

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110218\
 Data File : BG037763.D
 Acq On : 2 Nov 2018 20:38
 Operator : JU/SJ
 Sample : PB114420BS
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB114420BS

Manual Integrations
 APPROVED

Sohil
 11/5/2018 9:49:40 AM

Quant Time: Nov 03 01:46:06 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 18 14:06:42 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.10	152	50820	20.00	ng	0.00
21) Naphthalene-d8	10.92	136	231136	20.00	ng	0.00
39) Acenaphthene-d10	14.73	164	154242	20.00	ng	0.00
64) Phenanthrene-d10	17.48	188	349816	20.00	ng	0.00
76) Chrysene-d12	21.77	240	319030	20.00	ng	0.00
87) Perylene-d12	25.10	264	319277	20.00	ng	-0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	499204	164.23	ng	0.00
7) Phenol-d6	7.25	99	702262	152.78	ng	0.00
23) Nitrobenzene-d5	9.28	82	327875	79.46	ng	0.00
42) 2,4,6-Tribromophenol	16.23	330	267477	158.99	ng	0.00
45) 2-Fluorobiphenyl	13.36	172	810627	79.25	ng	0.00
79) Terphenyl-d14	20.08	244	1173087	78.57	ng	0.00

Target Compounds					Qvalue	
2) 1,4-Dioxane	3.52	88	60755	39.412	ng	97
3) Pyridine	3.93	79	138165	35.560	ng	96
4) n-Nitrosodimethylamine	3.85	42	58025	41.928	ng	# 97
6) Aniline	7.42	93	181920	32.443	ng	96
8) 2-Chlorophenol	7.66	128	155205	43.645	ng	97
9) Benzaldehyde	7.24	77	113897	39.613	ng	98
10) Phenol	7.27	94	205772	42.429	ng	100
11) bis(2-Chloroethyl)ether	7.52	93	140365	38.108	ng	96
12) 1,3-Dichlorobenzene	7.99	146	155254	39.217	ng	98
13) 1,4-Dichlorobenzene	8.14	146	159775	39.455	ng	98
14) 1,2-Dichlorobenzene	8.45	146	153520	39.411	ng	97
15) Benzyl Alcohol	8.33	79	136432	40.171	ng	95
16) 2,2'-oxybis(1-Chloropropan	8.62	45	264920	40.196	ng	98
17) 2-Methylphenol	8.53	107	138688	43.255	ng	97
18) Hexachloroethane	9.18	117	58005	42.016	ng	94
19) n-Nitroso-di-n-propylamine	8.90	70	117190	37.025	ng	96
20) 3+4-Methylphenols	8.86	107	192233	41.935	ng	99
22) Acetophenone	8.93	105	227877	39.311	ng	# 97
24) Nitrobenzene	9.32	77	175570	40.960	ng	95
25) Isophorone	9.83	82	331330	43.949	ng	97
26) 2-Nitrophenol	10.03	139	87247	45.183	ng	98
27) 2,4-Dimethylphenol	10.07	122	163188	47.333	ng	98
28) bis(2-Chloroethoxy)methane	10.31	93	203801	41.261	ng	97
29) 2,4-Dichlorophenol	10.56	162	166576	43.394	ng	98
30) 1,2,4-Trichlorobenzene	10.77	180	171145	40.998	ng	93
31) Naphthalene	10.97	128	497883	43.855	ng	99
32) Benzoic acid	10.19	122	87876	39.243	ng	98
33) 4-Chloroaniline	11.08	127	132260	24.795	ng	97
34) Hexachlorobutadiene	11.24	225	103143	41.689	ng	98
35) Caprolactam	11.86	113	58465m	37.975	ng	
36) 4-Chloro-3-methylphenol	12.18	107	180817	41.767	ng	98
37) 2-Methylnaphthalene	12.57	142	381245	46.068	ng	98
38) 1-Methylnaphthalene	12.79	142	360369	44.839	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.93	216	203840	40.972	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.89	237	258900	108.942	ng	98
43) 2,4,6-Trichlorophenol	13.16	196	143883	43.289	ng	98
44) 2,4,5-Trichlorophenol	13.24	196	156278	43.184	ng	98
46) 1,1'-Biphenyl	13.57	154	488205	38.219	ng	99
47) 2-Chloronaphthalene	13.62	162	393193	40.686	ng	99
48) 2-Nitroaniline	13.82	65	124729	42.561	ng	96
49) Acenaphthylene	14.46	152	643662	44.277	ng	99
50) Dimethylphthalate	14.18	163	492469	41.271	ng	99
51) 2,6-Dinitrotoluene	14.31	165	112108	42.470	ng	95
52) Acenaphthene	14.80	154	374201	41.796	ng	99
53) 3-Nitroaniline	14.65	138	95785	32.686	ng	# 95
54) 2,4-Dinitrophenol	14.84	184	82380	73.995	ng	96
55) Dibenzofuran	15.13	168	604450	40.749	ng	100
56) 4-Nitrophenol	14.94	139	237353	109.425	ng	97
57) 2,4-Dinitrotoluene	15.10	165	157582	45.478	ng	93
58) Fluorene	15.78	166	486805	43.824	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.35	232	137102	44.248	ng	99
60) Diethylphthalate	15.54	149	501170	40.910	ng	98
61) 4-Chlorophenyl-phenylether	15.77	204	258985	40.000	ng	97
62) 4-Nitroaniline	15.81	138	119789	41.138	ng	97
63) Azobenzene	16.06	77	473342	40.497	ng	98
65) 4,6-Dinitro-2-methylphenol	15.86	198	77223	43.119	ng	99
66) n-Nitrosodiphenylamine	15.98	169	440635	41.875	ng	100
67) 4-Bromophenyl-phenylether	16.66	248	161074	41.929	ng	98
68) Hexachlorobenzene	16.77	284	163682	41.111	ng	99
69) Atrazine	16.92	200	164471	43.231	ng	99
70) Pentachlorophenol	17.12	266	183479	68.799	ng	95
71) Phenanthrene	17.52	178	772393	43.445	ng	99
72) Anthracene	17.62	178	791413	45.360	ng	99
73) Carbazole	17.89	167	709518	40.654	ng	100
74) Di-n-butylphthalate	18.42	149	837781	43.152	ng	99
75) Fluoranthene	19.53	202	900583	44.662	ng	100
77) Benzidine	19.71	184	427645	39.422	ng	99
78) Pyrene	19.89	202	889341	42.643	ng	100
80) Butylbenzylphthalate	20.76	149	377753	44.258	ng	97
81) Benzo(a)anthracene	21.75	228	843474	44.698	ng	99
82) 3,3'-Dichlorobenzidine	21.66	252	228825	31.285	ng	97
83) Chrysene	21.81	228	791766	43.635	ng	100
84) Bis(2-ethylhexyl)phthalate	21.62	149	533858	44.622	ng	99
85) Di-n-octyl phthalate	22.86	149	902416	46.255	ng	98
86) Indeno(1,2,3-cd)pyrene	28.93	276	757726	38.934	ng	98
88) Benzo(b)fluoranthene	24.03	252	806127	46.054	ng	99
89) Benzo(k)fluoranthene	24.10	252	802397	46.789	ng	98
90) Benzo(a)pyrene	24.94	252	778543	46.808	ng	99
91) Dibenzo(a,h)anthracene	28.99	278	572549	37.318	ng	98
92) Benzo(g,h,i)perylene	30.13	276	613138	39.501	ng	98

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

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