

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022318\
 Data File : BF103190.D
 Acq On : 23 Feb 2018 11:06
 Operator : SJ/JU
 Sample : J1161-03
 Misc : MDL-S-1ppm
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MDL-SOIL-03-QT1-2018

Manual Integrations
 APPROVED

Quant Time: Feb 23 12:41:53 2018
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 23 02:43:46 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)	
1) 1,4-Dichlorobenzene-d4	6.87	152	171432	20.00	ng	0.00	
21) Naphthalene-d8	8.15	136	759387	20.00	ng	0.00	
38) Acenaphthene-d10	9.91	164	353698	20.00	ng	0.00	
63) Phenanthrene-d10	11.40	188	627738	20.00	ng	0.00	
75) Chrysene-d12	14.04	240	473909	20.00	ng	0.00	
86) Perylene-d12	15.54	264	462119	20.00	ng	0.00	Sohil
System Monitoring Compounds							
5) 2-Fluorophenol	5.49	112	1286019	113.46	ng	0.00	
7) Phenol-d6	6.50	99	1594640	114.97	ng	0.00	2/26/2018
23) Nitrobenzene-d5	7.44	82	1148408	86.36	ng	0.00	9:54:28 AM
41) 2,4,6-Tribromophenol	10.70	330	361094	120.61	ng	0.00	
44) 2-Fluorobiphenyl	9.23	172	1751985	85.41	ng	0.00	
78) Terphenyl-d14	12.99	244	1440734	73.08	ng	0.00	
Target Compounds						Qvalue	
2) 1,4-Dioxane	2.66	88	5958	0.995	ng	#	89
3) Pyridine	3.45	79	112	0.007	ng	#	9
4) n-Nitrosodimethylamine	3.49	42	3880m	0.602	ng		
6) Aniline	6.54	93	20262	1.012	ng		94
8) 2-Chlorophenol	6.66	128	13298	1.129	ng		90
9) Benzaldehyde	6.43	77	7318	0.164	ng		91
10) Phenol	6.52	94	18242	1.133	ng		81
11) bis(2-Chloroethyl)ether	6.60	93	14719	1.204	ng		88
12) 1,3-Dichlorobenzene	6.81	146	15558	1.156	ng		94
13) 1,4-Dichlorobenzene	6.89	146	15199	1.127	ng		97
14) 1,2-Dichlorobenzene	7.04	146	14948m	1.202	ng		
15) Benzyl Alcohol	7.03	79	27103	2.593	ng	#	39
16) 2,2'-oxybis(1-Chloropropan	7.14	45	23437	1.117	ng		96
17) 2-Methylphenol	7.12	107	11448	1.070	ng	#	92
18) Hexachloroethane	7.38	117	5669	1.154	ng		89
19) n-Nitroso-di-n-propylamine	7.27	70	10727	1.243	ng		95
20) 3+4-Methylphenols	7.29	107	15643	1.199	ng	#	87
22) Acetophenone	7.27	105	21186	1.195	ng	#	94
24) Nitrobenzene	7.45	77	18974	1.296	ng		96
25) Isophorone	7.44	82	1148395	43.850	ng	#	64
26) 2-Nitrophenol	7.77	139	5542	0.804	ng		92
27) 2,4-Dimethylphenol	7.81	122	9775	0.928	ng		92
28) bis(2-Chloroethoxy)methane	7.90	93	16415	1.066	ng		99
29) 2,4-Dichlorophenol	8.03	162	9357	0.928	ng		94
30) 1,2,4-Trichlorobenzene	8.09	180	11324	1.054	ng		98
31) Naphthalene	8.17	128	42965	1.156	ng		97
32) Benzoic acid	7.92	122	2774m	7.218	ng		
33) 4-Chloroaniline	8.23	127	14943	0.931	ng		96
34) Hexachlorobutadiene	8.29	225	6269	1.121	ng		90
35) Caprolactam	8.56	113	2782	0.785	ng	#	80
36) 4-Chloro-3-methylphenol	8.71	107	11613	1.021	ng		93
37) 2-Methylnaphthalene	8.87	142	27943	1.163	ng		98
39) 1,2,4,5-Tetrachlorobenzene	9.03	216	10664	1.103	ng		99
42) 2,4,6-Trichlorophenol	9.15	196	6006	0.923	ng		98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)

