

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071421\
 Data File : BF124566.D
 Acq On : 14 Jul 2021 16:11
 Operator : JU/CG
 Sample : M2940-09
 Misc : MDL-MDL-WATER 4ppm
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MDL-WATER-03-QT3-2021

Manual Integrations
 APPROVED

mohammad
 7/16/2021 10:55:10 AM

Quant Time: Jul 14 16:44:03 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF070721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 14 13:25:43 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.928	152	6750	20.000	ng	0.00
21) Naphthalene-d8	8.210	136	26241	20.000	ng	0.00
39) Acenaphthene-d10	9.963	164	14619	20.000	ng	0.00
64) Phenanthrene-d10	11.451	188	25876	20.000	ng	0.00
76) Chrysene-d12	14.092	240	18447	20.000	ng	0.00
86) Perylene-d12	15.592	264	14285	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.545	112	45079	115.708	ng	0.00
7) Phenol-d6	6.545	99	54700	115.415	ng	0.00
23) Nitrobenzene-d5	7.486	82	25686	61.792	ng	0.00
42) 2,4,6-Tribromophenol	10.757	330	12899	113.385	ng	0.00
45) 2-Fluorobiphenyl	9.286	172	63554	63.395	ng	0.00
79) Terphenyl-d14	13.045	244	68369	62.235	ng	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	2.799	88	354	2.778	ng	# 100
3) Pyridine	3.599	79	584	1.787	ng	# 51
4) n-Nitrosodimethylamine	3.481	42	925	3.482	ng	# 67
6) Aniline	6.587	93	2211	4.428	ng	98
8) 2-Chlorophenol	6.710	128	1773	4.059	ng	94
9) Benzaldehyde	6.475	77	1112	3.999	ng	93
10) Phenol	6.557	94	1896	4.098	ng	95
11) bis(2-Chloroethyl)ether	6.657	93	1465	4.084	ng	96
12) 1,3-Dichlorobenzene	6.869	146	2021	4.228	ng	# 87
13) 1,4-Dichlorobenzene	6.945	146	2009	4.114	ng	87
14) 1,2-Dichlorobenzene	7.098	146	1935m	4.170	ng	
15) Benzyl Alcohol	7.051	79	985	3.004	ng	# 86
16) 2,2'-oxybis(1-Chloropr...	7.192	45	2487	4.022	ng	75
17) 2-Methylphenol	7.157	107	1253	3.919	ng	94
18) Hexachloroethane	7.439	117	666	3.622	ng	# 65
19) n-Nitroso-di-n-propyla...	7.328	70	963	3.614	ng	# 78
20) 3+4-Methylphenols	7.310	107	1748	4.153	ng	# 74
22) Acetophenone	7.328	105	2571	4.317	ng	# 93
24) Nitrobenzene	7.504	77	1555	3.773	ng	96
25) Isophorone	7.734	82	2677	3.508	ng	# 86
26) 2-Nitrophenol	7.816	139	729	3.279	ng	# 73
27) 2,4-Dimethylphenol	7.845	122	1920	5.211	ng	98
28) bis(2-Chloroethoxy)met...	7.951	93	1828	4.107	ng	93
29) 2,4-Dichlorophenol	8.051	162	1436	3.779	ng	93
30) 1,2,4-Trichlorobenzene	8.145	180	1750	4.161	ng	98
31) Naphthalene	8.228	128	5516	4.064	ng	99
32) Benzoic acid	7.886	122	281	1.094	ng	88
33) 4-Chloroaniline	8.269	127	2061	4.018	ng	96
34) Hexachlorobutadiene	8.345	225	801	3.374	ng	95
35) Caprolactam	8.604	113	110	1.144	ng	# 35
36) 4-Chloro-3-methylphenol	8.739	107	1359	3.616	ng	# 74
37) 2-Methylnaphthalene	8.922	142	3556	3.883	ng	96
38) 1-Methylnaphthalene	9.016	142	3537	4.112	ng	98
40) 1,2,4,5-Tetrachloroben...	9.081	216	1550	3.975	ng	# 95
41) Hexachlorocyclopentadiene	9.075	237	1243	5.103	ng	87
43) 2,4,6-Trichlorophenol	9.192	196	742	2.636	ng	# 81

