

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG051823\
 Data File : BG057457.D
 Acq On : 18 May 2023 11:52
 Operator : CG/JU
 Sample : 02213-07
 Misc : LOD-MDL-WATER-4PPM
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LOD-MDL-WATER-01-QT2-2023

Manual Integrations
 APPROVED

Reviewed By :Christian
 Giraldo

Quant Time: May 18 13:16:11 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG042723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 12 22:04:03 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.363	152	16652	20.000	ng	0.00
21) Naphthalene-d8	11.194	136	69875	20.000	ng	0.00
39) Acenaphthene-d10	14.972	164	56792	20.000	ng	0.00
64) Phenanthrene-d10	17.715	188	143687	20.000	ng	0.00
76) Chrysene-d12	22.021	240	133725	20.000	ng	-0.01
86) Perylene-d12	25.516	264	140828	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.872	112	99798	102.520	ng	0.00
7) Phenol-d6	7.499	99	149065	100.840	ng	0.00
23) Nitrobenzene-d5	9.538	82	109394	65.708	ng	0.00
42) 2,4,6-Tribromophenol	16.458	330	85069	105.568	ng	0.00
45) 2-Fluorobiphenyl	13.597	172	300881	71.353	ng	0.00
79) Terphenyl-d14	20.282	244	565714	69.037	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.675	88	1465	3.640	ng	91
3) Pyridine	4.122	79	3487	2.696	ng	# 75
4) n-Nitrosodimethylamine	4.016	42	1605	2.640	ng	# 98
6) Aniline	7.676	93	6427	3.427	ng	# 89
8) 2-Chlorophenol	7.916	128	3770	3.665	ng	96
10) Phenol	7.523	94	5154	3.577	ng	87
11) bis(2-Chloroethyl)ether	7.758	93	3813	3.509	ng	89
12) 1,3-Dichlorobenzene	8.251	146	4479	3.945	ng	96
13) 1,4-Dichlorobenzene	8.392	146	4392	3.851	ng	97
14) 1,2-Dichlorobenzene	8.721	146	4218m	3.832	ng	
15) Benzyl Alcohol	8.592	79	3367	2.734	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.874	45	4434	3.238	ng	88
17) 2-Methylphenol	8.798	107	3311	3.187	ng	97
18) Hexachloroethane	9.455	117	1477	3.589	ng	97
19) n-Nitroso-di-n-propyla...	9.156	70	3514	3.118	ng	# 86
20) 3+4-Methylphenols	9.127	107	4528	3.183	ng	86
22) Acetophenone	9.185	105	7046	3.523	ng	# 90
24) Nitrobenzene	9.579	77	5862	3.489	ng	93
25) Isophorone	10.096	82	10130	3.138	ng	98
26) 2-Nitrophenol	10.301	139	2104	3.271	ng	92
27) 2,4-Dimethylphenol	10.343	122	3648	3.239	ng	94
28) bis(2-Chloroethoxy)met...	10.572	93	5362	3.315	ng	100
29) 2,4-Dichlorophenol	10.859	162	4079	3.415	ng	# 79
30) 1,2,4-Trichlorobenzene	11.059	180	5055	3.669	ng	94
31) Naphthalene	11.247	128	13148	3.520	ng	# 95
33) 4-Chloroaniline	11.359	127	5496	3.276	ng	# 92
34) Hexachlorobutadiene	11.517	225	3725	3.644	ng	# 76
35) Caprolactam	12.099	113	1179	2.889	ng	# 91
36) 4-Chloro-3-methylphenol	12.451	107	4390	3.234	ng	91
37) 2-Methylnaphthalene	12.827	142	10404	3.542	ng	97
38) 1-Methylnaphthalene	13.045	142	9505m	3.433	ng	
40) 1,2,4,5-Tetrachloroben...	13.186	216	7604	3.772	ng	94
41) Hexachlorocyclopentadiene	13.156	237	289	6.570	ng	# 27
43) 2,4,6-Trichlorophenol	13.427	196	3853	3.137	ng	94
44) 2,4,5-Trichlorophenol	13.515	196	4477	3.099	ng	# 92
46) 1,1'-Biphenyl	13.808	154	15529	3.685	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
47) 2-Chloronaphthalene	13.867	162	11454	3.655	ng	95	05/19/2023
48) 2-Nitroaniline	14.061	65	3065	2.876	ng	93	Supervised By :Jagrut Upadhyay
49) Acenaphthylene	14.701	152	17707	3.478	ng	94	
50) Dimethylphthalate	14.402	163	15509	3.323	ng	97	
51) 2,6-Dinitrotoluene	14.537	165	3128	3.283	ng	90	
52) Acenaphthene	15.036	154	11537	3.626	ng	97	
53) 3-Nitroaniline	14.878	138	2755	2.886	ng	# 77	05/22/2023
55) Dibenzofuran	15.365	168	19301	3.611	ng	97	
56) 4-Nitrophenol	15.207	139	1054	1.660	ng	# 81	
57) 2,4-Dinitrotoluene	15.330	165	4409	3.238	ng	# 92	
58) Fluorene	16.011	166	15296	3.474	ng	98	
59) 2,3,4,6-Tetrachlorophenol	15.600	232	4407	3.364	ng	# 93	
60) Diethylphthalate	15.753	149	15574	3.286	ng	94	
61) 4-Chlorophenyl-phenyle...	15.994	204	9138	3.445	ng	98	
62) 4-Nitroaniline	16.035	138	2759	2.865	ng	# 86	
63) Azobenzene	16.287	77	14135	3.091	ng	96	
65) 4,6-Dinitro-2-methylph...	16.099	198	1545	1.801	ng	91	
66) n-Nitrosodiphenylamine	16.205	169	13493	3.457	ng	98	
67) 4-Bromophenyl-phenylether	16.887	248	6050	3.449	ng	92	
68) Hexachlorobenzene	17.022	284	6720	3.942	ng	# 92	
69) Atrazine	17.133	200	5997	4.028	ng	99	
70) Pentachlorophenol	17.409	266	369	0.376	ng	# 20	
71) Phenanthrene	17.756	178	26790	3.639	ng	98	
72) Anthracene	17.844	178	26704	3.535	ng	99	
73) Carbazole	18.114	167	22356	3.515	ng	97	
74) Di-n-butylphthalate	18.625	149	23447	3.204	ng	98	
75) Fluoranthene	19.747	202	32394	3.537	ng	98	
77) Benzidine	19.912	184	13652	4.566	ng	# 94	
78) Pyrene	20.106	202	33345	3.432	ng	96	
80) Butylbenzylphthalate	20.958	149	10454	3.391	ng	# 87	
81) Benzo(a)anthracene	21.997	228	32346	3.395	ng	97	
82) 3,3'-Dichlorobenzidine	21.897	252	12903	4.228	ng	92	
83) Chrysene	22.068	228	31447	3.508	ng	99	
84) Bis(2-ethylhexyl)phtha...	21.845	149	14820	3.249	ng	100	
85) Di-n-octyl phthalate	23.131	149	24390	3.207	ng	99	
87) Indeno(1,2,3-cd)pyrene	29.558	276	34366	3.348	ng	# 95	
88) Benzo(b)fluoranthene	24.394	252	33025	3.587	ng	98	
89) Benzo(k)fluoranthene	24.470	252	32091	3.430	ng	98	
90) Benzo(a)pyrene	25.352	252	28452	3.164	ng	99	
91) Dibenzo(a,h)anthracene	29.605	278	29650	3.469	ng	# 95	
92) Benzo(g,h,i)perylene	30.838	276	28811	3.452	ng	97	

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