43) 2,4,5-Trichlorophenol	9.20	196	7087	0.947	ng	# 89
45) 1,1'-Biphenyl	9.33	154	39663	1.338	ng	99
46) 2-Chloronaphthalene	9.36	162	26312	1.122	ng	97
47) 2-Nitroaniline	9.46	65	8235	0.948	ng	# 72
48) Acenaphthylene	9.77	152	40946	1.135	ng	98
49) Dimethylphthalate	9.62	163	31629	1.115	ng	99 Sohil
50) 2,6-Dinitrotoluene	9.69	165	5463	0.917	ng	# 82
51) Acenaphthene	9.94	154	24503	1.165	ng	95
52) 3-Nitroaniline	9.87	138	6434	0.852	ng	# 86
54) Dibenzofuran	10.12	168	35807	1.240	ng	99 2/26/2018 9:54:28 AM
55) 4-Nitrophenol	10.12	139	17340m	3.703	ng	
56) 2,4-Dinitrotoluene	10.10	165	6558	0.871	ng	# 84
57) Fluorene	10.46	166	28461	1.157	ng	# 99
58) 2,3,4,6-Tetrachlorophenol	10.24	232	4716	0.938	ng	# 96
59) Diethylphthalate	10.32	149	31730	1.169	ng	97
60) 4-Chlorophenyl-phenylether	10.45	204	12140	1.164	ng	94
61) 4-Nitroaniline	10.48	138	5690	0.754	ng	96
62) Azobenzene	10.61	77	31549	1.142	ng	96
64) 4,6-Dinitro-2-methylphenol	10.52	198	802	5.134	ng	# 46
65) n-Nitrosodiphenylamine	10.56	169	24548	1.123	ng	97
66) 4-Bromophenyl-phenylether	10.94	248	6727	1.029	ng	93
67) Hexachlorobenzene	11.01	284	7367	1.038	ng	94
68) Atrazine	11.09	200	6542	0.996	ng	95
69) Pentachlorophenol	11.22	266	1046	0.305	ng	# 83
70) Phenanthrene	11.42	178	41660	1.206	ng	97
71) Anthracene	11.47	178	40693	1.149	ng	98
72) Carbazole	11.64	167	37336	1.058	ng	100
73) Di-n-butylphthalate	11.95	149	49239	1.115	ng	96
74) Fluoranthene	12.62	202	38492	1.109	ng	97
76) Benzidine	12.74	184	9231	0.521	ng	# 96
77) Pyrene	12.84	202	42166	1.141	ng	97
79) Butylbenzylphthalate	13.45	149	18438	0.944	ng	92
80) Benzo(a)anthracene	14.03	228	32809	1.067	ng	97
81) 3,3'-Dichlorobenzidine	14.00	252	10258	0.935	ng	# 97
82) Chrysene	14.07	228	33040	1.107	ng	99
83) Bis(2-ethylhexyl)phthalate	14.00	149	28138	1.068	ng	# 97
84) Di-n-octyl phthalate	14.63	149	47832	1.124	ng	96
85) Indeno(1,2,3-cd)pyrene	17.06	276	27163	1.149	ng	96
87) Benzo(b)fluoranthene	15.10	252	30506m	1.004	ng	
88) Benzo(k)fluoranthene	15.13	252	29834	1.043	ng	99
89) Benzo(a)pyrene	15.47	252	27247	1.004	ng	# 96
90) Dibenzo(a,h)anthracene	17.06	278	21600	1.030	ng	# 89
91) Benzo(g,h,i)perylene	17.52	276	21432	1.046	ng	96

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

