

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082521\
 Data File : BM031896.D
 Acq On : 25 Aug 2021 17:31
 Operator : CG/JU
 Sample : M2262-08
 Misc : LOQ-WATER-5ppm
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 LOQ-WATER-02-QT2-2021

Manual Integrations
 APPROVED

mohammad
 8/26/2021 11:41:18 AM

Quant Time: Aug 25 18:05:09 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM082521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 25 16:00:28 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.522	152	33698	20.000	ng	0.00
21) Naphthalene-d8	10.281	136	138938	20.000	ng	0.00
39) Acenaphthene-d10	14.157	164	89444	20.000	ng	0.00
64) Phenanthrene-d10	16.915	188	178455	20.000	ng	0.00
76) Chrysene-d12	21.121	240	134059	20.000	ng	0.00
86) Perylene-d12	23.262	264	126341	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.157	112	214290	105.130	ng	0.00
7) Phenol-d6	6.722	99	290388	100.689	ng	0.00
23) Nitrobenzene-d5	8.675	82	333136	103.858	ng	0.00
42) 2,4,6-Tribromophenol	15.663	330	109123	103.211	ng	0.00
45) 2-Fluorobiphenyl	12.775	172	775164	114.339	ng	0.00
79) Terphenyl-d14	19.568	244	960317	103.403	ng	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.175	88	4233	5.094	ng	# 92
3) Pyridine	3.569	79	8387	4.043	ng	# 80
4) n-Nitrosodimethylamine	3.481	42	6047	4.291	ng	# 51
6) Aniline	6.869	93	15629	5.101	ng	94
8) 2-Chlorophenol	7.093	128	10095	4.254	ng	93
9) Benzaldehyde	6.693	77	12238	6.398	ng	96
10) Phenol	6.746	94	12471	4.373	ng	93
11) bis(2-Chloroethyl)ether	6.975	93	10067	4.839	ng	88
12) 1,3-Dichlorobenzene	7.410	146	12553	4.789	ng	97
13) 1,4-Dichlorobenzene	7.557	146	12732	4.741	ng	96
14) 1,2-Dichlorobenzene	7.863	146	12769	4.883	ng	95
15) Benzyl Alcohol	7.775	79	10180	4.138	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.051	45	12893	4.619	ng	91
17) 2-Methylphenol	7.969	107	8425	4.140	ng	94
18) Hexachloroethane	8.569	117	4913	4.416	ng	91
19) n-Nitroso-di-n-propyla...	8.328	70	9458	4.257	ng	# 95
20) 3+4-Methylphenols	8.298	107	12271	4.383	ng	95
22) Acetophenone	8.340	105	18223	4.825	ng	# 99
24) Nitrobenzene	8.716	77	15322m	4.718	ng	
25) Isophorone	9.234	82	24486	4.420	ng	98
26) 2-Nitrophenol	9.416	139	5597	4.305	ng	92
27) 2,4-Dimethylphenol	9.481	122	9959	5.090	ng	88
28) bis(2-Chloroethoxy)met...	9.722	93	13158	4.863	ng	# 96
29) 2,4-Dichlorophenol	9.951	162	9247	4.024	ng	95
30) 1,2,4-Trichlorobenzene	10.145	180	11478	4.553	ng	97
31) Naphthalene	10.328	128	35704	4.599	ng	98
32) Benzoic acid	9.575	122	3724	2.234	ng	92
33) 4-Chloroaniline	10.463	127	13953	4.495	ng	97
34) Hexachlorobutadiene	10.604	225	7503	4.499	ng	95
35) Caprolactam	11.257	113	2791	3.985	ng	# 70
36) 4-Chloro-3-methylphenol	11.586	107	10975	4.027	ng	92
37) 2-Methylnaphthalene	11.951	142	25805	4.464	ng	91
38) 1-Methylnaphthalene	12.169	142	23454	4.289	ng	99
40) 1,2,4,5-Tetrachloroben...	12.328	216	12811	4.889	ng	98
41) Hexachlorocyclopentadiene	12.292	237	5273	4.127	ng	97
43) 2,4,6-Trichlorophenol	12.580	196	7978m	4.240	ng	

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.657	196	8665	4.262	ng	# 93
46) 1,1'-Biphenyl	12.986	154	34359	4.964	ng	92
47) 2-Chloronaphthalene	13.022	162	25373	4.656	ng	96
48) 2-Nitroaniline	13.251	65	7621	4.018	ng	91
49) Acenaphthylene	13.880	152	39573	4.514	ng	99
50) Dimethylphthalate	13.633	163	32067	4.523	ng	# 99
51) 2,6-Dinitrotoluene	13.757	165	6226	3.987	ng	95
52) Acenaphthene	14.222	154	23711	4.442	ng	95
53) 3-Nitroaniline	14.092	138	5521	3.635	ng	98
54) 2,4-Dinitrophenol	14.316	184	2399	2.713	ng	# 79
55) Dibenzofuran	14.563	168	38781	4.314	ng	99
56) 4-Nitrophenol	14.439	139	3998	3.331	ng	# 76
57) 2,4-Dinitrotoluene	14.557	165	8384	3.770	ng	# 96
58) Fluorene	15.216	166	31175	4.230	ng	94
59) 2,3,4,6-Tetrachlorophenol	14.798	232	7059	3.977	ng	97
60) Diethylphthalate	15.010	149	32742	4.231	ng	96
61) 4-Chlorophenyl-phenyle...	15.221	204	15541	4.248	ng	99
62) 4-Nitroaniline	15.269	138	5194	3.541	ng	93
63) Azobenzene	15.510	77	34870	4.087	ng	96
65) 4,6-Dinitro-2-methylph...	15.327	198	3694	3.005	ng	97
66) n-Nitrosodiphenylamine	15.439	169	26102	4.389	ng	98
67) 4-Bromophenyl-phenylether	16.116	248	8620	4.325	ng	96
68) Hexachlorobenzene	16.215	284	9325	4.426	ng	96
69) Atrazine	16.404	200	9087	4.532	ng	95
70) Pentachlorophenol	16.574	266	4642	3.508	ng	96
71) Phenanthrene	16.957	178	45400	4.321	ng	98
72) Anthracene	17.051	178	44749	4.354	ng	99
73) Carbazole	17.333	167	38349	3.951	ng	98
74) Di-n-butylphthalate	17.898	149	45774	3.757	ng	99
75) Fluoranthene	18.986	202	45973	4.007	ng	98
77) Benzidine	19.192	184	21324	6.081	ng	96
78) Pyrene	19.351	202	46654	4.161	ng	99
80) Butylbenzylphthalate	20.274	149	18387	3.813	ng	96
81) Benzo(a)anthracene	21.103	228	41600	4.314	ng	96
82) 3,3'-Dichlorobenzidine	21.056	252	14596	5.264	ng	95
83) Chrysene	21.156	228	40233	4.328	ng	96
84) Bis(2-ethylhexyl)phtha...	21.050	149	25230	3.637	ng	95
85) Di-n-octyl phthalate	21.909	149	43125	4.125	ng	97
87) Indeno(1,2,3-cd)pyrene	25.380	276	41618	4.761	ng	98
88) Benzo(b)fluoranthene	22.633	252	40168	4.199	ng	98
89) Benzo(k)fluoranthene	22.674	252	39612	4.369	ng	96
90) Benzo(a)pyrene	23.174	252	38310	4.624	ng	99
91) Dibenzo(a,h)anthracene	25.391	278	36146	4.774	ng	97
92) Benzo(g,h,i)perylene	26.021	276	35693	4.985	ng	96

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

