

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG100418\
 Data File : BG037137.D
 Acq On : 3 Oct 2018 17:23
 Operator : JU/SJ
 Sample : J5212-05MS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 CG-SOIL-2MS

Manual Integrations
 APPROVED

Sohil
 10/4/2018 12:59:18 PM

Quant Time: Oct 03 18:11:16 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	62077	20.00	ng	0.00
21) Naphthalene-d8	10.84	136	247898	20.00	ng	-0.01
38) Acenaphthene-d10	14.66	164	149856	20.00	ng	-0.01
63) Phenanthrene-d10	17.41	188	356281	20.00	ng	0.00
75) Chrysene-d12	21.68	240	352352	20.00	ng	-0.01
86) Perylene-d12	24.88	264	359985	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.62	112	466383	144.08	ng	0.00
7) Phenol-d6	7.21	99	627733	125.34	ng	0.00
23) Nitrobenzene-d5	9.20	82	423557	91.50	ng	0.00
41) 2,4,6-Tribromophenol	16.15	330	232792	129.84	ng	-0.01
44) 2-Fluorobiphenyl	13.29	172	851568	84.64	ng	-0.01
78) Terphenyl-d14	20.01	244	1245148	92.92	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.53	88	65281	44.074	ng	96
3) Pyridine	3.92	79	175264	40.981	ng	98
4) n-Nitrosodimethylamine	3.85	42	99776	52.491	ng	# 99
6) Aniline	7.37	93	188186	28.959	ng	98
8) 2-Chlorophenol	7.61	128	186128	49.523	ng	95
9) Benzaldehyde	7.18	77	95708	28.921	ng	97
10) Phenol	7.24	94	261832	50.658	ng	97
11) bis(2-Chloroethyl)ether	7.46	93	194801	40.745	ng	99
12) 1,3-Dichlorobenzene	7.93	146	188329	40.872	ng	98
13) 1,4-Dichlorobenzene	8.08	146	195974	41.560	ng	98
14) 1,2-Dichlorobenzene	8.39	146	183981	40.262	ng	96
15) Benzyl Alcohol	8.28	79	171300	42.155	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.56	45	392492	42.761	ng	97
17) 2-Methylphenol	8.48	107	166895	46.356	ng	98
18) Hexachloroethane	9.12	117	76157	43.888	ng	97
19) n-Nitroso-di-n-propylamine	8.85	70	162503	39.006	ng	98
20) 3+4-Methylphenols	8.81	107	229478	45.817	ng	97
22) Acetophenone	8.86	105	283506	48.666	ng	# 97
24) Nitrobenzene	9.24	77	241616	48.875	ng	100
25) Isophorone	9.76	82	450531	53.263	ng	99
26) 2-Nitrophenol	9.96	139	105968	53.778	ng	99
27) 2,4-Dimethylphenol	10.01	122	175354	57.129	ng	99
28) bis(2-Chloroethoxy)methane	10.24	93	266733	51.105	ng	98
29) 2,4-Dichlorophenol	10.49	162	189544	53.270	ng	96
30) 1,2,4-Trichlorobenzene	10.70	180	198347	44.544	ng	98
31) Naphthalene	10.89	128	578216	49.021	ng	99
32) Benzoic acid	10.01	122	175354	71.345	ng	# 14
33) 4-Chloroaniline	11.00	127	97911	18.257	ng	99
34) Hexachlorobutadiene	11.17	225	134318	43.619	ng	96
35) Caprolactam	11.78	113	63783m	43.553	ng	
36) 4-Chloro-3-methylphenol	12.12	107	202988	47.812	ng	99
37) 2-Methylnaphthalene	12.49	142	431110	47.990	ng	98
39) 1,2,4,5-Tetrachlorobenzene	12.86	216	234614	44.888	ng	97
40) Hexachlorocyclopentadiene	12.83	237	279884	104.119	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.10	196	148268	51.084	ng	97
43) 2,4,5-Trichlorophenol	13.18	196	165702	53.541	ng	97
45) 1,1'-Biphenyl	13.50	154	550958	48.006	ng	98
46) 2-Chloronaphthalene	13.55	162	420809	46.624	ng	100
47) 2-Nitroaniline	13.75	65	158345	50.722	ng	96
48) Acenaphthylene	14.39	152	704656	49.823	ng	99
49) Dimethylphthalate	14.12	163	901065	74.700	ng	98
50) 2,6-Dinitrotoluene	14.24	165	123626	46.974	ng	99
51) Acenaphthene	14.73	154	399698	46.760	ng	99
52) 3-Nitroaniline	14.57	138	78802	28.415	ng	97
53) 2,4-Dinitrophenol	14.78	184	49247	43.760	ng	92
54) Dibenzofuran	15.06	168	655384	45.329	ng	98
55) 4-Nitrophenol	14.89	139	176196	99.452	ng	99
56) 2,4-Dinitrotoluene	15.03	165	177680	48.383	ng	94
57) Fluorene	15.71	166	523555	48.448	ng	97
58) 2,3,4,6-Tetrachlorophenol	15.29	232	154730	49.284	ng	97
59) Diethylphthalate	15.48	149	545453	44.834	ng	100
60) 4-Chlorophenyl-phenylether	15.70	204	290119	42.392	ng	99
61) 4-Nitroaniline	15.74	138	133881	45.895	ng	95
62) Azobenzene	16.00	77	565765	48.398	ng	99
64) 4,6-Dinitro-2-methylphenol	15.79	198	91692	49.226	ng	95
65) n-Nitrosodiphenylamine	15.92	169	472687	48.057	ng	98
66) 4-Bromophenyl-phenylether	16.59	248	182360	44.999	ng	98
67) Hexachlorobenzene	16.72	284	183849	44.048	ng	98
68) Atrazine	16.87	200	197638	49.625	ng	98
69) Pentachlorophenol	17.06	266	209881	79.869	ng	98
70) Phenanthrene	17.45	178	843340	49.120	ng	99
71) Anthracene	17.54	178	860963	50.714	ng	99
72) Carbazole	17.81	167	774732	46.805	ng	100
73) Di-n-butylphthalate	18.36	149	895467	48.714	ng	99
74) Fluoranthene	19.46	202	1010919	48.915	ng	99
76) Benzidine	19.64	184	320471	30.087	ng	99
77) Pyrene	19.82	202	994330	47.026	ng	99
79) Butylbenzylphthalate	20.70	149	417281	49.512	ng	96
80) Benzo(a)anthracene	21.66	228	991960	48.227	ng	100
81) 3,3'-Dichlorobenzidine	21.57	252	217267	27.751	ng	96
82) Chrysene	21.72	228	931887	46.892	ng	98
83) Bis(2-ethylhexyl)phthalate	21.55	149	585657	49.743	ng	99
84) Di-n-octyl phthalate	22.76	149	995912	51.376	ng	98
85) Indeno(1,2,3-cd)pyrene	28.53	276	1099856	51.538	ng	98
87) Benzo(b)fluoranthene	23.86	252	966636	50.386	ng	97
88) Benzo(k)fluoranthene	23.93	252	949291	49.321	ng	99
89) Benzo(a)pyrene	24.73	252	959392	51.727	ng	100
90) Dibenzo(a,h)anthracene	28.59	278	891095	51.264	ng	96
91) Benzo(g,h,i)perylene	29.66	276	915885	52.500	ng	97

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

