

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG110217\
 Data File : BG030673.D
 Acq On : 2 Nov 2017 21:54
 Operator : SJ/JU
 Sample : PB103830BS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB103830BS

Manual Integrations
 APPROVED

Sohil
 11/3/2017 1:24:55 PM

Quant Time: Nov 03 02:11:31 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG101617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 01 17:41:24 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.15	152	62382	20.00	ng	0.00
21) Naphthalene-d8	10.97	136	280841	20.00	ng	0.00
38) Acenaphthene-d10	14.77	164	199990	20.00	ng	0.00
63) Phenanthrene-d10	17.51	188	507732	20.00	ng	0.00
75) Chrysene-d12	21.78	240	554077	20.00	ng	-0.01
86) Perylene-d12	25.08	264	574204	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.71	112	382743	102.50	ng	0.00
7) Phenol-d6	7.31	99	517077	98.65	ng	0.00
23) Nitrobenzene-d5	9.32	82	285596	64.62	ng	-0.01
41) 2,4,6-Tribromophenol	16.26	330	284068	110.02	ng	0.00
44) 2-Fluorobiphenyl	13.39	172	855608	62.75	ng	0.00
78) Terphenyl-d14	20.10	244	1509946	62.00	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.56	88	46030	30.381	ng	# 83
3) Pyridine	3.97	79	125582	28.463	ng	# 87
4) n-Nitrosodimethylamine	3.89	42	56731	27.507	ng	# 95
6) Aniline	7.47	93	110391	17.008	ng	# 93
8) 2-Chlorophenol	7.72	128	131607	30.747	ng	# 92
9) Benzaldehyde	7.28	77	57696	21.884	ng	# 95
10) Phenol	7.34	94	176912	31.307	ng	# 94
11) bis(2-Chloroethyl)ether	7.57	93	122850	29.839	ng	# 91
12) 1,3-Dichlorobenzene	8.04	146	136026	28.306	ng	# 97
13) 1,4-Dichlorobenzene	8.19	146	138371	28.203	ng	# 95
14) 1,2-Dichlorobenzene	8.51	146	135021	28.532	ng	# 95
15) Benzyl Alcohol	8.39	79	115831	31.358	ng	# 92
16) 2,2'-oxybis(1-Chloropropan	8.68	45	189566	27.112	ng	# 95
17) 2-Methylphenol	8.59	107	118576	31.588	ng	# 94
18) Hexachloroethane	9.25	117	49111	29.373	ng	# 97
19) n-Nitroso-di-n-propylamine	8.95	70	99745	27.987	ng	# 96
20) 3+4-Methylphenols	8.92	107	167889	31.741	ng	# 94
22) Acetophenone	8.98	105	191990	28.367	ng	# 94
24) Nitrobenzene	9.36	77	146036	29.389	ng	# 91
25) Isophorone	9.88	82	279908	30.339	ng	# 94
26) 2-Nitrophenol	10.07	139	73841	31.092	ng	# 86
27) 2,4-Dimethylphenol	10.13	122	138874	32.320	ng	# 89
28) bis(2-Chloroethoxy)methane	10.36	93	178165	31.493	ng	# 98
29) 2,4-Dichlorophenol	10.61	162	155402	33.915	ng	# 92
30) 1,2,4-Trichlorobenzene	10.83	180	151657	30.636	ng	# 98
31) Naphthalene	11.02	128	416710	29.772	ng	# 99
32) Benzoic acid	10.25	122	82413	22.906	ng	# 87
33) 4-Chloroaniline	11.13	127	89625	15.223	ng	# 95
34) Hexachlorobutadiene	11.30	225	92993	30.214	ng	# 97
35) Caprolactam	11.88	113	56127m	36.857	ng	# 94
36) 4-Chloro-3-methylphenol	12.23	107	163726	32.488	ng	# 88
37) 2-Methylnaphthalene	12.61	142	336990	33.035	ng	# 93
39) 1,2,4,5-Tetrachlorobenzene	12.98	216	191636	26.246	ng	# 98
40) Hexachlorocyclopentadiene	12.95	237	201720	73.339	ng	# 92