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.222	196	911	3.138	ng	# 87
46) 1,1'-Biphenyl	9.380	154	4698	4.231	ng	96
47) 2-Chloronaphthalene	9.410	162	3314	3.807	ng	96
48) 2-Nitroaniline	9.498	65	683	3.263	ng	97
49) Acenaphthylene	9.822	152	5286	3.891	ng	98
50) Dimethylphthalate	9.675	163	4165	4.095	ng	# 97
51) 2,6-Dinitrotoluene	9.739	165	600	2.992	ng	93
52) Acenaphthene	9.998	154	3356	3.864	ng	95
53) 3-Nitroaniline	9.904	138	854	3.817	ng	# 95
54) 2,4-Dinitrophenol	10.004	184	130	1.243	ng	# 1
55) Dibenzofuran	10.169	168	4827	3.780	ng	95
56) 4-Nitrophenol	10.045	139	659	3.539	ng	# 86
57) 2,4-Dinitrotoluene	10.139	165	983	3.581	ng	# 76
58) Fluorene	10.510	166	4207	4.259	ng	98
59) 2,3,4,6-Tetrachlorophenol	10.280	232	783	3.492	ng	# 93
60) Diethylphthalate	10.380	149	4287	4.031	ng	97
61) 4-Chlorophenyl-phenyle...	10.504	204	1724	3.845	ng	96
62) 4-Nitroaniline	10.510	138	886	3.929	ng	# 81
63) Azobenzene	10.669	77	3478	3.741	ng	91
65) 4,6-Dinitro-2-methylph...	10.545	198	348	2.354	ng	# 22
66) n-Nitrosodiphenylamine	10.622	169	3188	3.784	ng	95
67) 4-Bromophenyl-phenylether	10.992	248	925	3.457	ng	# 77
68) Hexachlorobenzene	11.057	284	983	3.586	ng	# 90
69) Atrazine	11.139	200	1093	4.236	ng	95
70) Pentachlorophenol	11.251	266	487	2.809	ng	# 78
71) Phenanthrene	11.474	178	5440	3.853	ng	96
72) Anthracene	11.527	178	5936	4.191	ng	98
73) Carbazole	11.674	167	5040	3.851	ng	98
74) Di-n-butylphthalate	12.010	149	7149	4.055	ng	98
75) Fluoranthene	12.663	202	5604	3.935	ng	96
77) Benzidine	12.786	184	3333	4.833	ng	99
78) Pyrene	12.892	202	6129	3.999	ng	99
80) Butylbenzylphthalate	13.510	149	2770	3.738	ng	96
81) Benzo(a)anthracene	14.080	228	4725	3.862	ng	99
82) 3,3'-Dichlorobenzidine	14.045	252	1715	5.352	ng	# 93
83) Chrysene	14.115	228	4686	3.998	ng	96
84) Bis(2-ethylhexyl)phtha...	14.068	149	3731	3.750	ng	# 94
85) Di-n-octyl phthalate	14.686	149	5093	3.538	ng	# 97
87) Indeno(1,2,3-cd)pyrene	17.086	276	2861	3.453	ng	# 83
88) Benzo(b)fluoranthene	15.145	252	4184m	4.081	ng	
89) Benzo(k)fluoranthene	15.174	252	3699	3.932	ng	99
90) Benzo(a)pyrene	15.515	252	3345	3.952	ng	96
91) Dibenzo(a,h)anthracene	17.109	278	2505	3.552	ng	95
92) Benzo(g,h,i)perylene	17.551	276	2409	3.499	ng	# 70

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

