

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072121\
 Data File : BM031044.D
 Acq On : 21 Jul 2021 15:19
 Operator : CG/JU
 Sample : M2940-09
 Misc : MDL-MDL-WATER- 8PPM
 ALS Vial : 9 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

mohammad
 7/22/2021 1:30:01 PM

Quant Time: Jul 21 15:55:11 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM072021.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 21 11:41:39 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.634	152	72631	20.00	ng	0.00
21) Naphthalene-d8	10.399	136	315413	20.00	ng	-0.01
39) Acenaphthene-d10	14.263	164	221423	20.00	ng	-0.01
64) Phenanthrene-d10	17.010	188	489645	20.00	ng	-0.01
76) Chrysene-d12	21.198	240	553497	20.00	ng	-0.01
86) Perylene-d12	23.380	264	575637	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.252	112	490199	121.19	ng	0.00
7) Phenol-d6	6.816	99	710592	121.19	ng	0.00
23) Nitrobenzene-d5	8.781	82	517069	82.31	ng	-0.01
42) 2,4,6-Tribromophenol	15.763	330	342936	118.61	ng	0.00
45) 2-Fluorobiphenyl	12.887	172	1225876	82.61	ng	0.00
79) Terphenyl-d14	19.651	244	2359189	84.33	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.240	88	11469	7.06	ng	92
3) Pyridine	3.628	79	31014	6.75	ng	# 89
4) n-Nitrosodimethylamine	3.546	42	16328	6.25	ng	# 87
6) Aniline	6.975	93	54237	8.62	ng	99
8) 2-Chlorophenol	7.205	128	35503	7.46	ng	96
9) Benzaldehyde	6.787	77	33513	9.59	ng	97
10) Phenol	6.840	94	43012	7.57	ng	92
11) bis(2-Chloroethyl)ether	7.075	93	33308	7.92	ng	98
12) 1,3-Dichlorobenzene	7.522	146	38985	7.53	ng	96
13) 1,4-Dichlorobenzene	7.669	146	40273	7.64	ng	97
14) 1,2-Dichlorobenzene	7.981	146	38345	7.45	ng	95
15) Benzyl Alcohol	7.875	79	35892	7.27	ng	96
16) 2,2'-oxybis(1-Chloropr...	8.169	45	43489	8.27	ng	97
17) 2-Methylphenol	8.075	107	29119	7.33	ng	94
18) Hexachloroethane	8.699	117	15097	7.42	ng	88
19) n-Nitroso-di-n-propyla...	8.434	70	30334	7.33	ng	98
20) 3+4-Methylphenols	8.399	107	42006	7.47	ng	96
22) Acetophenone	8.452	105	61218	7.84	ng	# 96
24) Nitrobenzene	8.822	77	48618	7.66	ng	99
25) Isophorone	9.346	82	78938	7.03	ng	98
26) 2-Nitrophenol	9.528	139	18239	6.83	ng	# 87
27) 2,4-Dimethylphenol	9.593	122	36219	8.61	ng	97
28) bis(2-Chloroethoxy)met...	9.834	93	47247	8.34	ng	96
29) 2,4-Dichlorophenol	10.051	162	35911	7.21	ng	96
30) 1,2,4-Trichlorobenzene	10.269	180	40298	7.37	ng	93
31) Naphthalene	10.451	128	122282	7.53	ng	98
32) Benzoic acid	9.657	122	17593	4.91	ng	96
33) 4-Chloroaniline	10.569	127	53377	7.62	ng	99
34) Hexachlorobutadiene	10.740	225	24751	7.05	ng	98
35) Caprolactam	11.334	113	11359	6.79	ng	97
36) 4-Chloro-3-methylphenol	11.687	107	40361	7.11	ng	92
37) 2-Methylnaphthalene	12.069	142	91660	7.33	ng	98
38) 1-Methylnaphthalene	12.287	142	85720	7.20	ng	98
40) 1,2,4,5-Tetrachloroben...	12.440	216	47627	7.72	ng	# 98
41) Hexachlorocyclopentadiene	12.422	237	31264	9.11	ng	97
43) 2,4,6-Trichlorophenol	12.687	196	29569	6.82	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
44) 2,4,5-Trichlorophenol	12.751	196	33227	6.93	ng	#	91
46) 1,1'-Biphenyl	13.092	154	125060	7.97	ng		96
47) 2-Chloronaphthalene	13.134	162	92968	7.56	ng		93
48) 2-Nitroaniline	13.339	65	26697	6.81	ng		89
49) Acenaphthylene	13.987	152	146093	7.43	ng		100
50) Dimethylphthalate	13.734	163	123358	7.43	ng		100
51) 2,6-Dinitrotoluene	13.845	165	25094	6.75	ng	#	78
52) Acenaphthene	14.328	154	89895	7.27	ng		99
53) 3-Nitroaniline	14.175	138	26011	6.92	ng		94
54) 2,4-Dinitrophenol	14.381	184	11494	5.24	ng		94
55) Dibenzofuran	14.663	168	147893	7.29	ng		93
56) 4-Nitrophenol	14.481	139	22844	6.87	ng		84
57) 2,4-Dinitrotoluene	14.634	165	32766	6.33	ng	#	89
58) Fluorene	15.316	166	118264	6.91	ng		99
59) 2,3,4,6-Tetrachlorophenol	14.892	232	30138m	6.74	ng		
60) Diethylphthalate	15.104	149	126837	7.12	ng		98
61) 4-Chlorophenyl-phenyle...	15.322	204	59016	6.96	ng	#	86
62) 4-Nitroaniline	15.339	138	27518	6.64	ng		93
63) Azobenzene	15.610	77	120555	6.80	ng		92
65) 4,6-Dinitro-2-methylph...	15.398	198	17145	5.58	ng		88
66) n-Nitrosodiphenylamine	15.533	169	105379	7.32	ng		99
67) 4-Bromophenyl-phenylether	16.216	248	34989	6.91	ng		91
68) Hexachlorobenzene	16.316	284	41888	7.34	ng		97
69) Atrazine	16.492	200	39408	7.46	ng		97
70) Pentachlorophenol	16.663	266	24798	6.64	ng		96
71) Phenanthrene	17.057	178	199418	7.35	ng		100
72) Anthracene	17.145	178	201572	7.54	ng		99
73) Carbazole	17.416	167	185189	7.00	ng		97
74) Di-n-butylphthalate	17.998	149	220469	7.09	ng		99
75) Fluoranthene	19.069	202	238201	7.02	ng		99
77) Benzidine	19.269	184	134018	9.81	ng		99
78) Pyrene	19.433	202	253626	7.34	ng		97
80) Butylbenzylphthalate	20.357	149	106256	7.20	ng		96
81) Benzo(a)anthracene	21.180	228	266013	7.25	ng		98
82) 3,3'-Dichlorobenzidine	21.127	252	97932	9.74	ng		99
83) Chrysene	21.233	228	250915	7.04	ng		98
84) Bis(2-ethylhexyl)phtha...	21.133	149	170212	7.44	ng		95
85) Di-n-octyl phthalate	22.004	149	293114	7.49	ng		98
87) Indeno(1,2,3-cd)pyrene	25.539	276	292831	7.12	ng		99
88) Benzo(b)fluoranthene	22.727	252	276940	7.26	ng		98
89) Benzo(k)fluoranthene	22.768	252	259086	7.26	ng		98
90) Benzo(a)pyrene	23.280	252	261529	7.62	ng		98
91) Dibenzo(a,h)anthracene	25.556	278	248011	6.97	ng		99
92) Benzo(g,h,i)perylene	26.198	276	243166	7.17	ng		99

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

