

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052218\
 Data File : BM015259.D
 Acq On : 23 May 2018 02:44
 Operator : SJ/JU
 Sample : J2133-09
 Misc : MDL 8 PPM
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-WATER-03-QT2-2018

Manual Integrations
 APPROVED

Sohil
 5/23/2018 10:11:37 AM

Quant Time: May 23 05:49:46 2018
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM052218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 22 15:47:28 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.40	152	228310	20.00	ng	0.00
21) Naphthalene-d8	11.24	136	959000	20.00	ng	0.00
38) Acenaphthene-d10	15.02	164	516580	20.00	ng	0.00
63) Phenanthrene-d10	17.73	188	1036103	20.00	ng	0.00
75) Chrysene-d12	21.83	240	965347	20.00	ng	0.00
86) Perylene-d12	24.27	264	919644	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.89	112	1802257	129.15	ng	0.00
7) Phenol-d6	7.55	99	2364199	130.10	ng	0.00
23) Nitrobenzene-d5	9.59	82	1507605	86.12	ng	0.00
41) 2,4,6-Tribromophenol	16.49	330	903663	139.50	ng	0.00
44) 2-Fluorobiphenyl	13.65	172	3434051	82.64	ng	0.00
78) Terphenyl-d14	20.29	244	4256293	97.36	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.62	88	43659	7.573	ng	# 100
3) Pyridine	4.08	79	99661m	6.748	ng	
4) n-Nitrosodimethylamine	3.98	42	51790	6.481	ng	# 68
6) Aniline	7.72	93	171094	7.642	ng	95
8) 2-Chlorophenol	7.96	128	117160	7.439	ng	99
9) Benzaldehyde	7.53	77	68813	5.987	ng	96
10) Phenol	7.57	94	138112	7.484	ng	86
11) bis(2-Chloroethyl)ether	7.81	93	114391	7.815	ng	91
12) 1,3-Dichlorobenzene	8.29	146	143234	7.977	ng	96
13) 1,4-Dichlorobenzene	8.44	146	145173	7.996	ng	97
14) 1,2-Dichlorobenzene	8.76	146	139522	8.097	ng	95
15) Benzyl Alcohol	8.65	79	97000	7.382	ng	95
16) 2,2'-oxybis(1-Chloropropan	8.93	45	187392	8.115	ng	95
17) 2-Methylphenol	8.85	107	97415	7.941	ng	# 93
18) Hexachloroethane	9.50	117	48882	7.582	ng	93
19) n-Nitroso-di-n-propylamine	9.22	70	91455	7.896	ng	# 88
20) 3+4-Methylphenols	9.17	107	131805	8.092	ng	95
22) Acetophenone	9.25	105	189658	8.244	ng	# 96
24) Nitrobenzene	9.63	77	148367	8.139	ng	97
25) Isophorone	10.15	82	213726	7.108	ng	95
26) 2-Nitrophenol	10.35	139	51976	6.355	ng	# 76
27) 2,4-Dimethylphenol	10.39	122	87898	5.926	ng	91
28) bis(2-Chloroethoxy)methane	10.63	93	146996	7.466	ng	98
29) 2,4-Dichlorophenol	10.89	162	99454	7.081	ng	99
30) 1,2,4-Trichlorobenzene	11.10	180	134096	8.103	ng	98
31) Naphthalene	11.29	128	391711	7.859	ng	99
32) Benzoic acid	10.48	122	61695	7.190	ng	99
33) 4-Chloroaniline	11.40	127	152327	7.658	ng	# 93
34) Hexachlorobutadiene	11.56	225	82610	8.395	ng	99
35) Caprolactam	12.16	113	24488	6.160	ng	91
36) 4-Chloro-3-methylphenol	12.49	107	103360	7.203	ng	92
37) 2-Methylnaphthalene	12.87	142	267822	7.753	ng	98
39) 1,2,4,5-Tetrachlorobenzene	13.23	216	140619	7.728	ng	# 100
40) Hexachlorocyclopentadiene	13.20	237	49855	9.188	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.46	196	74776	6.838	ng	97
43) 2,4,5-Trichlorophenol	13.54	196	87197	7.385	ng	# 90
45) 1,1'-Biphenyl	13.86	154	370515	8.198	ng	98
46) 2-Chloronaphthalene	13.91	162	276179	7.939	ng	94
47) 2-Nitroaniline	14.11	65	64163	6.474	ng	88
48) Acenaphthylene	14.75	152	405431	7.847	ng	99
49) Dimethylphthalate	14.46	163	331563	8.042	ng	99
50) 2,6-Dinitrotoluene	14.59	165	63200	7.348	ng	90
51) Acenaphthene	15.08	154	258963	8.046	ng	98
52) 3-Nitroaniline	14.92	138	59354	6.554	ng	# 76
53) 2,4-Dinitrophenol	15.13	184	18360	9.825	ng	95
54) Dibenzofuran	15.41	168	391631	7.717	ng	95
55) 4-Nitrophenol	15.21	139	38919	5.298	ng	# 54
56) 2,4-Dinitrotoluene	15.37	165	75294	6.393	ng	# 82
57) Fluorene	16.05	166	311872	7.777	ng	100
58) 2,3,4,6-Tetrachlorophenol	15.63	232	67015	6.619	ng	# 100
59) Diethylphthalate	15.80	149	320272	7.881	ng	96
60) 4-Chlorophenyl-phenylether	16.03	204	159774	7.821	ng	95
61) 4-Nitroaniline	16.07	138	57499	6.072	ng	# 66
62) Azobenzene	16.33	77	289948	7.510	ng	91
64) 4,6-Dinitro-2-methylphenol	16.13	198	34341	6.005	ng	96
65) n-Nitrosodiphenylamine	16.24	169	262943	7.898	ng	98
66) 4-Bromophenyl-phenylether	16.92	248	91170	7.876	ng	# 89
67) Hexachlorobenzene	17.04	284	108126	8.111	ng	# 89
68) Atrazine	17.17	200	52775	4.997	ng	98
69) Pentachlorophenol	17.39	266	48776	6.573	ng	97
70) Phenanthrene	17.77	178	460604	7.805	ng	99
71) Anthracene	17.87	178	433922	7.376	ng	99
72) Carbazole	18.13	167	402120	7.721	ng	99
73) Di-n-butylphthalate	18.65	149	451806	7.368	ng	# 99
74) Fluoranthene	19.75	202	476488	7.362	ng	97
76) Benzidine	19.92	184	115366	3.817	ng	98
77) Pyrene	20.10	202	502198	8.243	ng	100
79) Butylbenzylphthalate	20.95	149	176513	7.095	ng	89
80) Benzo(a)anthracene	21.82	228	470686	7.737	ng	100
81) 3,3'-Dichlorobenzidine	21.74	252	161931	7.189	ng	99
82) Chrysene	21.87	228	457319	7.982	ng	99
83) Bis(2-ethylhexyl)phthalate	21.70	149	259297	7.166	ng	97
84) Di-n-octyl phthalate	22.63	149	387405	6.121	ng	# 94
85) Indeno(1,2,3-cd)pyrene	26.79	276	486763	7.022	ng	# 91
87) Benzo(b)fluoranthene	23.52	252	455687	7.831	ng	# 97
88) Benzo(k)fluoranthene	23.57	252	427120	7.760	ng	99
89) Benzo(a)pyrene	24.16	252	404044	7.444	ng	99
90) Dibenzo(a,h)anthracene	26.79	278	410176	7.408	ng	100
91) Benzo(g,h,i)perylene	27.57	276	396683	7.355	ng	97

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

