

Data Path : Z:\HPCHEM1\BNA G\DATA\BG102017\
 Data File : BG029380.D
 Acq On : 20 Oct 2017 20:06
 Operator : SJ/JU
 Sample : I5981-06MSD
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 NB-01-101817-BMSD

Manual Integrations
 APPROVED

Sohil
 10/23/2017 3:40:48 PM

Quant Time: Oct 21 22:51:25 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG101617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 16 17:32:15 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.17	152	57056	20.00	ng	0.00
21) Naphthalene-d8	10.98	136	250313	20.00	ng	0.00
38) Acenaphthene-d10	14.79	164	175375	20.00	ng	0.00
63) Phenanthrene-d10	17.52	188	455421	20.00	ng	0.00
75) Chrysene-d12	21.79	240	494447	20.00	ng	0.00
86) Perylene-d12	25.10	264	528987	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.72	112	476743	139.59	ng	0.00
7) Phenol-d6	7.32	99	597347	124.60	ng	0.00
23) Nitrobenzene-d5	9.33	82	398403	101.14	ng	0.00
41) 2,4,6-Tribromophenol	16.27	330	403532	178.22	ng	0.00
44) 2-Fluorobiphenyl	13.41	172	1060810	88.72	ng	0.00
78) Terphenyl-d14	20.11	244	2001148	92.07	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.59	88	56963	41.106	ng	# 85
3) Pyridine	3.99	79	134266	33.272	ng	97
4) n-Nitrosodimethylamine	3.90	42	84980	45.050	ng	# 93
6) Aniline	7.49	93	154579	26.039	ng	94
8) 2-Chlorophenol	7.73	128	177429	45.322	ng	93
9) Benzaldehyde	7.29	77	47411	19.662	ng	95
10) Phenol	7.35	94	215086	41.615	ng	94
11) bis(2-Chloroethyl)ether	7.58	93	172277	45.750	ng	92
12) 1,3-Dichlorobenzene	8.06	146	191116	43.483	ng	96
13) 1,4-Dichlorobenzene	8.20	146	192617	42.924	ng	95
14) 1,2-Dichlorobenzene	8.52	146	189552	43.794	ng	97
15) Benzyl Alcohol	8.40	79	174336	51.603	ng	94
16) 2,2'-oxybis(1-Chloropropan	8.69	45	289294	45.238	ng	94
17) 2-Methylphenol	8.60	107	162115	47.218	ng	95
18) Hexachloroethane	9.26	117	67015	43.823	ng	97
19) n-Nitroso-di-n-propylamine	8.97	70	146742	45.017	ng	# 95
20) 3+4-Methylphenols	8.93	107	225747	46.664	ng	95
22) Acetophenone	8.99	105	276178	45.782	ng	# 95
24) Nitrobenzene	9.37	77	207202	46.784	ng	92
25) Isophorone	9.90	82	415952	50.583	ng	# 95
26) 2-Nitrophenol	10.09	139	107788	50.921	ng	# 88
27) 2,4-Dimethylphenol	10.14	122	185880	48.535	ng	93
28) bis(2-Chloroethoxy)methane	10.38	93	249419	49.465	ng	95
29) 2,4-Dichlorophenol	10.63	162	210382	51.514	ng	91
30) 1,2,4-Trichlorobenzene	10.84	180	206818	46.875	ng	99
31) Naphthalene	11.04	128	572429	45.886	ng	98
32) Benzoic acid	10.26	122	105030	32.753	ng	87
33) 4-Chloroaniline	11.14	127	39133	7.457	ng	# 93
34) Hexachlorobutadiene	11.32	225	127573	46.504	ng	97
35) Caprolactam	11.91	113	68407m	50.400	ng	
36) 4-Chloro-3-methylphenol	12.25	107	234789	52.271	ng	# 89
37) 2-Methylnaphthalene	12.63	142	450564	49.555	ng	94
39) 1,2,4,5-Tetrachlorobenzene	12.99	216	252543	39.442	ng	98
40) Hexachlorocyclopentadiene	12.97	237	292494	118.223	ng	94

