

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042315\
 Data File : BG016667.D
 Acq On : 23 Apr 2015 23:46
 Operator : TP/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040EC

Quant Time: Apr 24 02:27:41 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 23 19:19:11 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	52897	20.00	ng	0.00
21) Naphthalene-d8	10.38	136	236012	20.00	ng	0.00
38) Acenaphthene-d10	14.25	164	133252	20.00	ng	0.00
63) Phenanthrene-d10	17.00	188	270986	20.00	ng	0.00
75) Chrysene-d12	21.18	240	255497	20.00	ng	0.00
86) Perylene-d12	23.38	264	236335	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.19	112	277629	85.17	ng	0.00
7) Phenol-d6	6.80	99	387409	84.72	ng	0.00
23) Nitrobenzene-d5	8.76	82	324831	86.97	ng	0.00
41) 2,4,6-Tribromophenol	15.75	330	82863	83.82	ng	0.00
44) 2-Fluorobiphenyl	12.87	172	707723	85.33	ng	0.00
78) Terphenyl-d14	19.63	244	910822	82.12	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.05	88	62141	44.01	ng	100
3) Pyridine	3.45	79	175207	43.70	ng	96
4) n-Nitrosodimethylamine	3.38	42	49172	41.57	ng	89
6) Aniline	6.93	93	264901	42.24	ng	98
8) 2-Chlorophenol	7.17	128	168403	41.80	ng	98
9) Benzaldehyde	6.74	77	69376	37.13	ng	95
10) Phenol	6.82	94	202386	41.99	ng	98
11) bis(2-Chloroethyl)ether	7.03	93	154902	41.17	ng	97
12) 1,3-Dichlorobenzene	7.48	146	171846	41.24	ng	99
13) 1,4-Dichlorobenzene	7.63	146	178631	41.94	ng	99
14) 1,2-Dichlorobenzene	7.95	146	168070	41.33	ng	98
15) Benzyl Alcohol	7.85	79	122885	42.82	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.14	45	144067	40.91	ng	99
17) 2-Methylphenol	8.06	107	137828	42.77	ng	98
18) Hexachloroethane	8.67	117	64005	42.62	ng	99
19) n-Nitroso-di-n-propylamine	8.41	70	120233	42.18	ng	97
20) 3+4-Methylphenols	8.39	107	191731	42.31	ng	98
22) Acetophenone	8.42	105	223819	43.21	ng	# 99
24) Nitrobenzene	8.80	77	169990	43.34	ng	97
25) Isophorone	9.32	82	332149	42.91	ng	100
26) 2-Nitrophenol	9.51	139	100517	44.16	ng	98
27) 2,4-Dimethylphenol	9.58	122	162917	42.86	ng	97
28) bis(2-Chloroethoxy)methane	9.81	93	198603	42.52	ng	97
29) 2,4-Dichlorophenol	10.05	162	139054	43.04	ng	99
30) 1,2,4-Trichlorobenzene	10.25	180	143948	42.04	ng	97
31) Naphthalene	10.43	128	521928	41.94	ng	99
32) Benzoic acid	9.74	122	63610	34.03	ng	97
33) 4-Chloroaniline	10.55	127	241235	42.08	ng	97
34) Hexachlorobutadiene	10.72	225	64789	42.36	ng	98
35) Caprolactam	11.32	113	74108	42.55	ng	94
36) 4-Chloro-3-methylphenol	11.70	107	160852	42.82	ng	96
37) 2-Methylnaphthalene	12.06	142	373141	41.52	ng	99
39) 1,2,4,5-Tetrachlorobenzene	12.43	216	134209	42.28	ng	99
40) Hexachlorocyclopentadiene	12.40	237	45110	40.61	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	12.68	196	99173	42.75	nq	99
43) 2,4,5-Trichlorophenol	12.76	196	105021	42.78	nq	98
45) 1,1'-Biphenyl	13.08	154	443030	42.33	nq	99
46) 2-Chloronaphthalene	13.12	162	340348	42.31	nq	98
47) 2-Nitroaniline	13.34	65	96518	43.99	nq	98
48) Acenaphthylene	13.97	152	599583	42.56	nq	99
49) Dimethylphthalate	13.72	163	417277	41.86	nq	100
50) 2,6-Dinitrotoluene	13.84	165	102575	41.93	nq	95
51) Acenaphthene	14.31	154	352185	41.65	nq	99
52) 3-Nitroaniline	14.17	138	125457	42.44	nq	99
53) 2,4-Dinitrophenol	14.39	184	43451	37.17	nq	97
54) Dibenzofuran	14.65	168	491524	41.21	nq	99
55) 4-Nitrophenol	14.51	139	93791	42.34	nq	95
56) 2,4-Dinitrotoluene	14.63	165	144711	43.19	nq	99
57) Fluorene	15.30	166	432812	41.51	nq	99
58) 2,3,4,6-Tetrachlorophenol	14.89	232	76795	41.35	nq	99
59) Diethylphthalate	15.08	149	433821	41.76	nq	100
60) 4-Chlorophenyl-phenylether	15.29	204	176574	40.55	nq	99
61) 4-Nitroaniline	15.34	138	148336	45.36	nq	98
62) Azobenzene	15.59	77	394964	41.63	nq	99
64) 4,6-Dinitro-2-methylphenol	15.40	198	77719	42.38	nq	97
65) n-Nitrosodiphenylamine	15.52	169	395257	42.54	nq	99
66) 4-Bromophenyl-phenylether	16.19	248	93999	41.51	nq	96
67) Hexachlorobenzene	16.30	284	98153	42.14	nq	99
68) Atrazine	16.47	200	105094	42.14	nq	98
69) Pentachlorophenol	16.66	266	45617	40.34	nq	98
70) Phenanthrene	17.04	178	657319	42.22	nq	100
71) Anthracene	17.13	178	662355	42.77	nq	99
72) Carbazole	17.40	167	676575	43.61	nq	99
73) Di-n-butylphthalate	17.97	149	821097	44.41	nq	99
74) Fluoranthene	19.06	202	712112	43.17	nq	98
76) Benzidine	19.25	184	355735	40.88	nq	100
77) Pyrene	19.42	202	734587	40.93	nq	99
79) Butylbenzylphthalate	20.32	149	378714	42.89	nq	97
80) Benzo(a)anthracene	21.17	228	641027	41.38	nq	98
81) 3,3'-Dichlorobenzidine	21.10	252	214863	42.55	nq	97
82) Chrysene	21.22	228	612670	41.25	nq	100
83) Bis(2-ethylhexyl)phthalate	21.10	149	523966	43.62	nq	98
84) Di-n-octyl phthalate	21.96	149	876221	43.79	nq	99
85) Indeno(1,2,3-cd)pyrene	25.63	276	673025	41.58	nq	100
87) Benzo(b)fluoranthene	22.73	252	597296	41.61	nq	98
88) Benzo(k)fluoranthene	22.77	252	604526	43.15	nq	99
89) Benzo(a)pyrene	23.29	252	570055	42.14	nq	99
90) Dibenzo(a,h)anthracene	25.64	278	553895	42.11	nq	97
91) Benzo(g,h,i)perylene	26.32	276	562562	42.03	nq	99

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

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