

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080621\
 Data File : BF124959.D
 Acq On : 6 Aug 2021 12:31
 Operator : JU/CG
 Sample : M2262-03
 Misc : MDL-MDL-SOIL 8ppm
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MDL-SOIL-03-QT2-2021

Manual Integrations
 APPROVED

mohammad
 8/9/2021 12:16:07 PM

Quant Time: Aug 06 12:54:51 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF080421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 06 11:27:24 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.828	152	7081	20.000	ng	0.00
21) Naphthalene-d8	8.110	136	26077	20.000	ng	# 0.00
39) Acenaphthene-d10	9.863	164	14507	20.000	ng	0.00
64) Phenanthrene-d10	11.351	188	26135	20.000	ng	0.00
76) Chrysene-d12	13.986	240	20292	20.000	ng	0.00
86) Perylene-d12	15.445	264	17167	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.445	112	44310	102.506	ng	0.00
7) Phenol-d6	6.457	99	55432	101.699	ng	0.00
23) Nitrobenzene-d5	7.392	82	33303	66.809	ng	0.00
42) 2,4,6-Tribromophenol	10.657	330	13514	104.270	ng	0.00
45) 2-Fluorobiphenyl	9.186	172	68119	68.708	ng	0.00
79) Terphenyl-d14	12.939	244	74757	62.144	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.587	88	1381	8.540	ng	# 100
3) Pyridine	3.393	79	2159m	5.745	ng	
4) n-Nitrosodimethylamine	3.275	42	1673	6.269	ng	88
6) Aniline	6.492	93	5393	9.367	ng	95
8) 2-Chlorophenol	6.610	128	3740	7.888	ng	87
9) Benzaldehyde	6.381	77	3083	10.397	ng	97
10) Phenol	6.469	94	4454	7.920	ng	97
11) bis(2-Chloroethyl)ether	6.563	93	3498	8.291	ng	99
12) 1,3-Dichlorobenzene	6.769	146	4274	8.282	ng	96
13) 1,4-Dichlorobenzene	6.845	146	4362	8.455	ng	92
14) 1,2-Dichlorobenzene	6.998	146	4188	8.662	ng	95
15) Benzyl Alcohol	6.969	79	3320	8.444	ng	93
16) 2,2'-oxybis(1-Chloropr...	7.104	45	5866	8.678	ng	93
17) 2-Methylphenol	7.075	107	2744	7.731	ng	93
18) Hexachloroethane	7.339	117	1601	7.926	ng	87
19) n-Nitroso-di-n-propyla...	7.233	70	2460	7.606	ng	96
20) 3+4-Methylphenols	7.228	107	3851	8.173	ng	91
22) Acetophenone	7.233	105	5371	8.410	ng	# 90
24) Nitrobenzene	7.410	77	3916	8.119	ng	95
25) Isophorone	7.645	82	7020	8.251	ng	# 92
26) 2-Nitrophenol	7.728	139	1815	7.410	ng	# 89
27) 2,4-Dimethylphenol	7.757	122	3059	8.810	ng	# 73
28) bis(2-Chloroethoxy)met...	7.857	93	4280	8.839	ng	# 95
29) 2,4-Dichlorophenol	7.963	162	2996	7.745	ng	95
30) 1,2,4-Trichlorobenzene	8.051	180	3487	8.043	ng	96
31) Naphthalene	8.133	128	11465	8.557	ng	96
32) Benzoic acid	7.857	122	752	10.528	ng	# 76
33) 4-Chloroaniline	8.180	127	4419	8.851	ng	96
34) Hexachlorobutadiene	8.251	225	2164	8.359	ng	# 80
35) Caprolactam	8.516	113	676	6.581	ng	# 87
36) 4-Chloro-3-methylphenol	8.657	107	3176	7.849	ng	95
37) 2-Methylnaphthalene	8.822	142	7651	8.493	ng	99
38) 1-Methylnaphthalene	8.922	142	6753	8.012	ng	96
40) 1,2,4,5-Tetrachloroben...	8.986	216	3386	8.132	ng	96
41) Hexachlorocyclopentadiene	8.975	237	865	10.481	ng	88
43) 2,4,6-Trichlorophenol	9.098	196	1906m	6.876	ng	

