

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050818\
 Data File : BF105158.D
 Acq On : 8 May 2018 21:39
 Operator : JU/SJ
 Sample : J2133-03
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MDL-SOIL-03-QT2-2018

Manual Integrations
 APPROVED

Sohil
 5/9/2018 6:51:01 PM

Quant Time: May 09 09:10:45 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 08 16:33:54 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	203314	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	817054	20.00	ng	0.00
38) Acenaphthene-d10	9.83	164	401968	20.00	ng	0.00
63) Phenanthrene-d10	11.31	188	788441	20.00	ng	0.00
75) Chrysene-d12	14.01	240	668297	20.00	ng	0.00
86) Perylene-d12	15.56	264	586327	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.40	112	1566928	126.97	ng	0.00
7) Phenol-d6	6.42	99	1879077	127.87	ng	0.00
23) Nitrobenzene-d5	7.36	82	1119035	92.82	ng	0.00
41) 2,4,6-Tribromophenol	10.62	330	603886	136.17	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	2007268	77.51	ng	0.00
78) Terphenyl-d14	12.91	244	2270751	78.93	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.52	88	12471	1.911	ng	87
3) Pyridine	3.35	79	31802	1.817	ng	# 78
4) n-Nitrosodimethylamine	3.19	42	14109	1.591	ng	# 77
6) Aniline	6.46	93	44017	2.051	ng	96
8) 2-Chlorophenol	6.57	128	29063	2.220	ng	92
9) Benzaldehyde	6.35	77	12758	1.133	ng	97
10) Phenol	6.43	94	32033	1.959	ng	98
11) bis(2-Chloroethyl)ether	6.53	93	30582	2.269	ng	91
12) 1,3-Dichlorobenzene	6.73	146	36741	2.328	ng	96
13) 1,4-Dichlorobenzene	6.81	146	36529	2.315	ng	97
14) 1,2-Dichlorobenzene	6.96	146	34126	2.345	ng	96
15) Benzyl Alcohol	6.93	79	25465	2.264	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.07	45	52701	2.104	ng	97
17) 2-Methylphenol	7.03	107	24153	2.137	ng	95
18) Hexachloroethane	7.30	117	11421	2.249	ng	96
19) n-Nitroso-di-n-propylamine	7.20	70	21013	2.160	ng	91
20) 3+4-Methylphenols	7.19	107	31526	2.373	ng	92
22) Acetophenone	7.20	105	47106	2.473	ng	# 97
24) Nitrobenzene	7.37	77	35235	2.578	ng	98
25) Isophorone	7.60	82	51137	1.897	ng	95
26) 2-Nitrophenol	7.69	139	9119	11.249	ng	# 94
27) 2,4-Dimethylphenol	7.72	122	18909	1.578	ng	96
28) bis(2-Chloroethoxy)methane	7.82	93	35155	2.113	ng	99
29) 2,4-Dichlorophenol	7.92	162	22303	1.990	ng	94
30) 1,2,4-Trichlorobenzene	8.01	180	29657	2.294	ng	94
31) Naphthalene	8.09	128	91734	2.332	ng	99
32) Benzoic acid	7.76	122	7467	10.176	ng	91
33) 4-Chloroaniline	8.14	127	33389	2.041	ng	95
34) Hexachlorobutadiene	8.21	225	17870	2.386	ng	96
35) Caprolactam	8.46	113	5569	1.636	ng	# 78
36) 4-Chloro-3-methylphenol	8.61	107	22056	1.890	ng	95
37) 2-Methylnaphthalene	8.78	142	61295	2.235	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.95	216	30960	2.420	ng	# 97
40) Hexachlorocyclopentadiene	8.93	237	11842	1.712	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.06	196	15132	1.939	nq	96
43) 2,4,5-Trichlorophenol	9.09	196	17036	2.089	nq	92
45) 1,1'-Biphenyl	9.25	154	88443	2.700	nq	97
46) 2-Chloronaphthalene	9.27	162	59571	2.434	nq	97
47) 2-Nitroaniline	9.36	65	11092	7.182	nq	86
48) Acenaphthylene	9.69	152	89448	2.366	nq	100
49) Dimethylphthalate	9.54	163	67094	2.466	nq	99
50) 2,6-Dinitrotoluene	9.60	165	10550	1.787	nq #	87
51) Acenaphthene	9.86	154	60703	2.491	nq	97
52) 3-Nitroaniline	9.77	138	11993	1.928	nq #	94
53) 2,4-Dinitrophenol	9.87	184	1353	10.193	nq #	26
54) Dibenzofuran	10.03	168	85473	2.398	nq	98
55) 4-Nitrophenol	9.92	139	7252	5.526	nq	94
56) 2,4-Dinitrotoluene	10.00	165	12401	5.069	nq	98
57) Fluorene	10.37	166	67335	2.576	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.14	232	13436	1.846	nq #	87
59) Diethylphthalate	10.25	149	63887	2.389	nq	98
60) 4-Chlorophenyl-phenylether	10.37	204	32428	2.665	nq	91
61) 4-Nitroaniline	10.37	138	11287	1.824	nq	91
62) Azobenzene	10.53	77	61032	2.188	nq	94
64) 4,6-Dinitro-2-methylphenol	10.41	198	3235	9.910	nq	97
65) n-Nitrosodiphenylamine	10.49	169	54414	2.214	nq	93
66) 4-Bromophenyl-phenylether	10.86	248	18596	2.211	nq #	88
67) Hexachlorobenzene	10.92	284	22506	2.323	nq #	83
68) Atrazine	11.00	200	10396	1.350	nq	97
69) Pentachlorophenol	11.11	266	9171	4.765	nq	95
70) Phenanthrene	11.33	178	101841	2.450	nq	97
71) Anthracene	11.39	178	98170	2.324	nq	98
72) Carbazole	11.54	167	88185	2.317	nq	97
73) Di-n-butylphthalate	11.88	149	82647	2.030	nq	98
74) Fluoranthene	12.52	202	97990	2.262	nq	97
76) Benzidine	12.65	184	10863m	0.435	nq	
77) Pyrene	12.75	202	104763	2.155	nq	97
79) Butylbenzylphthalate	13.40	149	21851	7.965	nq #	97
80) Benzo(a)anthracene	14.00	228	86382	2.131	nq	99
81) 3,3'-Dichlorobenzidine	13.96	252	19306	1.267	nq	88
82) Chrysene	14.03	228	93919	2.250	nq	98
83) Bis(2-ethylhexyl)phthalate	14.00	149	34519	1.718	nq	98
84) Di-n-octyl phthalate	14.69	149	38811	7.787	nq #	92
85) Indeno(1,2,3-cd)pyrene	17.00	276	53519	1.618	nq	95
87) Benzo(b)fluoranthene	15.13	252	71289	1.987	nq	96
88) Benzo(k)fluoranthene	15.16	252	74006	2.128	nq #	95
89) Benzo(a)pyrene	15.49	252	60561	1.856	nq	97
90) Dibenzo(a,h)anthracene	17.02	278	44071	1.739	nq	96
91) Benzo(g,h,i)perylene	17.43	276	43681	1.762	nq	96

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

