

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111020\
 Data File : BF122296.D
 Acq On : 10 Nov 2020 13:52
 Operator : JU/CG
 Sample : L4214-07
 Misc : LOD-MDL-WATER 4ppm
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LOD-MDL-WATER-01-QT4-2020

Manual Integrations
 APPROVED

mohammad
 11/11/2020 10:32:05 AM

Quant Time: Nov 10 14:50:14 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 10 13:10:47 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	323780	20.00	ng	0.00
21) Naphthalene-d8	8.151	136	1159697	20.00	ng	0.00
39) Acenaphthene-d10	9.904	164	615592	20.00	ng	0.00
64) Phenanthrene-d10	11.386	188	1123862	20.00	ng	0.00
76) Chrysene-d12	14.027	240	1020087	20.00	ng	0.00
86) Perylene-d12	15.492	264	982243	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	2698355	127.97	ng	0.00
7) Phenol-d6	6.492	99	3500786	126.91	ng	0.00
23) Nitrobenzene-d5	7.434	82	2252105	118.89	ng	0.00
42) 2,4,6-Tribromophenol	10.698	330	971294	147.01	ng	0.00
45) 2-Fluorobiphenyl	9.233	172	3626459	92.04	ng	0.00
79) Terphenyl-d14	12.980	244	4235830	87.97	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.652	88	40185	4.16	ng	87
3) Pyridine	3.457	79	99944	3.76	ng	93
4) n-Nitrosodimethylamine	3.334	42	47320	3.77	ng	# 96
6) Aniline	6.534	93	145365	4.02	ng	97
8) 2-Chlorophenol	6.651	128	88790	4.06	ng	94
9) Benzaldehyde	6.416	77	78002	5.01	ng	96
10) Phenol	6.504	94	107112m	3.94	ng	
11) bis(2-Chloroethyl)ether	6.598	93	89827	4.16	ng	95
12) 1,3-Dichlorobenzene	6.804	146	103465	4.17	ng	100
13) 1,4-Dichlorobenzene	6.881	146	106491	4.27	ng	99
14) 1,2-Dichlorobenzene	7.039	146	98584	4.24	ng	98
15) Benzyl Alcohol	6.998	79	65902	3.36	ng	96
16) 2,2'-oxybis(1-Chloropr...	7.139	45	127769	4.06	ng	98
17) 2-Methylphenol	7.104	107	73204	4.01	ng	100
18) Hexachloroethane	7.381	117	35190	4.05	ng	97
19) n-Nitroso-di-n-propyla...	7.275	70	68011	4.13	ng	95
20) 3+4-Methylphenols	7.257	107	92061	4.10	ng	# 69
22) Acetophenone	7.269	105	134114	4.44	ng	# 87
24) Nitrobenzene	7.445	77	105889	4.89	ng	99
25) Isophorone	7.681	82	178999	4.24	ng	99
26) 2-Nitrophenol	7.763	139	28628	3.91	ng	97
27) 2,4-Dimethylphenol	7.792	122	86542	4.34	ng	98
28) bis(2-Chloroethoxy)met...	7.892	93	110784	4.29	ng	99
29) 2,4-Dichlorophenol	7.998	162	71073	4.03	ng	97
30) 1,2,4-Trichlorobenzene	8.092	180	83500	4.29	ng	96
31) Naphthalene	8.169	128	273506	4.44	ng	98
32) Benzoic acid	7.834	122	20470	2.28	ng	99
33) 4-Chloroaniline	8.210	127	111030	4.13	ng	98
34) Hexachlorobutadiene	8.292	225	47970	4.29	ng	98
35) Caprolactam	8.539	113	20638	3.69	ng	90
36) 4-Chloro-3-methylphenol	8.681	107	73781	4.01	ng	99
37) 2-Methylnaphthalene	8.863	142	179061	4.40	ng	99
38) 1-Methylnaphthalene	8.963	142	165579	4.34	ng	99
40) 1,2,4,5-Tetrachloroben...	9.028	216	83440	4.40	ng	99
41) Hexachlorocyclopentadiene	9.022	237	42109	3.80	ng	99
43) 2,4,6-Trichlorophenol	9.133	196	49674	4.03	ng	100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.169	196	51725	4.01	ng	98
46) 1,1'-Biphenyl	9.328	154	223869	4.57	ng	98
47) 2-Chloronaphthalene	9.351	162	171595	4.41	ng	95
48) 2-Nitroaniline	9.433	65	37542	3.93	ng	94
49) Acenaphthylene	9.763	152	260719	4.36	ng	99
50) Dimethylphthalate	9.616	163	186893	4.27	ng	99
51) 2,6-Dinitrotoluene	9.675	165	34035	4.02	ng	92
52) Acenaphthene	9.933	154	163380	4.42	ng	100
53) 3-Nitroaniline	9.839	138	39304	4.02	ng	# 93
54) 2,4-Dinitrophenol	9.945	184	5882	2.28	ng	# 1
55) Dibenzofuran	10.104	168	239356	4.42	ng	99
56) 4-Nitrophenol	9.986	139	29012	3.47	ng	91
57) 2,4-Dinitrotoluene	10.075	165	38571	3.77	ng	92
58) Fluorene	10.451	166	191593	4.62	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.222	232	43029m	4.01	ng	
60) Diethylphthalate	10.322	149	187467	4.37	ng	99
61) 4-Chlorophenyl-phenyle...	10.445	204	90226	4.60	ng	98
62) 4-Nitroaniline	10.445	138	37975	3.95	ng	87
63) Azobenzene	10.604	77	192027	4.26	ng	98
65) 4,6-Dinitro-2-methylph...	10.480	198	9452	2.21	ng	78
66) n-Nitrosodiphenylamine	10.557	169	162153	4.23	ng	94
67) 4-Bromophenyl-phenylether	10.933	248	55537	4.21	ng	99
68) Hexachlorobenzene	10.998	284	59841	4.19	ng	99
69) Atrazine	11.080	200	45471	4.77	ng	96
70) Pentachlorophenol	11.186	266	27556	3.35	ng	98
71) Phenanthrene	11.410	178	284617	4.46	ng	100
72) Anthracene	11.463	178	287028	4.37	ng	97
73) Carbazole	11.610	167	256037	4.28	ng	98
74) Di-n-butylphthalate	11.951	149	262423	4.12	ng	99
75) Fluoranthene	12.598	202	283983	4.27	ng	96
77) Benzidine	12.716	184	135004	4.77	ng	# 95
78) Pyrene	12.827	202	299994	4.19	ng	98
80) Butylbenzylphthalate	13.451	149	88490	3.43	ng	94
81) Benzo(a)anthracene	14.015	228	279454	4.40	ng	98
82) 3,3'-Dichlorobenzidine	13.974	252	89722	3.80	ng	98
83) Chrysene	14.051	228	273739	4.28	ng	99
84) Bis(2-ethylhexyl)phtha...	14.015	149	138561	4.01	ng	98
85) Di-n-octyl phthalate	14.627	149	195017	3.33	ng	100
87) Indeno(1,2,3-cd)pyrene	16.939	276	229256	3.89	ng	98
88) Benzo(b)fluoranthene	15.057	252	248145	3.90	ng	100
89) Benzo(k)fluoranthene	15.086	252	250528	4.04	ng	98
90) Benzo(a)pyrene	15.421	252	218081	3.93	ng	99
91) Dibenzo(a,h)anthracene	16.962	278	192169	3.89	ng	97
92) Benzo(g,h,i)perylene	17.380	276	187652	3.91	ng	96

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

