

Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG062315\  
 Data File : BG017757.D  
 Acq On : 23 Jun 2015 12:59  
 Operator : TP/IZ  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 Client Sampled :  
 SSTDCCC040

Manual Integrations  
 APPROVED

apatel  
 6/25/2015 4:56:15 PM

Quant Time: Jun 24 00:50:36 2015  
 Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG061815.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Jun 19 06:05:07 2015  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	74273	20.00	ng	-0.01
21) Naphthalene-d8	10.39	136	353303	20.00	ng	-0.01
38) Acenaphthene-d10	14.26	164	217192	20.00	ng	-0.01
63) Phenanthrene-d10	17.01	188	455243	20.00	ng	-0.02
75) Chrysene-d12	21.22	240	454281	20.00	ng	-0.02
86) Perylene-d12	23.44	264	429881	20.00	ng	-0.02

## System Monitoring Compounds

5) 2-Fluorophenol	5.21	112	358589	84.73	ng	-0.02
7) Phenol-d6	6.78	99	518166	87.11	ng	-0.02
23) Nitrobenzene-d5	8.76	82	544293	88.62	ng	-0.01
41) 2,4,6-Tribromophenol	15.76	330	160116	79.50	ng	-0.01
44) 2-Fluorobiphenyl	12.88	172	1046294	78.24	ng	-0.02
78) Terphenyl-d14	19.66	244	1570170	78.54	ng	-0.02

## Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.17	88	83691	45.40	ng	96
3) Pyridine	3.55	79	253158	49.34	ng	97
4) n-Nitrosodimethylamine	3.47	42	115653	49.59	ng	# 99
6) Aniline	6.94	93	357345	43.69	ng	99
8) 2-Chlorophenol	7.17	128	215061	40.62	ng	91
9) Benzaldehyde	6.76	77	133122	38.60	ng	96
10) Phenol	6.81	94	283122	43.66	ng	97
11) bis(2-Chloroethyl)ether	7.04	93	216852	44.10	ng	97
12) 1,3-Dichlorobenzene	7.50	146	228226	38.76	ng	95
13) 1,4-Dichlorobenzene	7.64	146	238340	39.16	ng	97
14) 1,2-Dichlorobenzene	7.96	146	224520	39.08	ng	97
15) Benzyl Alcohol	7.85	79	224991	43.92	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.15	45	344612	45.92	ng	97
17) 2-Methylphenol	8.05	107	198993	41.22	ng	96
18) Hexachloroethane	8.68	117	95588	40.96	ng	91
19) n-Nitroso-di-n-propylamine	8.42	70	215533	43.05	ng	95
20) 3+4-Methylphenols	8.38	107	272221	41.80	ng	96
22) Acetophenone	8.43	105	329562	41.13	ng	# 96
24) Nitrobenzene	8.81	77	293588	44.48	ng	99
25) Isophorone	9.33	82	564497	41.26	ng	95
26) 2-Nitrophenol	9.51	139	130193	43.23	ng	94
27) 2,4-Dimethylphenol	9.58	122	221120	38.77	ng	95
28) bis(2-Chloroethoxy)methane	9.82	93	291867	42.33	ng	98
29) 2,4-Dichlorophenol	10.04	162	197608	39.31	ng	93
30) 1,2,4-Trichlorobenzene	10.26	180	209273	37.07	ng	99
31) Naphthalene	10.44	128	756122	39.59	ng	99
32) Benzoic acid	9.71	122	132835	43.32	ng	91
33) 4-Chloroaniline	10.56	127	335288	40.12	ng	97
34) Hexachlorobutadiene	10.73	225	120235	37.21	ng	94
35) Caprolactam	11.33	113	107273	38.42	ng	87
36) 4-Chloro-3-methylphenol	11.69	107	248215	39.91	ng	95
37) 2-Methylnaphthalene	12.07	142	534087	38.87	ng	99
39) 1,2,4,5-Tetrachlorobenzene	12.44	216	222311	38.16	ng	97
40) Hexachlorocyclopentadiene	12.42	237	140386	35.88	ng	94

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	12.68	196	157552	39.95	ng	99
43) 2,4,5-Trichlorophenol	12.75	196	174120	40.49	ng	97
45) 1,1'-Biphenyl	13.10	154	632704	39.43	ng	97
46) 2-Chloronaphthalene	13.13	162	507177	39.26	ng	99
47) 2-Nitroaniline	13.34	65	189962	49.20	ng	92
48) Acenaphthylene	13.98	152	914506	39.62	ng	98
49) Dimethylphthalate	13.74	163	632400	38.07	ng	99
50) 2,6-Dinitrotoluene	13.85	165	148518	43.36	ng	# 89
51) Acenaphthene	14.33	154	551762	39.61	ng	99
52) 3-Nitroaniline	14.18	138	176456	44.09	ng	92
53) 2,4-Dinitrophenol	14.38	184	66984	43.98	ng	# 83
54) Dibenzofuran	14.66	168	769326	39.22	ng	98
55) 4-Nitrophenol	14.49	139	140965	44.58	ng	95
56) 2,4-Dinitrotoluene	14.63	165	198809	40.83	ng	# 85
57) Fluorene	15.32	166	687804	38.80	ng	99
58) 2,3,4,6-Tetrachlorophenol	14.89	232	149151m	41.82	ng	
59) Diethylphthalate	15.11	149	674110	37.41	ng	100
60) 4-Chlorophenyl-phenylether	15.32	204	313105	38.48	ng	99
61) 4-Nitroaniline	15.35	138	198738	44.73	ng	97
62) Azobenzene	15.62	77	802629	42.51	ng	98
64) 4,6-Dinitro-2-methylphenol	15.40	198	113522	46.14	ng	91
65) n-Nitrosodiphenylamine	15.53	169	587636	40.40	ng	98
66) 4-Bromophenyl-phenylether	16.22	248	181319	38.56	ng	99
67) Hexachlorobenzene	16.32	284	197072	38.69	ng	98
68) Atrazine	16.50	200	199772	37.63	ng	99
69) Pentachlorophenol	16.67	266	105070	41.12	ng	99
70) Phenanthrene	17.06	178	1023391	39.85	ng	99
71) Anthracene	17.15	178	1017172	39.67	ng	98
72) Carbazole	17.43	167	994560	39.58	ng	100
73) Di-n-butylphthalate	18.00	149	1225362	39.06	ng	99
74) Fluoranthene	19.08	202	1125325	38.39	ng	97
76) Benzidine	19.28	184	515001	33.99	ng	99
77) Pyrene	19.45	202	1139073	39.40	ng	99
79) Butylbenzylphthalate	20.37	149	552103	40.38	ng	93
80) Benzo(a)anthracene	21.20	228	1023863	39.15	ng	99
81) 3,3'-Dichlorobenzidine	21.14	252	367171	37.64	ng	99
82) Chrysene	21.26	228	943815	39.32	ng	99
83) Bis(2-ethylhexyl)phthalate	21.15	149	763639	40.20	ng	98
84) Di-n-octyl phthalate	22.03	149	1255226	40.26	ng	97
85) Indeno(1,2,3-cd)pyrene	25.71	276	1121083	38.94	ng	97
87) Benzo(b)fluoranthene	22.77	252	1018338	39.35	ng	99
88) Benzo(k)fluoranthene	22.82	252	984670	39.80	ng	97
89) Benzo(a)pyrene	23.35	252	934529	39.31	ng	98
90) Dibenzo(a,h)anthracene	25.72	278	931321	38.70	ng	99
91) Benzo(g,h,i)perylene	26.40	276	900436	38.61	ng	97

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

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