

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042715\
 Data File : BG016730.D
 Acq On : 27 Apr 2015 11:11
 Operator : TP/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Apr 27 14:39:10 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	7.57	152	57313	20.00	ng	-0.03
21) Naphthalene-d8	10.36	136	255220	20.00	ng	-0.02
38) Acenaphthene-d10	14.24	164	148734	20.00	ng	-0.02
63) Phenanthrene-d10	16.99	188	310613	20.00	ng	0.00
75) Chrysene-d12	21.17	240	294983	20.00	ng	0.00
86) Perylene-d12	23.37	264	293536	20.00	ng	-0.02

System Monitoring Compounds						
5) 2-Fluorophenol	5.17	112	288174	81.59	ng	-0.02
7) Phenol-d6	6.78	99	410071	82.76	ng	-0.02
23) Nitrobenzene-d5	8.74	82	333561	82.58	ng	-0.02
41) 2,4,6-Tribromophenol	15.73	330	99056	89.77	ng	-0.02
44) 2-Fluorobiphenyl	12.85	172	776345	83.86	ng	-0.02
78) Terphenyl-d14	19.62	244	1012759	79.09	ng	-0.02

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.03	88	60212	39.35	ng	100
3) Pyridine	3.43	79	173616	39.97	ng	99
4) n-Nitrosodimethylamine	3.35	42	50547	39.44	ng	97
6) Aniline	6.91	93	277970	40.91	ng	98
8) 2-Chlorophenol	7.15	128	178759	40.95	ng	98
9) Benzaldehyde	6.72	77	77419	38.24	ng	97
10) Phenol	6.81	94	213078	40.80	ng	97
11) bis(2-Chloroethyl)ether	7.01	93	159180	39.05	ng	97
12) 1,3-Dichlorobenzene	7.46	146	184919	40.96	ng	97
13) 1,4-Dichlorobenzene	7.61	146	188229	40.79	ng	99
14) 1,2-Dichlorobenzene	7.93	146	180094	40.88	ng	100
15) Benzyl Alcohol	7.83	79	132041	42.47	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.12	45	144165	37.78	ng	98
17) 2-Methylphenol	8.05	107	147749	42.31	ng	96
18) Hexachloroethane	8.64	117	66900	41.12	ng	98
19) n-Nitroso-di-n-propylamine	8.39	70	125240	40.55	ng	97
20) 3+4-Methylphenols	8.38	107	205197	41.80	ng	95
22) Acetophenone	8.40	105	236322	42.19	ng	98
24) Nitrobenzene	8.78	77	173081	40.80	ng	97
25) Isophorone	9.30	82	360473	43.06	ng	99
26) 2-Nitrophenol	9.49	139	110143	44.75	ng	97
27) 2,4-Dimethylphenol	9.56	122	174205	42.38	ng	99
28) bis(2-Chloroethoxy)methane	9.79	93	205747	40.74	ng	98
29) 2,4-Dichlorophenol	10.03	162	157222	45.00	ng	98
30) 1,2,4-Trichlorobenzene	10.23	180	157649	42.58	ng	99
31) Naphthalene	10.41	128	553907	41.16	ng	99
32) Benzoic acid	9.73	122	107303	48.13	ng	100
33) 4-Chloroaniline	10.53	127	265506	42.82	ng	97
34) Hexachlorobutadiene	10.70	225	70341	42.53	ng	97
35) Caprolactam	11.32	113	92573	49.15	ng	97
36) 4-Chloro-3-methylphenol	11.69	107	179566	44.20	ng	98
37) 2-Methylnaphthalene	12.04	142	405607	41.73	ng	99
39) 1,2,4,5-Tetrachlorobenzene	12.41	216	152733	43.11	ng	99
40) Hexachlorocyclopentadiene	12.38	237	44657	36.84	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	12.67	196	118944	45.94	nq	97
43) 2,4,5-Trichlorophenol	12.75	196	125733	45.89	nq	96
45) 1,1'-Biphenyl	13.07	154	491142	42.05	nq	99
46) 2-Chloronaphthalene	13.10	162	378919	42.20	nq	98
47) 2-Nitroaniline	13.32	65	107490	43.89	nq	96
48) Acenaphthylene	13.95	152	663779	42.21	nq	100
49) Dimethylphthalate	13.71	163	474655	42.66	nq	99
50) 2,6-Dinitrotoluene	13.83	165	119464	43.75	nq	98
51) Acenaphthene	14.30	154	392346	41.57	nq	100
52) 3-Nitroaniline	14.16	138	146799	44.49	nq	94
53) 2,4-Dinitrophenol	14.38	184	50664	38.49	nq	93
54) Dibenzofuran	14.64	168	551784	41.44	nq	100
55) 4-Nitrophenol	14.51	139	106631	43.12	nq	96
56) 2,4-Dinitrotoluene	14.62	165	166151	44.43	nq	99
57) Fluorene	15.29	166	495450	42.57	nq	98
58) 2,3,4,6-Tetrachlorophenol	14.88	232	93643	45.17	nq	97
59) Diethylphthalate	15.07	149	499487	43.08	nq	100
60) 4-Chlorophenyl-phenylether	15.28	204	203609	41.89	nq	100
61) 4-Nitroaniline	15.33	138	169251	46.37	nq	99
62) Azobenzene	15.58	77	424411	40.08	nq	98
64) 4,6-Dinitro-2-methylphenol	15.39	198	91816	43.68	nq	95
65) n-Nitrosodiphenylamine	15.50	169	452845	42.52	nq	99
66) 4-Bromophenyl-phenylether	16.17	248	110385	42.53	nq	97
67) Hexachlorobenzene	16.29	284	112358	42.09	nq	98
68) Atrazine	16.46	200	127390	44.56	nq	98
69) Pentachlorophenol	16.64	266	56904	43.90	nq	96
70) Phenanthrene	17.03	178	741308	41.54	nq	98
71) Anthracene	17.11	178	749419	42.22	nq	100
72) Carbazole	17.40	167	763658	42.95	nq	99
73) Di-n-butylphthalate	17.95	149	932902	44.02	nq	99
74) Fluoranthene	19.05	202	807357	42.70	nq	99
76) Benzidine	19.24	184	418614	41.67	nq	99
77) Pyrene	19.41	202	823243	39.73	nq	99
79) Butylbenzylphthalate	20.32	149	444939	43.65	nq	95
80) Benzo(a)anthracene	21.16	228	745061	41.65	nq	99
81) 3,3'-Dichlorobenzidine	21.10	252	261904	44.93	nq	98
82) Chrysene	21.21	228	707026	41.23	nq	100
83) Bis(2-ethylhexyl)phthalate	21.09	149	616017	44.42	nq	99
84) Di-n-octyl phthalate	21.95	149	1051494	45.52	nq	100
85) Indeno(1,2,3-cd)pyrene	25.62	276	830989	44.46	nq	100
87) Benzo(b)fluoranthene	22.71	252	707373	39.67	nq	99
88) Benzo(k)fluoranthene	22.75	252	703212	40.41	nq	99
89) Benzo(a)pyrene	23.28	252	689749	41.05	nq	99
90) Dibenzo(a,h)anthracene	25.62	278	680869	41.67	nq	99
91) Benzo(g,h,i)perylene	26.30	276	683377	41.11	nq	97

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

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