

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083018\
 Data File : BF108656.D
 Acq On : 30 Aug 2018 22:33
 Operator : JU/SJ
 Sample : J4699-01MSD
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CG-SOIL-1MSD

Manual Integrations
 APPROVED

Sohil
 8/31/2018 12:18:29 PM

Quant Time: Aug 31 03:14:38 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF083018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 30 15:06:55 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.00	152	81626	20.00	ng	0.00
21) Naphthalene-d8	8.28	136	310631	20.00	ng	0.00
38) Acenaphthene-d10	10.04	164	140094	20.00	ng	0.00
63) Phenanthrene-d10	11.54	188	279718	20.00	ng	0.00
75) Chrysene-d12	14.19	240	276039	20.00	ng	0.00
86) Perylene-d12	15.75	264	232327	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.65	112	333152	70.90	ng	0.00
7) Phenol-d6	6.64	99	509122	75.65	ng	0.00
23) Nitrobenzene-d5	7.57	82	366260	59.85	ng	0.00
41) 2,4,6-Tribromophenol	10.84	330	146748	79.35	ng	0.00
44) 2-Fluorobiphenyl	9.35	172	650789	62.46	ng	0.00
78) Terphenyl-d14	13.12	244	735597	51.78	ng	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.90	88	60364	27.643	ng	# 84
3) Pyridine	3.71	79	134857	26.730	ng	# 86
4) n-Nitrosodimethylamine	3.68	42	68832	27.350	ng	# 80
6) Aniline	6.70	93	194998	26.852	ng	# 81
8) 2-Chlorophenol	6.80	128	197066	34.302	ng	# 94
9) Benzaldehyde	6.58	77	104242	25.061	ng	# 98
10) Phenol	6.66	94	263309	33.552	ng	# 85
11) bis(2-Chloroethyl)ether	6.73	93	239864	34.711	ng	# 98
12) 1,3-Dichlorobenzene	6.94	146	212920	31.964	ng	# 99
13) 1,4-Dichlorobenzene	7.02	146	240973	33.141	ng	# 98
14) 1,2-Dichlorobenzene	7.17	146	221453	33.162	ng	# 97
15) Benzyl Alcohol	7.16	79	136913	28.282	ng	# 93
16) 2,2'-oxybis(1-Chloropropan	7.27	45	357175	33.692	ng	# 71
17) 2-Methylphenol	7.26	107	149243	33.412	ng	# 69
18) Hexachloroethane	7.50	117	76592	30.629	ng	# 98
19) n-Nitroso-di-n-propylamine	7.40	70	142941	31.717	ng	# 91
20) 3+4-Methylphenols	7.42	107	170521	30.469	ng	# 24
22) Acetophenone	7.41	105	292786	37.757	ng	# 89
24) Nitrobenzene	7.58	77	206455	33.271	ng	# 92
25) Isophorone	7.81	82	378574	37.239	ng	# 96
26) 2-Nitrophenol	7.90	139	93939	33.441	ng	# 92
27) 2,4-Dimethylphenol	7.93	122	160988	39.406	ng	# 97
28) bis(2-Chloroethoxy)methane	8.02	93	236573	36.838	ng	# 97
29) 2,4-Dichlorophenol	8.15	162	165273	34.100	ng	# 97
30) 1,2,4-Trichlorobenzene	8.22	180	187017	34.506	ng	# 97
31) Naphthalene	8.31	128	596056	40.286	ng	# 100
32) Benzoic acid	8.04	122	21756m	8.304	ng	#
33) 4-Chloroaniline	8.37	127	180059	27.568	ng	# 97
34) Hexachlorobutadiene	8.41	225	120999	34.001	ng	# 99
35) Caprolactam	8.72	113	39452	30.548	ng	# 88
36) 4-Chloro-3-methylphenol	8.84	107	167275	32.156	ng	# 96
37) 2-Methylnaphthalene	9.00	142	378456	37.806	ng	# 99
39) 1,2,4,5-Tetrachlorobenzene	9.16	216	192710	38.747	ng	# 97
40) Hexachlorocyclopentadiene	9.14	237	43805	26.145	ng	# 95