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42) 2,4,6-Trichlorophenol	13.22	196	194714	48.442	nq	99
43) 2,4,5-Trichlorophenol	13.30	196	215833	52.286	nq	94
45) 1,1'-Biphenyl	13.62	154	620135	41.139	nq	96
46) 2-Chloronaphthalene	13.67	162	481142	43.759	nq	97
47) 2-Nitroaniline	13.86	65	168761	55.880	nq	# 86
48) Acenaphthylene	14.51	152	790722	45.951	nq	98
49) Dimethylphthalate	14.23	163	704802	51.508	nq	100
50) 2,6-Dinitrotoluene	14.35	165	159739	55.689	nq	# 82
51) Acenaphthene	14.85	154	503756	44.994	nq	99
52) 3-Nitroaniline	14.68	138	94624	30.571	nq	# 79
53) 2,4-Dinitrophenol	14.89	184	107861	69.257	nq	# 52
54) Dibenzofuran	15.18	168	808248	50.568	nq	99
55) 4-Nitrophenol	14.99	139	266398	104.396	nq	# 84
56) 2,4-Dinitrotoluene	15.14	165	236244	60.841	nq	# 77
57) Fluorene	15.83	166	655490	49.644	nq	98
58) 2,3,4,6-Tetrachlorophenol	15.40	232	215676	62.625	nq	96
59) Diethylphthalate	15.59	149	735690	53.949	nq	99
60) 4-Chlorophenyl-phenylether	15.81	204	365732	51.952	nq	96
61) 4-Nitroaniline	15.84	138	180410	56.763	nq	88
62) Azobenzene	16.11	77	630447	54.825	nq	97
64) 4,6-Dinitro-2-methylphenol	15.90	198	106425	41.813	nq	85
65) n-Nitrosodiphenylamine	16.03	169	630772	46.236	nq	98
66) 4-Bromophenyl-phenylether	16.71	248	256437	49.071	nq	99
67) Hexachlorobenzene	16.83	284	270398	45.679	nq	98
68) Atrazine	16.97	200	265712	54.585	nq	99
69) Pentachlorophenol	17.17	266	348810	102.577	nq	99
70) Phenanthrene	17.56	178	1144181	48.632	nq	97
71) Anthracene	17.66	178	1170364	49.541	nq	97
72) Carbazole	17.92	167	1105999	48.735	nq	98
73) Di-n-butylphthalate	18.47	149	1298649	49.755	nq	99
74) Fluoranthene	19.56	202	1447878	52.615	nq	96
76) Benzidine	19.73	184	234888	15.734	nq	98
77) Pyrene	19.92	202	1467222	48.917	nq	94
79) Butylbenzylphthalate	20.79	149	619291	48.463	nq	92
80) Benzo(a)anthracene	21.78	228	1474439	50.790	nq	98
81) 3,3'-Dichlorobenzidine	21.68	252	390272	32.571	nq	98
82) Chrysene	21.84	228	1377949	49.564	nq	95
83) Bis(2-ethylhexyl)phthalate	21.67	149	862232	47.015	nq	100
84) Di-n-octyl phthalate	22.91	149	1492165	46.685	nq	97
85) Indeno(1,2,3-cd)pyrene	28.89	276	1834934	53.648	nq	# 93
87) Benzo(b)fluoranthene	24.05	252	1562869	51.606	nq	97
88) Benzo(k)fluoranthene	24.12	252	1471041	48.756	nq	97
89) Benzo(a)pyrene	24.95	252	1489102	51.147	nq	100
90) Dibenzo(a,h)anthracene	28.96	278	1522916	51.667	nq	99
91) Benzo(g,h,i)perylene	30.09	276	1505435	54.969	nq	96

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

