

Data Path : Z:\HPCHEM1\BNA G\DATA\BG041316\
 Data File : BG021618.D
 Acq On : 13 Apr 2016 12:52
 Operator : UM/SJ
 Sample : SSTDICCC040
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICCC040

Manual Integrations
 APPROVED

Sohil
 4/14/2016 5:29:32 PM

Quant Time: Apr 13 13:50:19 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 13 13:39:40 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.22	152	30263	20.00	ng	0.00
21) Naphthalene-d8	11.04	136	138420	20.00	ng	0.00
38) Acenaphthene-d10	14.83	164	84004	20.00	ng	0.00
63) Phenanthrene-d10	17.56	188	190333	20.00	ng	0.00
75) Chrysene-d12	21.83	240	233740	20.00	ng	0.00
86) Perylene-d12	25.17	264	229289	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.76	112	144077	72.01	ng	0.00
7) Phenol-d6	7.36	99	213313	54.36	ng	0.00
23) Nitrobenzene-d5	9.38	82	213411	79.01	ng	0.00
41) 2,4,6-Tribromophenol	16.31	330	70305	55.77	ng	0.00
44) 2-Fluorobiphenyl	13.46	172	455024	92.01	ng	0.00
78) Terphenyl-d14	20.15	244	733889	69.47	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.64	88	32857	42.59	ng	91
3) Pyridine	4.04	79	94571	29.89	ng	92
4) n-Nitrosodimethylamine	3.96	42	45951	32.61	ng	# 88
6) Aniline	7.54	93	139264	27.05	ng	96
8) 2-Chlorophenol	7.78	128	85685	32.15	ng	95
9) Benzaldehyde	7.36	77	59018	30.62	ng	91
10) Phenol	7.39	94	115587	26.90	ng	95
11) bis(2-Chloroethyl)ether	7.63	93	90766	31.19	ng	93
12) 1,3-Dichlorobenzene	8.11	146	96704	39.68	ng	93
13) 1,4-Dichlorobenzene	8.26	146	98471	37.98	ng	97
14) 1,2-Dichlorobenzene	8.58	146	93575	35.58	ng	98
15) Benzyl Alcohol	8.45	79	80838	23.19	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.75	45	195453	31.31	ng	96
17) 2-Methylphenol	8.64	107	78619	25.71	ng	95
18) Hexachloroethane	9.31	117	35553	40.71	ng	# 74
19) n-Nitroso-di-n-propylamine	9.03	70	79744	22.35	ng	# 96
20) 3+4-Methylphenols	8.97	107	108195	23.49	ng	96
22) Acetophenone	9.04	105	151716	37.36	ng	# 99
24) Nitrobenzene	9.43	77	113971	39.89	ng	98
25) Isophorone	9.95	82	222041	29.07	ng	99
26) 2-Nitrophenol	10.14	139	43232	42.11	ng	97
27) 2,4-Dimethylphenol	10.19	122	94355	33.32	ng	96
28) bis(2-Chloroethoxy)methane	10.43	93	121912	36.52	ng	91
29) 2,4-Dichlorophenol	10.67	162	84419	35.44	ng	98
30) 1,2,4-Trichlorobenzene	10.89	180	91238	43.43	ng	92
31) Naphthalene	11.09	128	292266	39.25	ng	# 95
32) Benzoic acid	10.30	122	49935	15.65	ng	96
33) 4-Chloroaniline	11.19	127	122554	28.08	ng	97
34) Hexachlorobutadiene	11.36	225	58421	50.04	ng	96
35) Caprolactam	11.96	113	37559	13.50	ng	95
36) 4-Chloro-3-methylphenol	12.28	107	102921	23.55	ng	98
37) 2-Methylnaphthalene	12.68	142	198086	32.22	ng	98
39) 1,2,4,5-Tetrachlorobenzene	13.03	216	104357	51.43	ng	98
40) Hexachlorocyclopentadiene	13.02	237	60600	70.78	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.26	196	67351	43.95	nq	94
43) 2,4,5-Trichlorophenol	13.33	196	73363	39.35	nq	97
45) 1,1'-Biphenyl	13.67	154	281696	45.19	nq	97
46) 2-Chloronaphthalene	13.72	162	207367	44.72	nq	99
47) 2-Nitroaniline	13.90	65	69901	32.17	nq	96
48) Acenaphthylene	14.56	152	339943	37.94	nq	98
49) Dimethylphthalate	14.27	163	275559	29.62	nq	99
50) 2,6-Dinitrotoluene	14.40	165	54926	32.66	nq	96
51) Acenaphthene	14.89	154	218836	37.69	nq	97
52) 3-Nitroaniline	14.72	138	59804	25.98	nq	96
53) 2,4-Dinitrophenol	14.92	184	17001	24.63	nq	# 87
54) Dibenzofuran	15.22	168	294504	34.86	nq	99
55) 4-Nitrophenol	15.01	139	53594	21.78	nq	89
56) 2,4-Dinitrotoluene	15.17	165	74127	27.42	nq	98
57) Fluorene	15.87	166	260491	31.45	nq	100
58) 2,3,4,6-Tetrachlorophenol	15.44	232	58286m	29.53	nq	
59) Diethylphthalate	15.63	149	289348	25.75	nq	97
60) 4-Chlorophenyl-phenylether	15.86	204	128080	33.68	nq	95
61) 4-Nitroaniline	15.88	138	68704	21.54	nq	97
62) Azobenzene	16.15	77	251480	30.04	nq	99
64) 4,6-Dinitro-2-methylphenol	15.93	198	30443	41.86	nq	93
65) n-Nitrosodiphenylamine	16.07	169	228624	45.74	nq	96
66) 4-Bromophenyl-phenylether	16.75	248	82294	47.16	nq	98
67) Hexachlorobenzene	16.86	284	83592	46.37	nq	98
68) Atrazine	17.01	200	95820	32.99	nq	96
69) Pentachlorophenol	17.20	266	51018	41.99	nq	94
70) Phenanthrene	17.60	178	416342	38.64	nq	99
71) Anthracene	17.69	178	427106	38.36	nq	98
72) Carbazole	17.96	167	403861	32.78	nq	97
73) Di-n-butylphthalate	18.51	149	535191	30.89	nq	99
74) Fluoranthene	19.60	202	517642	30.58	nq	99
76) Benzidine	19.77	184	293671	41.90	nq	99
77) Pyrene	19.95	202	526303	35.63	nq	100
79) Butylbenzylphthalate	20.84	149	253449	41.15	nq	95
80) Benzo(a)anthracene	21.82	228	537698	39.43	nq	99
81) 3,3'-Dichlorobenzidine	21.72	252	426584	45.79	nq	100
82) Chrysene	21.89	228	512124	39.53	nq	99
83) Bis(2-ethylhexyl)phthalate	21.72	149	362235	44.92	nq	# 99
84) Di-n-octyl phthalate	22.97	149	622197	46.64	nq	100
85) Indeno(1,2,3-cd)pyrene	28.99	276	615555	46.23	nq	98
87) Benzo(b)fluoranthene	24.11	252	539341	39.98	nq	98
88) Benzo(k)fluoranthene	24.18	252	511384	37.96	nq	99
89) Benzo(a)pyrene	25.01	252	509456	39.18	nq	98
90) Dibenzo(a,h)anthracene	29.06	278	515958	40.33	nq	99
91) Benzo(g,h,i)perylene	30.18	276	505788	39.73	nq	98

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

