

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072018\
 Data File : BG035786.D
 Acq On : 20 Jul 2018 16:12
 Operator : JU/SJ
 Sample : PB111042BSD
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB111042BSD

Manual Integrations
 APPROVED

Sohil
 7/23/2018 2:18:57 PM

Quant Time: Jul 20 16:59:12 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 18 17:45:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.03	152	49172	20.00	ng	-0.02
21) Naphthalene-d8	10.83	136	178643	20.00	ng	0.00
38) Acenaphthene-d10	14.65	164	125050	20.00	ng	0.00
63) Phenanthrene-d10	17.39	188	338153	20.00	ng	0.00
75) Chrysene-d12	21.66	240	361578	20.00	ng	-0.02
86) Perylene-d12	24.86	264	350567	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.61	112	264084	89.16	ng	0.00
7) Phenol-d6	7.20	99	378677	85.69	ng	0.00
23) Nitrobenzene-d5	9.19	82	358950	83.94	ng	0.00
41) 2,4,6-Tribromophenol	16.14	330	155959	94.62	ng	0.00
44) 2-Fluorobiphenyl	13.28	172	827258	88.63	ng	0.00
78) Terphenyl-d14	20.00	244	1476790	90.03	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.52	88	45576	30.234	ng	# 74
3) Pyridine	3.91	79	132213	33.597	ng	# 72
4) n-Nitrosodimethylamine	3.83	42	53419	31.614	ng	# 74
6) Aniline	7.36	93	158552	27.682	ng	98
8) 2-Chlorophenol	7.60	128	126292	41.926	ng	96
9) Benzaldehyde	7.17	77	76307	28.144	ng	97
10) Phenol	7.23	94	186971	41.880	ng	97
11) bis(2-Chloroethyl)ether	7.45	93	127265	36.400	ng	92
12) 1,3-Dichlorobenzene	7.92	146	142334	39.129	ng	98
13) 1,4-Dichlorobenzene	8.06	146	139525	37.751	ng	98
14) 1,2-Dichlorobenzene	8.38	146	135713	38.223	ng	93
15) Benzyl Alcohol	8.27	79	144029	35.893	ng	88
16) 2,2'-oxybis(1-Chloropropan	8.55	45	143906	49.095	ng	76
17) 2-Methylphenol	8.47	107	123723	40.761	ng	# 91
18) Hexachloroethane	9.11	117	55272	39.112	ng	94
19) n-Nitroso-di-n-propylamine	8.83	70	117535	31.586	ng	# 83
20) 3+4-Methylphenols	8.80	107	168323	39.974	ng	93
22) Acetophenone	8.85	105	213918	38.507	ng	# 89
24) Nitrobenzene	9.23	77	173736	40.397	ng	90
25) Isophorone	9.75	82	326130	41.083	ng	100
26) 2-Nitrophenol	9.94	139	73788	48.310	ng	# 92
27) 2,4-Dimethylphenol	10.00	122	128709	49.782	ng	88
28) bis(2-Chloroethoxy)methane	10.23	93	174157	39.814	ng	95
29) 2,4-Dichlorophenol	10.48	162	142071	48.138	ng	93
30) 1,2,4-Trichlorobenzene	10.69	180	155293	42.911	ng	99
31) Naphthalene	10.88	128	381386	40.984	ng	98
32) Benzoic acid	10.17	122	61712	35.456	ng	98
33) 4-Chloroaniline	11.00	127	80889	19.944	ng	98
34) Hexachlorobutadiene	11.16	225	121541	43.069	ng	97
35) Caprolactam	11.77	113	48080m	37.962	ng	
36) 4-Chloro-3-methylphenol	12.11	107	170487	43.096	ng	97
37) 2-Methylnaphthalene	12.48	142	293007	42.263	ng	93
39) 1,2,4,5-Tetrachlorobenzene	12.85	216	203921	43.205	ng	99
40) Hexachlorocyclopentadiene	12.83	237	275816	111.428	ng	93

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.09	196	130543	47.295	ng	100
43) 2,4,5-Trichlorophenol	13.17	196	143909	46.606	ng	95
45) 1,1'-Biphenyl	13.49	154	409375	42.225	ng	99
46) 2-Chloronaphthalene	13.53	162	324787	43.068	ng	99
47) 2-Nitroaniline	13.73	65	110464	41.433	ng	91
48) Acenaphthylene	14.38	152	526829	41.705	ng	99
49) Dimethylphthalate	14.11	163	443072	40.440	ng	99
50) 2,6-Dinitrotoluene	14.23	165	102482	45.933	ng	91
51) Acenaphthene	14.72	154	309280	39.303	ng	99
52) 3-Nitroaniline	14.56	138	58548	25.675	ng	# 94
53) 2,4-Dinitrophenol	14.77	184	79701	70.334	ng	# 64
54) Dibenzofuran	15.05	168	519112	41.479	ng	94
55) 4-Nitrophenol	14.88	139	146457	90.184	ng	# 88
56) 2,4-Dinitrotoluene	15.02	165	148236	46.312	ng	# 82
57) Fluorene	15.70	166	436058	41.572	ng	96
58) 2,3,4,6-Tetrachlorophenol	15.28	232	138846	46.899	ng	92
59) Diethylphthalate	15.47	149	446507	39.634	ng	96
60) 4-Chlorophenyl-phenylether	15.69	204	247040	40.071	ng	96
61) 4-Nitroaniline	15.73	138	97130	40.720	ng	# 86
62) Azobenzene	15.98	77	443304	39.892	ng	96
64) 4,6-Dinitro-2-methylphenol	15.78	198	67406	35.809	ng	86
65) n-Nitrosodiphenylamine	15.90	169	391574	40.683	ng	97
66) 4-Bromophenyl-phenylether	16.58	248	170408	43.437	ng	96
67) Hexachlorobenzene	16.70	284	174651	43.229	ng	94
68) Atrazine	16.85	200	181706	45.851	ng	97
69) Pentachlorophenol	17.05	266	144285	58.633	ng	99
70) Phenanthrene	17.44	178	710071	42.083	ng	98
71) Anthracene	17.53	178	718018	42.736	ng	98
72) Carbazole	17.80	167	661324	41.447	ng	98
73) Di-n-butylphthalate	18.35	149	765768	40.613	ng	# 98
74) Fluoranthene	19.44	202	954192	41.983	ng	97
76) Benzidine	19.63	184	374264	39.371	ng	98
77) Pyrene	19.81	202	944080	41.293	ng	98
79) Butylbenzylphthalate	20.69	149	357010	43.666	ng	# 95
80) Benzo(a)anthracene	21.64	228	954276	42.351	ng	99
81) 3,3'-Dichlorobenzidine	21.56	252	228058	29.027	ng	# 98
82) Chrysene	21.71	228	901873	43.192	ng	98
83) Bis(2-ethylhexyl)phthalate	21.54	149	496332	44.654	ng	99
84) Di-n-octyl phthalate	22.74	149	851407	45.748	ng	# 93
85) Indeno(1,2,3-cd)pyrene	28.51	276	1025336	44.353	ng	# 86
87) Benzo(b)fluoranthene	23.84	252	931222	43.704	ng	# 96
88) Benzo(k)fluoranthene	23.92	252	891178	42.781	ng	# 97
89) Benzo(a)pyrene	24.71	252	889965	44.896	ng	# 98
90) Dibenzo(a,h)anthracene	28.57	278	861033	44.244	ng	97
91) Benzo(g,h,i)perylene	29.65	276	855776	44.938	ng	# 94

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

