

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012323\
 Data File : BM038508.D
 Acq On : 23 Jan 2023 13:16
 Operator : CG/JU
 Sample : N4955-08
 Misc : LOQ-WATER 5ppm
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 LOQ-WATER-02-QT4-2022

Quant Time: Jan 24 05:11:56 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM012023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jan 21 01:21:49 2023
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Christian
 Giraldo

01/24/2023

Supervised By :Jagrut
 Upadhyay

01/24/2023

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.963	152	59551	20.000	ng	0.00
21) Naphthalene-d8	10.775	136	188705	20.000	ng	0.00
39) Acenaphthene-d10	14.604	164	120137	20.000	ng	0.00
64) Phenanthrene-d10	17.345	188	275592	20.000	ng	# 0.00
76) Chrysene-d12	21.521	240	304650	20.000	ng	0.00
86) Perylene-d12	23.933	264	401687	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	324918	90.197	ng	0.00
7) Phenol-d6	7.128	99	499722	108.727	ng	0.00
23) Nitrobenzene-d5	9.139	82	319990	73.637	ng	0.00
42) 2,4,6-Tribromophenol	16.092	330	227544	131.528	ng	0.00
45) 2-Fluorobiphenyl	13.227	172	714511	73.262	ng	0.00
79) Terphenyl-d14	19.956	244	1157339	74.060	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.375	88	7604	4.500	ng	95
3) Pyridine	3.810	79	16747m	3.774	ng	
4) n-Nitrosodimethylamine	3.710	42	9611	4.562	ng	# 83
6) Aniline	7.292	93	24375	4.273	ng	# 96
8) 2-Chlorophenol	7.522	128	18292	4.971	ng	91
9) Benzaldehyde	7.110	77	17559m	6.441	ng	
10) Phenol	7.157	94	22619	4.925	ng	87
11) bis(2-Chloroethyl)ether	7.387	93	17914	4.944	ng	95
12) 1,3-Dichlorobenzene	7.851	146	20770	4.965	ng	95
13) 1,4-Dichlorobenzene	7.998	146	21318	5.043	ng	97
14) 1,2-Dichlorobenzene	8.316	146	20263	4.949	ng	94
15) Benzyl Alcohol	8.210	79	16330	4.752	ng	93
16) 2,2'-oxybis(1-Chloropr...	8.487	45	17302	4.547	ng	94
17) 2-Methylphenol	8.416	107	12656	3.893	ng	94
18) Hexachloroethane	9.045	117	7794	4.960	ng	94
19) n-Nitroso-di-n-propyla...	8.775	70	13352	4.100	ng	# 94
20) 3+4-Methylphenols	8.739	107	17107	3.927	ng	93
22) Acetophenone	8.798	105	24735	4.758	ng	# 99
24) Nitrobenzene	9.181	77	24192m	5.502	ng	
25) Isophorone	9.704	82	33508	4.673	ng	97
26) 2-Nitrophenol	9.892	139	7718	4.422	ng	98
27) 2,4-Dimethylphenol	9.945	122	9350	3.219	ng	95
28) bis(2-Chloroethoxy)met...	10.192	93	19118	4.557	ng	95
29) 2,4-Dichlorophenol	10.428	162	15382	4.514	ng	97
30) 1,2,4-Trichlorobenzene	10.639	180	19771	5.073	ng	97
31) Naphthalene	10.828	128	48397	4.721	ng	98
32) Benzoic acid	10.028	122	5601m	7.765	ng	
33) 4-Chloroaniline	10.951	127	18043	4.217	ng	96
34) Hexachlorobutadiene	11.098	225	13870	5.411	ng	93
35) Caprolactam	11.727	113	2304	2.789	ng	92
36) 4-Chloro-3-methylphenol	12.069	107	13535	4.368	ng	92
37) 2-Methylnaphthalene	12.433	142	31351	4.255	ng	# 90
38) 1-Methylnaphthalene	12.651	142	29457m	4.239	ng	
40) 1,2,4,5-Tetrachloroben...	12.798	216	20891	4.623	ng	99
41) Hexachlorocyclopentadiene	12.769	237	9398	3.784	ng	99
43) 2,4,6-Trichlorophenol	13.039	196	12364	4.274	ng	97