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.28	196	106706	36.230	ng	99
43) 2,4,5-Trichlorophenol	9.33	196	129687	37.760	ng	# 94
45) 1,1'-Biphenyl	9.46	154	459221	37.653	ng	96
46) 2-Chloronaphthalene	9.49	162	350820	36.743	ng	97
47) 2-Nitroaniline	9.60	65	113066	35.406	ng	84
48) Acenaphthylene	9.91	152	551998	39.489	ng	99
49) Dimethylphthalate	9.75	163	486394	44.090	ng	# 99
50) 2,6-Dinitrotoluene	9.82	165	87226	35.871	ng	# 88
51) Acenaphthene	10.08	154	301101	36.026	ng	99
52) 3-Nitroaniline	10.02	138	71976	27.361	ng	96
53) 2,4-Dinitrophenol	10.13	184	19162	23.175	ng	# 46
54) Dibenzofuran	10.25	168	474462	36.279	ng	95
55) 4-Nitrophenol	10.20	139	80014	49.601	ng	# 84
56) 2,4-Dinitrotoluene	10.24	165	112779	35.027	ng	# 92
57) Fluorene	10.60	166	375080	37.008	ng	98
58) 2,3,4,6-Tetrachlorophenol	10.37	232	113477	36.881	ng	# 78
59) Diethylphthalate	10.45	149	393216	35.379	ng	98
60) 4-Chlorophenyl-phenylether	10.58	204	193632	33.668	ng	94
61) 4-Nitroaniline	10.64	138	77205	30.330	ng	93
62) Azobenzene	10.74	77	380625	34.454	ng	92
64) 4,6-Dinitro-2-methylphenol	10.65	198	18230	14.688	ng	88
65) n-Nitrosodiphenylamine	10.70	169	339266	36.507	ng	95
66) 4-Bromophenyl-phenylether	11.07	248	118333	34.421	ng	# 88
67) Hexachlorobenzene	11.14	284	124403	33.589	ng	# 87
68) Atrazine	11.22	200	109389	41.055	ng	95
69) Pentachlorophenol	11.34	266	117748	59.829	ng	99
70) Phenanthrene	11.57	178	532069	37.685	ng	98
71) Anthracene	11.62	178	592153	39.916	ng	99
72) Carbazole	11.78	167	525447	36.630	ng	99
73) Di-n-butylphthalate	12.08	149	598619	36.272	ng	99
74) Fluoranthene	12.76	202	648526	40.807	ng	94
76) Benzidine	12.89	184	388392	49.985	ng	99
77) Pyrene	12.99	202	651360	31.390	ng	99
79) Butylbenzylphthalate	13.59	149	279022	31.978	ng	# 96
80) Benzo(a)anthracene	14.18	228	621288	36.920	ng	100
81) 3,3'-Dichlorobenzidine	14.15	252	231638	37.956	ng	# 94
82) Chrysene	14.22	228	590789	36.507	ng	99
83) Bis(2-ethylhexyl)phthalate	14.14	149	380818	33.955	ng	# 99
84) Di-n-octyl phthalate	14.76	149	672144	40.761	ng	96
85) Indeno(1,2,3-cd)pyrene	17.37	276	411985	36.613	ng	# 89
87) Benzo(b)fluoranthene	15.28	252	564994	39.589	ng	# 96
88) Benzo(k)fluoranthene	15.31	252	530005	37.489	ng	# 95
89) Benzo(a)pyrene	15.68	252	483125	37.459	ng	# 97
90) Dibenzo(a,h)anthracene	17.38	278	352608	34.231	ng	# 92
91) Benzo(g,h,i)perylene	17.87	276	322433	32.206	ng	# 88

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

