07	93	36694	9.34	ng	97
12) 1,3-Dichlorobenzene	7.53	146	40477	10.71	ng	95
13) 1,4-Dichlorobenzene	7.67	146	40153	10.21	ng	99
14) 1,2-Dichlorobenzene	7.98	146	37429	10.00	ng	97
15) Benzyl Alcohol	7.88	79	25325	7.94	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.18	45	34741	8.02	ng	93
17) 2-Methylphenol	8.09	107	30310	9.15	ng	91
18) Hexachloroethane	8.71	117	13837	9.31	ng	85
19) n-Nitroso-di-n-propylamine	8.44	70	27172	8.36	ng	# 97
20) 3+4-Methylphenols	8.42	107	40933	8.91	ng	98
22) Acetophenone	8.46	105	51031	9.54	ng	# 99
24) Nitrobenzene	8.84	77	38479	9.10	ng	98
25) Isophorone	9.35	82	75378	8.63	ng	98
26) 2-Nitrophenol	9.55	139	19982	9.89	ng	94
27) 2,4-Dimethylphenol	9.62	122	35714	9.60	ng	97
28) bis(2-Chloroethoxy)methane	9.84	93	46327	9.26	ng	98
29) 2,4-Dichlorophenol	10.09	162	28153	9.50	ng	95
30) 1,2,4-Trichlorobenzene	10.29	180	32250	10.41	ng	98
31) Naphthalene	10.47	128	121893	10.07	ng	99
32) Benzoic acid	9.79	122	4525m	2.32	ng	
33) 4-Chloroaniline	10.59	127	53674	9.45	ng	98
34) Hexachlorobutadiene	10.75	225	14432	10.66	ng	96
35) Caprolactam	11.34	113	16029	8.62	ng	93
36) 4-Chloro-3-methylphenol	11.73	107	34933	9.00	ng	97
37) 2-Methylnaphthalene	12.09	142	87839	10.08	ng	97
39) 1,2,4,5-Tetrachlorobenzene	12.47	216	32284	10.81	ng	97
40) Hexachlorocyclopentadiene	12.44	237	4975	3.85	ng	87

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	12.71	196	20444	9.21	nq	96
43) 2,4,5-Trichlorophenol	12.80	196	22908	9.63	nq	96
45) 1,1'-Biphenyl	13.11	154	108163	10.40	nq	99
46) 2-Chloronaphthalene	13.15	162	81202	10.22	nq	98
47) 2-Nitroaniline	13.37	65	21232	8.36	nq	91
48) Acenaphthylene	14.00	152	145724	10.29	nq	99
49) Dimethylphthalate	13.74	163	100809	10.07	nq	97
50) 2,6-Dinitrotoluene	13.86	165	24049	9.91	nq	96
51) Acenaphthene	14.35	154	86371	10.21	nq	97
52) 3-Nitroaniline	14.20	138	28568	9.21	nq	99
53) 2,4-Dinitrophenol	14.42	184	3063	2.75	nq	94
54) Dibenzofuran	14.68	168	123275	10.45	nq	98
55) 4-Nitrophenol	14.53	139	17405	6.85	nq	96
56) 2,4-Dinitrotoluene	14.66	165	32712	9.78	nq	98
57) Fluorene	15.33	166	108386	10.41	nq	99
58) 2,3,4,6-Tetrachlorophenol	14.92	232	16861	9.14	nq	98
59) Diethylphthalate	15.11	149	106760	9.97	nq	98
60) 4-Chlorophenyl-phenylether	15.33	204	44614	10.51	nq	98
61) 4-Nitroaniline	15.36	138	33418	9.41	nq	98
62) Azobenzene	15.62	77	100866	9.03	nq	96
64) 4,6-Dinitro-2-methylphenol	15.42	198	13318	7.66	nq	97
65) n-Nitrosodiphenylamine	15.55	169	100073	10.18	nq	96
66) 4-Bromophenyl-phenylether	16.22	248	22766	9.84	nq	98
67) Hexachlorobenzene	16.33	284	24493	10.16	nq	97
68) Atrazine	16.49	200	28007	10.17	nq	100
69) Pentachlorophenol	16.69	266	6381	4.51	nq	94
70) Phenanthrene	17.07	178	173768	10.55	nq	98
71) Anthracene	17.15	178	170121	10.42	nq	98
72) Carbazole	17.43	167	174239	10.57	nq	100
73) Di-n-butylphthalate	17.99	149	213687	10.64	nq	99
74) Fluoranthene	19.08	202	192948	11.28	nq	99
76) Benzidine	19.28	184	117689	9.27	nq	99
77) Pyrene	19.45	202	209293	9.80	nq	98
79) Butylbenzylphthalate	20.35	149	100894	9.59	nq	98
80) Benzo(a)anthracene	21.19	228	191284	10.51	nq	98
81) 3,3'-Dichlorobenzidine	21.13	252	62083	10.41	nq	98
82) Chrysene	21.24	228	185315	10.54	nq	96
83) Bis(2-ethylhexyl)phthalate	21.12	149	141635	9.90	nq	98
84) Di-n-octyl phthalate	21.98	149	239781	10.30	nq	99
85) Indeno(1,2,3-cd)pyrene	25.67	276	192611	10.49	nq	99
87) Benzo(b)fluoranthene	22.75	252	178351	10.58	nq	97
88) Benzo(k)fluoranthene	22.80	252	167641	10.07	nq	98
89) Benzo(a)pyrene	23.32	252	162838	10.26	nq	98
90) Dibenzo(a,h)anthracene	25.68	278	158351	10.24	nq	96
91) Benzo(g,h,i)perylene	26.36	276	159594	10.30	nq	99

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

