

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG092718\
 Data File : BG037031.D
 Acq On : 28 Sep 2018 1:44
 Operator : JU/SJ
 Sample : PB113286BS
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB113286BS

Manual Integrations
 APPROVED

Sohil
 9/28/2018 4:24:37 PM

Quant Time: Sep 28 04:48:24 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 24 10:41:23 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.04	152	39867	20.00	ng	0.00
21) Naphthalene-d8	10.84	136	158628	20.00	ng	-0.01
38) Acenaphthene-d10	14.66	164	103206	20.00	ng	-0.01
63) Phenanthrene-d10	17.41	188	273360	20.00	ng	0.00
75) Chrysene-d12	21.68	240	286200	20.00	ng	-0.01
86) Perylene-d12	24.88	264	293656	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.62	112	311096	149.65	ng	0.00
7) Phenol-d6	7.21	99	443241	137.81	ng	0.00
23) Nitrobenzene-d5	9.21	82	287356	97.01	ng	0.00
41) 2,4,6-Tribromophenol	16.16	330	187781	152.07	ng	-0.01
44) 2-Fluorobiphenyl	13.29	172	618180	89.21	ng	-0.01
78) Terphenyl-d14	20.02	244	1128517	103.68	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.52	88	43314	45.535	ng	98
3) Pyridine	3.91	79	119006	43.329	ng	97
4) n-Nitrosodimethylamine	3.83	42	68757	56.324	ng	# 90
6) Aniline	7.37	93	127091	30.453	ng	99
8) 2-Chlorophenol	7.61	128	118555	49.117	ng	93
9) Benzaldehyde	7.18	77	74449	35.031	ng	99
10) Phenol	7.23	94	164489	49.554	ng	92
11) bis(2-Chloroethyl)ether	7.46	93	116325	37.885	ng	96
12) 1,3-Dichlorobenzene	7.93	146	125429	42.386	ng	99
13) 1,4-Dichlorobenzene	8.08	146	123471	40.772	ng	98
14) 1,2-Dichlorobenzene	8.39	146	123689	42.147	ng	99
15) Benzyl Alcohol	8.28	79	112772	43.212	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.56	45	255640	43.367	ng	97
17) 2-Methylphenol	8.48	107	106036	45.860	ng	96
18) Hexachloroethane	9.12	117	50023	44.887	ng	97
19) n-Nitroso-di-n-propylamine	8.84	70	103813	38.800	ng	99
20) 3+4-Methylphenols	8.81	107	149292	46.413	ng	96
22) Acetophenone	8.86	105	187152	50.206	ng	# 96
24) Nitrobenzene	9.25	77	152913	48.339	ng	99
25) Isophorone	9.76	82	290792	53.726	ng	99
26) 2-Nitrophenol	9.96	139	67178	53.278	ng	99
27) 2,4-Dimethylphenol	10.01	122	114336	58.213	ng	98
28) bis(2-Chloroethoxy)methane	10.25	93	165723	49.620	ng	99
29) 2,4-Dichlorophenol	10.49	162	122043	53.602	ng	96
30) 1,2,4-Trichlorobenzene	10.70	180	130878	45.933	ng	95
31) Naphthalene	10.90	128	380374	50.396	ng	98
32) Benzoic acid	10.15	122	48973m	34.418	ng	
33) 4-Chloroaniline	11.01	127	73943	21.547	ng	97
34) Hexachlorobutadiene	11.17	225	88121	44.721	ng	98
35) Caprolactam	11.78	113	49408m	52.724	ng	
36) 4-Chloro-3-methylphenol	12.13	107	146045	53.758	ng	99
37) 2-Methylnaphthalene	12.50	142	287440	50.003	ng	98
39) 1,2,4,5-Tetrachlorobenzene	12.87	216	163066	45.301	ng	98
40) Hexachlorocyclopentadiene	12.84	237	203426	109.882	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.10	196	104924	52.491	nq	97
43) 2,4,5-Trichlorophenol	13.18	196	119602	56.113	nq	98
45) 1,1'-Biphenyl	13.50	154	382270	48.363	nq	99
46) 2-Chloronaphthalene	13.55	162	287206	46.205	nq	98
47) 2-Nitroaniline	13.75	65	117264	54.542	nq	91
48) Acenaphthylene	14.39	152	505367	51.884	nq	98
49) Dimethylphthalate	14.12	163	411218	49.500	nq	100
50) 2,6-Dinitrotoluene	14.24	165	90866	50.132	nq	98
51) Acenaphthene	14.73	154	285201	48.447	nq	98
52) 3-Nitroaniline	14.58	138	60545	31.700	nq	97
53) 2,4-Dinitrophenol	14.78	184	49017	59.837	nq	92
54) Dibenzofuran	15.06	168	484709	48.678	nq	99
55) 4-Nitrophenol	14.89	139	143541	117.642	nq	96
56) 2,4-Dinitrotoluene	15.03	165	131383	51.947	nq	95
57) Fluorene	15.72	166	392199	52.697	nq	95
58) 2,3,4,6-Tetrachlorophenol	15.29	232	120686	55.816	nq	98
59) Diethylphthalate	15.48	149	418975	50.004	nq	98
60) 4-Chlorophenyl-phenylether	15.70	204	218790	46.420	nq	98
61) 4-Nitroaniline	15.74	138	104028	51.781	nq	98
62) Azobenzene	16.00	77	429922	53.401	nq	99
64) 4,6-Dinitro-2-methylphenol	15.79	198	60505	42.336	nq	92
65) n-Nitrosodiphenylamine	15.92	169	356204	47.200	nq	99
66) 4-Bromophenyl-phenylether	16.60	248	140983	45.342	nq	99
67) Hexachlorobenzene	16.72	284	144129	45.007	nq	100
68) Atrazine	16.87	200	158004	51.708	nq	100
69) Pentachlorophenol	17.06	266	167966	83.307	nq	97
70) Phenanthrene	17.45	178	680370	51.649	nq	99
71) Anthracene	17.55	178	698971	53.661	nq	99
72) Carbazole	17.82	167	633975	49.919	nq	99
73) Di-n-butylphthalate	18.36	149	743375	52.707	nq	100
74) Fluoranthene	19.46	202	858681	54.153	nq	99
76) Benzidine	19.65	184	339825	39.278	nq	99
77) Pyrene	19.82	202	845371	49.223	nq	100
79) Butylbenzylphthalate	20.70	149	356799	52.121	nq	99
80) Benzo(a)anthracene	21.66	228	847878	50.750	nq	100
81) 3,3'-Dichlorobenzidine	21.58	252	195610	30.760	nq	98
82) Chrysene	21.73	228	804129	49.815	nq	99
83) Bis(2-ethylhexyl)phthalate	21.56	149	494536	51.713	nq	99
84) Di-n-octyl phthalate	22.77	149	835812	53.083	nq	100
85) Indeno(1,2,3-cd)pyrene	28.54	276	942126	54.351	nq	# 94
87) Benzo(b)fluoranthene	23.87	252	820756	52.446	nq	99
88) Benzo(k)fluoranthene	23.93	252	822036	52.357	nq	99
89) Benzo(a)pyrene	24.73	252	809704	53.517	nq	99
90) Dibenzo(a,h)anthracene	28.60	278	758690	53.505	nq	99
91) Benzo(g,h,i)perylene	29.67	276	781951	54.947	nq	98

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