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.139	196	2276	8.111	ng	91
46) 1,1'-Biphenyl	9.286	154	9351	8.650	ng	98
47) 2-Chloronaphthalene	9.310	162	7121	8.280	ng	90
48) 2-Nitroaniline	9.404	65	2030	8.035	ng #	84
49) Acenaphthylene	9.727	152	11408	8.732	ng	94
50) Dimethylphthalate	9.580	163	8287	8.430	ng #	97
51) 2,6-Dinitrotoluene	9.645	165	1617	7.387	ng #	87
52) Acenaphthene	9.898	154	6448	8.141	ng	100
53) 3-Nitroaniline	9.816	138	1755	7.612	ng #	64
54) 2,4-Dinitrophenol	9.927	184	299	7.937	ng #	42
55) Dibenzofuran	10.069	168	10545	8.519	ng	98
56) 4-Nitrophenol	9.969	139	1004	6.606	ng #	78
57) 2,4-Dinitrotoluene	10.051	165	2321	7.809	ng #	80
58) Fluorene	10.410	166	8032	8.463	ng	95
59) 2,3,4,6-Tetrachlorophenol	10.186	232	1626	7.723	ng #	85
60) Diethylphthalate	10.280	149	8428	8.354	ng	99
61) 4-Chlorophenyl-phenyle...	10.404	204	4030	8.992	ng	91
62) 4-Nitroaniline	10.421	138	1521	6.795	ng	91
63) Azobenzene	10.563	77	7849	7.956	ng	98
65) 4,6-Dinitro-2-methylph...	10.457	198	998	6.095	ng	85
66) n-Nitrosodiphenylamine	10.521	169	6508	7.683	ng	96
67) 4-Bromophenyl-phenylether	10.892	248	2183	7.626	ng	92
68) Hexachlorobenzene	10.957	284	2558	8.214	ng	96
69) Atrazine	11.045	200	2236	8.785	ng	94
70) Pentachlorophenol	11.157	266	572	7.303	ng #	84
71) Phenanthrene	11.374	178	11946	8.426	ng	98
72) Anthracene	11.427	178	12312	8.657	ng	96
73) Carbazole	11.580	167	10201	7.821	ng	95
74) Di-n-butylphthalate	11.910	149	13938	8.409	ng #	97
75) Fluoranthene	12.563	202	11899	8.113	ng	96
77) Benzidine	12.680	184	6542	12.049	ng	97
78) Pyrene	12.792	202	12358	7.732	ng	96
80) Butylbenzylphthalate	13.404	149	5139	7.225	ng	93
81) Benzo(a)anthracene	13.974	228	10541	7.971	ng	96
82) 3,3'-Dichlorobenzidine	13.939	252	3826	10.974	ng	96
83) Chrysene	14.015	228	9795	7.662	ng	99
84) Bis(2-ethylhexyl)phtha...	13.962	149	7721	7.825	ng	98
85) Di-n-octyl phthalate	14.574	149	12063	7.749	ng #	98
87) Indeno(1,2,3-cd)pyrene	16.892	276	7897	7.725	ng #	91
88) Benzo(b)fluoranthene	15.015	252	10278	8.527	ng	96
89) Benzo(k)fluoranthene	15.045	252	8641m	7.880	ng	
90) Benzo(a)pyrene	15.380	252	8175	7.964	ng	96
91) Dibenzo(a,h)anthracene	16.903	278	6720	7.476	ng #	95
92) Benzo(g,h,i)perylene	17.333	276	6130	7.110	ng #	85

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