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44) 2,4,5-Trichlorophenol	13.116	196	13901	4.103	ng	95
46) 1,1'-Biphenyl	13.439	154	43195	4.489	ng	98
47) 2-Chloronaphthalene	13.480	162	35805	4.767	ng	95
48) 2-Nitroaniline	13.698	65	8646	3.920	ng	92
49) Acenaphthylene	14.327	152	49063	4.175	ng	99
50) Dimethylphthalate	14.057	163	42325	4.577	ng	99
51) 2,6-Dinitrotoluene	14.186	165	7670	3.941	ng	90
52) Acenaphthene	14.663	154	32035	4.470	ng	98
53) 3-Nitroaniline	14.521	138	6452	3.336	ng	# 89
54) 2,4-Dinitrophenol	14.733	184	4031	7.353	ng	90
55) Dibenzofuran	14.998	168	54958	4.587	ng	96
56) 4-Nitrophenol	14.827	139	5037	3.163	ng	95
57) 2,4-Dinitrotoluene	14.974	165	9625	3.648	ng	# 87
58) Fluorene	15.645	166	42412	4.365	ng	93
59) 2,3,4,6-Tetrachlorophenol	15.227	232	12242	4.420	ng	# 96
60) Diethylphthalate	15.415	149	39179	4.326	ng	98
61) 4-Chlorophenyl-phenyle...	15.639	204	23502	4.607	ng	98
62) 4-Nitroaniline	15.680	138	6187	5.983	ng	88
63) Azobenzene	15.933	77	40675	4.239	ng	100
65) 4,6-Dinitro-2-methylph...	15.733	198	6028	3.370	ng	90
66) n-Nitrosodiphenylamine	15.857	169	34955	4.225	ng	97
67) 4-Bromophenyl-phenylether	16.533	248	14043	4.594	ng	99
68) Hexachlorobenzene	16.645	284	17255	5.231	ng	95
69) Atrazine	16.804	200	12536	4.542	ng	99
70) Pentachlorophenol	16.998	266	8887	3.784	ng	93
71) Phenanthrene	17.386	178	69978	4.547	ng	99
72) Anthracene	17.480	178	68022	4.344	ng	97
73) Carbazole	17.751	167	59349	4.303	ng	98
74) Di-n-butylphthalate	18.304	149	60205	3.974	ng	# 98
75) Fluoranthene	19.398	202	81799	4.422	ng	96
77) Benzidine	19.592	184	14612	2.105	ng	100
78) Pyrene	19.762	202	86769	4.090	ng	99
80) Butylbenzylphthalate	20.645	149	25674	3.500	ng	95
81) Benzo(a)anthracene	21.503	228	99434	4.606	ng	99
82) 3,3'-Dichlorobenzidine	21.439	252	29509	4.256	ng	# 97
83) Chrysene	21.556	228	95894	4.542	ng	99
84) Bis(2-ethylhexyl)phtha...	21.415	149	36957	3.476	ng	96
85) Di-n-octyl phthalate	22.344	149	70655	3.738	ng	96
87) Indeno(1,2,3-cd)pyrene	26.444	276	140107	4.714	ng	98
88) Benzo(b)fluoranthene	23.197	252	112495	4.643	ng	98
89) Benzo(k)fluoranthene	23.244	252	109259	4.358	ng	99
90) Benzo(a)pyrene	23.827	252	102449	4.264	ng	96
91) Dibenzo(a,h)anthracene	26.462	278	121529	4.912	ng	# 96
92) Benzo(g,h,i)perylene	27.226	276	126992	5.011	ng	96

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

