

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080218\
 Data File : BG036053.D
 Acq On : 2 Aug 2018 10:04
 Operator : JU/SJ
 Sample : PB111656BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB111656BS

Manual Integrations
 APPROVED

Sohil
 8/3/2018 12:58:25 PM

Quant Time: Aug 02 11:06:23 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 31 12:30:06 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	29525	20.00	ng	0.00
21) Naphthalene-d8	10.82	136	104851	20.00	ng	-0.01
38) Acenaphthene-d10	14.65	164	82512	20.00	ng	0.00
63) Phenanthrene-d10	17.39	188	242223	20.00	ng	0.00
75) Chrysene-d12	21.65	240	265885	20.00	ng	0.00
86) Perylene-d12	24.84	264	261012	20.00	ng	-0.02

System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	234281	131.74	ng	0.00
7) Phenol-d6	7.19	99	339352	127.90	ng	0.00
23) Nitrobenzene-d5	9.18	82	223321	88.98	ng	0.00
41) 2,4,6-Tribromophenol	16.14	330	179303	164.87	ng	0.00
44) 2-Fluorobiphenyl	13.27	172	545148	88.52	ng	-0.01
78) Terphenyl-d14	20.00	244	1110233	92.04	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.51	88	26752	29.556	ng	# 74
3) Pyridine	3.91	79	76192	32.245	ng	# 68
4) n-Nitrosodimethylamine	3.83	42	33961	33.473	ng	# 79
6) Aniline	7.34	93	97960	28.485	ng	95
8) 2-Chlorophenol	7.59	128	77405	42.796	ng	94
9) Benzaldehyde	7.16	77	43921	26.978	ng	93
10) Phenol	7.22	94	112963	42.141	ng	97
11) bis(2-Chloroethyl)ether	7.44	93	75351	35.893	ng	87
12) 1,3-Dichlorobenzene	7.91	146	85237	39.025	ng	90
13) 1,4-Dichlorobenzene	8.05	146	86347	38.909	ng	99
14) 1,2-Dichlorobenzene	8.37	146	84031	39.416	ng	95
15) Benzyl Alcohol	8.26	79	88669	36.801	ng	92
16) 2,2'-oxybis(1-Chloropropan	8.54	45	84080	47.772	ng	79
17) 2-Methylphenol	8.47	107	74992	41.147	ng	# 92
18) Hexachloroethane	9.10	117	33635	39.640	ng	91
19) n-Nitroso-di-n-propylamine	8.82	70	72257	32.340	ng	88
20) 3+4-Methylphenols	8.79	107	104140	41.188	ng	93
22) Acetophenone	8.84	105	130320	39.969	ng	# 93
24) Nitrobenzene	9.22	77	109512	43.385	ng	95
25) Isophorone	9.74	82	212752	45.662	ng	98
26) 2-Nitrophenol	9.93	139	44333	49.453	ng	# 85
27) 2,4-Dimethylphenol	9.99	122	81050	53.411	ng	84
28) bis(2-Chloroethoxy)methane	10.22	93	107590	41.907	ng	97
29) 2,4-Dichlorophenol	10.48	162	88976	51.365	ng	96
30) 1,2,4-Trichlorobenzene	10.69	180	96440	45.403	ng	98
31) Naphthalene	10.87	128	239586	43.866	ng	98
32) Benzoic acid	10.17	122	25904	25.357	ng	91
33) 4-Chloroaniline	10.99	127	62246	26.148	ng	98
34) Hexachlorobutadiene	11.15	225	76171	45.988	ng	98
35) Caprolactam	11.77	113	34012m	45.754	ng	
36) 4-Chloro-3-methylphenol	12.11	107	118782	51.157	ng	90
37) 2-Methylnaphthalene	12.47	142	192357	47.273	ng	# 92
39) 1,2,4,5-Tetrachlorobenzene	12.84	216	133440	42.848	ng	99
40) Hexachlorocyclopentadiene	12.82	237	164762	100.879	ng	90

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.08	196	87446	48.014	ng	98
43) 2,4,5-Trichlorophenol	13.17	196	101022	49.583	ng	98
45) 1,1'-Biphenyl	13.48	154	276033	43.150	ng	99
46) 2-Chloronaphthalene	13.52	162	217856	43.782	ng	99
47) 2-Nitroaniline	13.73	65	76966	43.752	ng	94
48) Acenaphthylene	14.37	152	368939	44.262	ng	97
49) Dimethylphthalate	14.10	163	319556	44.203	ng	99
50) 2,6-Dinitrotoluene	14.22	165	73122	49.670	ng	94
51) Acenaphthene	14.71	154	217836	41.953	ng	97
52) 3-Nitroaniline	14.56	138	47934	31.857	ng #	95
53) 2,4-Dinitrophenol	14.76	184	61131	80.689	ng #	61
54) Dibenzofuran	15.04	168	373566	45.238	ng	95
55) 4-Nitrophenol	14.87	139	90456	84.416	ng #	84
56) 2,4-Dinitrotoluene	15.01	165	111066	52.588	ng #	87
57) Fluorene	15.69	166	310746	44.898	ng	96
58) 2,3,4,6-Tetrachlorophenol	15.27	232	105436	53.974	ng	91
59) Diethylphthalate	15.46	149	331522	44.598	ng	98
60) 4-Chlorophenyl-phenylether	15.68	204	183058	45.000	ng	95
61) 4-Nitroaniline	15.72	138	70112	44.546	ng #	85
62) Azobenzene	15.97	77	324811	44.298	ng	93
64) 4,6-Dinitro-2-methylphenol	15.77	198	61675	45.740	ng	85
65) n-Nitrosodiphenylamine	15.90	169	280768	40.723	ng	96
66) 4-Bromophenyl-phenylether	16.58	248	124776	44.401	ng	98
67) Hexachlorobenzene	16.70	284	133820	46.241	ng	95
68) Atrazine	16.85	200	129546	45.636	ng	96
69) Pentachlorophenol	17.04	266	121190	68.752	ng	98
70) Phenanthrene	17.43	178	529756	43.830	ng	98
71) Anthracene	17.52	178	544891	45.276	ng	98
72) Carbazole	17.79	167	486374	42.555	ng	99
73) Di-n-butylphthalate	18.34	149	559385	41.417	ng #	98
74) Fluoranthene	19.44	202	696278	42.768	ng	96
76) Benzidine	19.62	184	254284	36.377	ng	98
77) Pyrene	19.80	202	696323	41.418	ng	98
79) Butylbenzylphthalate	20.68	149	257635	42.853	ng	96
80) Benzo(a)anthracene	21.63	228	697078	42.071	ng	99
81) 3,3'-Dichlorobenzidine	21.55	252	181088	31.344	ng #	98
82) Chrysene	21.70	228	671015	43.702	ng	99
83) Bis(2-ethylhexyl)phthalate	21.53	149	361064	44.175	ng	98
84) Di-n-octyl phthalate	22.73	149	622251	45.468	ng	96
85) Indeno(1,2,3-cd)pyrene	28.48	276	753889	44.348	ng #	53
87) Benzo(b)fluoranthene	23.83	252	686401	43.267	ng #	96
88) Benzo(k)fluoranthene	23.90	252	658198	42.438	ng #	97
89) Benzo(a)pyrene	24.70	252	659290	44.671	ng #	97
90) Dibenzo(a,h)anthracene	28.55	278	638364	44.057	ng #	94
91) Benzo(g,h,i)perylene	29.63	276	631630	44.548	ng #	94

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

