

Data Path : Z:\HPCHEM1\BNA G\DATA\BG040915\
 Data File : BG016304.D
 Acq On : 10 Apr 2015 00:19
 Operator : TP/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Apr 10 03:46:45 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG040615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 07 04:17:17 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	81203	20.00	ng	0.00
21) Naphthalene-d8	10.48	136	408070	20.00	ng	0.00
38) Acenaphthene-d10	14.34	164	248192	20.00	ng	0.00
63) Phenanthrene-d10	17.08	188	495204	20.00	ng	0.00
75) Chrysene-d12	21.26	240	406156	20.00	ng	0.00
86) Perylene-d12	23.50	264	391032	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.28	112	402305	78.19	ng	0.00
7) Phenol-d6	6.89	99	636947	82.46	ng	0.02
23) Nitrobenzene-d5	8.86	82	531926	75.51	ng	0.00
41) 2,4,6-Tribromophenol	15.83	330	139168	82.91	ng	0.00
44) 2-Fluorobiphenyl	12.96	172	1083612	74.34	ng	0.00
78) Terphenyl-d14	19.70	244	1284639	74.03	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.17	88	80377	34.36	ng	98
3) Pyridine	3.57	79	254468	36.40	ng	96
4) n-Nitrosodimethylamine	3.49	42	75469	36.32	ng	98
6) Aniline	7.03	93	417178	37.82	ng	99
8) 2-Chlorophenol	7.27	128	253614	41.04	ng	97
9) Benzaldehyde	6.84	77	150652	33.30	ng	98
10) Phenol	6.91	94	340364	41.18	ng	100
11) bis(2-Chloroethyl)ether	7.13	93	240848	36.55	ng	100
12) 1,3-Dichlorobenzene	7.59	146	237834	38.12	ng	98
13) 1,4-Dichlorobenzene	7.73	146	244849	37.83	ng	97
14) 1,2-Dichlorobenzene	8.05	146	235572	38.06	ng	98
15) Benzyl Alcohol	7.94	79	207158	38.47	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.23	45	259711	34.99	ng	98
17) 2-Methylphenol	8.15	107	227154	40.99	ng	97
18) Hexachloroethane	8.77	117	92382	37.31	ng	95
19) n-Nitroso-di-n-propylamine	8.50	70	201464	36.40	ng	98
20) 3+4-Methylphenols	8.48	107	320041	41.69	ng	96
22) Acetophenone	8.52	105	350710	37.06	ng	# 99
24) Nitrobenzene	8.90	77	280821	37.28	ng	95
25) Isophorone	9.41	82	563089	36.00	ng	99
26) 2-Nitrophenol	9.60	139	152153	43.32	ng	98
27) 2,4-Dimethylphenol	9.67	122	269319	41.16	ng	100
28) bis(2-Chloroethoxy)methane	9.90	93	332900	37.28	ng	99
29) 2,4-Dichlorophenol	10.14	162	230009	44.43	ng	97
30) 1,2,4-Trichlorobenzene	10.35	180	211153	39.06	ng	99
31) Naphthalene	10.54	128	807649	37.98	ng	100
32) Benzoic acid	9.83	122	180961	43.42	ng	96
33) 4-Chloroaniline	10.65	127	402736	40.36	ng	99
34) Hexachlorobutadiene	10.82	225	93682	39.46	ng	99
35) Caprolactam	11.42	113	132066	40.52	ng	99
36) 4-Chloro-3-methylphenol	11.79	107	295047	43.14	ng	98
37) 2-Methylnaphthalene	12.15	142	593685	38.81	ng	99
39) 1,2,4,5-Tetrachlorobenzene	12.52	216	209390	38.32	ng	99
40) Hexachlorocyclopentadiene	12.50	237	92002	36.78	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	12.77	196	172591	42.50	nq	98
43) 2,4,5-Trichlorophenol	12.85	196	187618	43.25	nq	99
45) 1,1'-Biphenyl	13.17	154	707677	37.17	nq	99
46) 2-Chloronaphthalene	13.21	162	546772	37.70	nq	99
47) 2-Nitroaniline	13.42	65	186531	39.90	nq	99
48) Acenaphthylene	14.06	152	973192	37.74	nq	99
49) Dimethylphthalate	13.80	163	670845	36.88	nq	99
50) 2,6-Dinitrotoluene	13.92	165	174298	39.62	nq	97
51) Acenaphthene	14.40	154	586618	38.07	nq	98
52) 3-Nitroaniline	14.25	138	221073	39.35	nq	98
53) 2,4-Dinitrophenol	14.46	184	69876	32.87	nq	99
54) Dibenzofuran	14.74	168	801948	37.50	nq	100
55) 4-Nitrophenol	14.57	139	174161	37.51	nq	99
56) 2,4-Dinitrotoluene	14.71	165	238595	39.82	nq	93
57) Fluorene	15.38	166	707476	37.50	nq	100
58) 2,3,4,6-Tetrachlorophenol	14.97	232	138706	41.43	nq	96
59) Diethylphthalate	15.17	149	711208	36.73	nq	99
60) 4-Chlorophenyl-phenylether	15.38	204	288184	37.34	nq	98
61) 4-Nitroaniline	15.42	138	243122	38.12	nq	97
62) Azobenzene	15.67	77	739369	35.88	nq	99
64) 4,6-Dinitro-2-methylphenol	15.47	198	119039	35.27	nq	97
65) n-Nitrosodiphenylamine	15.60	169	658939	39.36	nq	99
66) 4-Bromophenyl-phenylether	16.27	248	155002	39.20	nq	97
67) Hexachlorobenzene	16.39	284	161919	39.40	nq	96
68) Atrazine	16.55	200	178289	38.21	nq	100
69) Pentachlorophenol	16.74	266	98036	39.07	nq	97
70) Phenanthrene	17.12	178	1044576	37.50	nq	99
71) Anthracene	17.21	178	1051227	38.08	nq	99
72) Carbazole	17.49	167	1021055	36.86	nq	99
73) Di-n-butylphthalate	18.05	149	1262066	37.23	nq	99
74) Fluoranthene	19.14	202	1044041	36.37	nq	98
76) Benzidine	19.33	184	549809	31.92	nq	99
77) Pyrene	19.50	202	1065164	37.66	nq	99
79) Butylbenzylphthalate	20.40	149	557640	40.13	nq	95
80) Benzo(a)anthracene	21.24	228	904368	37.68	nq	100
81) 3,3'-Dichlorobenzidine	21.18	252	310588	39.35	nq	99
82) Chrysene	21.29	228	879827	37.97	nq	99
83) Bis(2-ethylhexyl)phthalate	21.17	149	750624	39.90	nq	98
84) Di-n-octyl phthalate	22.05	149	1260116	41.12	nq	99
85) Indeno(1,2,3-cd)pyrene	25.79	276	970210	40.30	nq	99
87) Benzo(b)fluoranthene	22.82	252	866351	37.83	nq	100
88) Benzo(k)fluoranthene	22.87	252	842312	37.58	nq	98
89) Benzo(a)pyrene	23.40	252	814270	37.94	nq	99
90) Dibenzo(a,h)anthracene	25.81	278	802130	38.40	nq	98
91) Benzo(g,h,i)perylene	26.49	276	811878	38.91	nq	99

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

