

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020924\
 Data File : BM044040.D
 Acq On : 09 Feb 2024 15:30
 Operator : MA/JU
 Sample : 04699-07
 Misc : LOD-MDL-WATER-8ppm
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 LOD-MDL-WATER-01-QT4-2023

Quant Time: Feb 09 16:18:30 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM020824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 09 02:42:01 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.910	152	601786	20.000	ng	0.00
21) Naphthalene-d8	10.710	136	2560261	20.000	ng	0.00
39) Acenaphthene-d10	14.551	164	1508151	20.000	ng	0.00
64) Phenanthrene-d10	17.303	188	2968050	20.000	ng	0.00
76) Chrysene-d12	21.503	240	2299184	20.000	ng	0.00
86) Perylene-d12	23.897	264	2438909	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	3299143	79.736	ng	0.00
7) Phenol-d6	7.075	99	4343644	82.509	ng	0.00
23) Nitrobenzene-d5	9.080	82	2929028	60.707	ng	0.00
42) 2,4,6-Tribromophenol	16.045	330	1338433	81.966	ng	0.00
45) 2-Fluorobiphenyl	13.180	172	6210291	58.116	ng	0.00
79) Terphenyl-d14	19.944	244	7740580	59.019	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.387	88	103893	6.327	ng	100
3) Pyridine	3.787	79	236114	5.446	ng	98
4) n-Nitrosodimethylamine	3.710	42	94629	5.753	ng	# 99
6) Aniline	7.239	93	377076	7.285	ng	100
8) 2-Chlorophenol	7.469	128	260038	6.018	ng	98
9) Benzaldehyde	7.057	77	234682m	8.675	ng	
10) Phenol	7.104	94	330356	6.218	ng	99
11) bis(2-Chloroethyl)ether	7.345	93	272942	6.204	ng	98
12) 1,3-Dichlorobenzene	7.792	146	290129	5.914	ng	98
13) 1,4-Dichlorobenzene	7.945	146	294121	5.948	ng	100
14) 1,2-Dichlorobenzene	8.257	146	277219	5.884	ng	99
15) Benzyl Alcohol	8.151	79	214936	6.270	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.451	45	351925	6.310	ng	98
17) 2-Methylphenol	8.351	107	208272	5.972	ng	98
18) Hexachloroethane	8.980	117	109231	5.900	ng	98
19) n-Nitroso-di-n-propyla...	8.722	70	209610	6.484	ng	98
20) 3+4-Methylphenols	8.681	107	284821	5.934	ng	97
22) Acetophenone	8.739	105	424967	6.539	ng	# 98
24) Nitrobenzene	9.116	77	304467	6.162	ng	97
25) Isophorone	9.639	82	567755	6.111	ng	98
26) 2-Nitrophenol	9.827	139	124074	5.416	ng	99
27) 2,4-Dimethylphenol	9.886	122	188681	6.442	ng	98
28) bis(2-Chloroethoxy)met...	10.133	93	373281	6.213	ng	99
29) 2,4-Dichlorophenol	10.357	162	225514	5.848	ng	98
30) 1,2,4-Trichlorobenzene	10.569	180	254693	5.873	ng	99
31) Naphthalene	10.757	128	858397	6.075	ng	100
32) Benzoic acid	9.969	122	106032	3.594	ng	98
33) 4-Chloroaniline	10.880	127	347272	6.490	ng	99
34) Hexachlorobutadiene	11.033	225	141012	5.681	ng	98
35) Caprolactam	11.651	113	70064	5.646	ng	97
36) 4-Chloro-3-methylphenol	11.998	107	243642	5.835	ng	99
37) 2-Methylnaphthalene	12.374	142	566768	6.022	ng	98
38) 1-Methylnaphthalene	12.592	142	554666	5.997	ng	100
40) 1,2,4,5-Tetrachloroben...	12.733	216	270458	6.170	ng	99
41) Hexachlorocyclopentadiene	12.710	237	120602	6.304	ng	98
43) 2,4,6-Trichlorophenol	12.980	196	168445	5.646	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.051	196	185424	5.664	ng	98
46) 1,1'-Biphenyl	13.386	154	766306	6.439	ng	100
47) 2-Chloronaphthalene	13.427	162	566639	6.011	ng	99
48) 2-Nitroaniline	13.639	65	152753	5.760	ng	93
49) Acenaphthylene	14.274	152	935675	6.653	ng	100
50) Dimethylphthalate	14.021	163	700282	6.263	ng	100
51) 2,6-Dinitrotoluene	14.139	165	138879	5.629	ng	93
52) Acenaphthene	14.615	154	534492	6.141	ng	99
53) 3-Nitroaniline	14.468	138	149921	5.655	ng	92
54) 2,4-Dinitrophenol	14.668	184	49855	8.719	ng	98
55) Dibenzofuran	14.951	168	866649	6.433	ng	99
56) 4-Nitrophenol	14.768	139	107987	5.010	ng	95
57) 2,4-Dinitrotoluene	14.921	165	169748	5.264	ng	# 93
58) Fluorene	15.598	166	671935	6.455	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.174	232	157288	6.123	ng	100
60) Diethylphthalate	15.386	149	685027	5.928	ng	99
61) 4-Chlorophenyl-phenyle...	15.604	204	322780	6.399	ng	99
62) 4-Nitroaniline	15.627	138	140605	5.268	ng	94
63) Azobenzene	15.898	77	670207	6.181	ng	99
65) 4,6-Dinitro-2-methylph...	15.680	198	75340	4.260	ng	96
66) n-Nitrosodiphenylamine	15.815	169	566707	6.285	ng	99
67) 4-Bromophenyl-phenylether	16.498	248	144607	4.928	ng	97
68) Hexachlorobenzene	16.592	284	212678	5.916	ng	98
69) Atrazine	16.774	200	181239	7.423	ng	99
70) Pentachlorophenol	16.945	266	107667	4.604	ng	99
71) Phenanthrene	17.345	178	1012476	6.460	ng	100
72) Anthracene	17.433	178	977257	6.298	ng	99
73) Carbazole	17.709	167	910680	6.035	ng	99
74) Di-n-butylphthalate	18.292	149	1103572	5.626	ng	99
75) Fluoranthene	19.362	202	1045375	5.938	ng	99
77) Benzidine	19.562	184	405072	11.337	ng	99
78) Pyrene	19.727	202	1086975	6.139	ng	99
80) Butylbenzylphthalate	20.650	149	411202	5.224	ng	99
81) Benzo(a)anthracene	21.486	228	979951	6.127	ng	99
82) 3,3'-Dichlorobenzidine	21.427	252	325199	6.197	ng	99
83) Chrysene	21.538	228	906230	5.998	ng	99
84) Bis(2-ethylhexyl)phtha...	21.439	149	643944	5.545	ng	99
85) Di-n-octyl phthalate	22.374	149	990546	4.884	ng	98
87) Indeno(1,2,3-cd)pyrene	26.379	276	1001841	5.804	ng	100
88) Benzo(b)fluoranthene	23.168	252	904666	5.934	ng	99
89) Benzo(k)fluoranthene	23.215	252	876963	6.053	ng	100
90) Benzo(a)pyrene	23.791	252	766559	5.756	ng	99
91) Dibenzo(a,h)anthracene	26.403	278	850285	5.889	ng	100
92) Benzo(g,h,i)perylene	27.132	276	808372	5.520	ng	99

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

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