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.21	196	140311	30.611	ng	99
43) 2,4,5-Trichlorophenol	13.29	196	154117	32.740	ng	93
45) 1,1'-Biphenyl	13.61	154	467786	27.213	ng	100
46) 2-Chloronaphthalene	13.65	162	364363	29.060	ng	97
47) 2-Nitroaniline	13.85	65	109562	31.813	ng	# 81
48) Acenaphthylene	14.49	152	590315	30.082	ng	99
49) Dimethylphthalate	14.22	163	499919	32.038	ng	99
50) 2,6-Dinitrotoluene	14.34	165	110959	33.922	ng	# 77
51) Acenaphthene	14.83	154	365626	28.637	ng	99
52) 3-Nitroaniline	14.67	138	71239	20.183	ng	# 83
53) 2,4-Dinitrophenol	14.88	184	113070	64.201	ng	# 51
54) Dibenzofuran	15.17	168	606287	33.264	ng	99
55) 4-Nitrophenol	14.97	139	191158	65.691	ng	# 78
56) 2,4-Dinitrotoluene	15.13	165	162318	36.657	ng	# 75
57) Fluorene	15.81	166	473463	31.445	ng	99
58) 2,3,4,6-Tetrachlorophenol	15.39	232	149164	37.981	ng	# 94
59) Diethylphthalate	15.57	149	497088	31.966	ng	99
60) 4-Chlorophenyl-phenylether	15.80	204	267833	33.363	ng	94
61) 4-Nitroaniline	15.83	138	118929	32.814	ng	83
62) Azobenzene	16.09	77	434679	33.148	ng	96
64) 4,6-Dinitro-2-methylphenol	15.89	198	81804	30.381	ng	80
65) n-Nitrosodiphenylamine	16.01	169	439509	28.897	ng	99
66) 4-Bromophenyl-phenylether	16.69	248	180906	31.051	ng	98
67) Hexachlorobenzene	16.82	284	193288	29.289	ng	95
68) Atrazine	16.95	200	177188	32.650	ng	99
69) Pentachlorophenol	17.16	266	225253	59.417	ng	99
70) Phenanthrene	17.55	178	814952	31.070	ng	98
71) Anthracene	17.64	178	835159	31.709	ng	99
72) Carbazole	17.91	167	773773	30.583	ng	99
73) Di-n-butylphthalate	18.45	149	894261	30.732	ng	99
74) Fluoranthene	19.55	202	1027918	33.506	ng	97
76) Benzidine	19.72	184	356333	21.300	ng	96
77) Pyrene	19.91	202	1051859	31.295	ng	96
79) Butylbenzylphthalate	20.78	149	424271	29.628	ng	86
80) Benzo(a)anthracene	21.76	228	1048688	32.236	ng	99
81) 3,3'-Dichlorobenzidine	21.67	252	241338	17.974	ng	98
82) Chrysene	21.83	228	985978	31.648	ng	98
83) Bis(2-ethylhexyl)phthalate	21.64	149	589119	28.666	ng	100
84) Di-n-octyl phthalate	22.88	149	1016699	28.386	ng	# 95
85) Indeno(1,2,3-cd)pyrene	28.86	276	1249049	32.588	ng	# 87
87) Benzo(b)fluoranthene	24.03	252	1067287	32.466	ng	99
88) Benzo(k)fluoranthene	24.10	252	1038490	31.709	ng	98
89) Benzo(a)pyrene	24.93	252	1026662	32.486	ng	99
90) Dibenzo(a,h)anthracene	28.92	278	1047794	32.749	ng	96
91) Benzo(g,h,i)perylene	30.05	276	1047666	35.242	ng	98

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

