

Data Path : Z:\HPCHEM1\BNA N\DATA\BN020518\
 Data File : BN000019.D
 Acq On : 05 Feb 2018 15:16
 Operator : SJ/JU
 Sample : MDL-W-03
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MDL-W-03

Manual Integrations
 APPROVED

Sohil
 2/6/2018 5:15:44 PM

Quant Time: Feb 06 16:59:41 2018
 Quant Method : Z:\HPCHEM1\BNA N\METHODS\8270-BN020118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 01 19:48:39 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.48	152	220925	20.00	ng	-0.01
21) Naphthalene-d8	11.32	136	1089009	20.00	ng	-0.01
38) Acenaphthene-d10	15.09	164	656991	20.00	ng	0.00
63) Phenanthrene-d10	17.80	188	1557607	20.00	ng	0.00
75) Chrysene-d12	21.93	240	1971166	20.00	ng	-0.02
86) Perylene-d12	24.43	264	2126173	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.95	112	1857733	123.98	ng	0.00
7) Phenol-d6	7.59	99	2721423	119.57	ng	0.00
23) Nitrobenzene-d5	9.65	82	1639975	91.06	ng	-0.01
41) 2,4,6-Tribromophenol	16.55	330	835067	126.44	ng	0.00
44) 2-Fluorobiphenyl	13.72	172	4595659	93.68	ng	-0.01
78) Terphenyl-d14	20.37	244	6798680	91.04	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.66	88	28218	4.760	ng	96
3) Pyridine	4.13	79	70027	3.895	ng	95
4) n-Nitrosodimethylamine	4.02	42	21595	4.083	ng	# 85
6) Aniline	7.78	93	128877	4.154	ng	90
8) 2-Chlorophenol	8.03	128	73880	4.457	ng	92
9) Benzaldehyde	7.59	77	71752	3.545	ng	100
10) Phenol	7.62	94	103761	4.460	ng	92
11) bis(2-Chloroethyl)ether	7.88	93	87135	4.695	ng	87
12) 1,3-Dichlorobenzene	8.36	146	88173	4.791	ng	97
13) 1,4-Dichlorobenzene	8.51	146	92076	4.872	ng	92
14) 1,2-Dichlorobenzene	8.84	146	91180	4.883	ng	93
15) Benzyl Alcohol	8.70	79	59031	3.532	ng	# 88
16) 2,2'-oxybis(1-Chloropropan	9.01	45	91378	4.552	ng	77
17) 2-Methylphenol	8.90	107	63228	3.769	ng	96
18) Hexachloroethane	9.58	117	28274	4.363	ng	# 76
19) n-Nitroso-di-n-propylamine	9.28	70	56072	3.656	ng	# 84
20) 3+4-Methylphenols	9.23	107	84490	3.596	ng	94
22) Acetophenone	9.30	105	143581	4.464	ng	# 89
24) Nitrobenzene	9.69	77	101226	5.240	ng	# 94
25) Isophorone	10.22	82	155371	3.642	ng	97
26) 2-Nitrophenol	10.41	139	17796	2.635	ng	# 89
27) 2,4-Dimethylphenol	10.46	122	77231	4.415	ng	90
28) bis(2-Chloroethoxy)methane	10.70	93	107049	4.351	ng	97
29) 2,4-Dichlorophenol	10.95	162	46986	3.131	ng	90
30) 1,2,4-Trichlorobenzene	11.17	180	84547	4.811	ng	95
31) Naphthalene	11.37	128	296968	4.862	ng	99
32) Benzoic acid	10.48	122	18163m	8.245	ng	
33) 4-Chloroaniline	11.46	127	109481	3.866	ng	95
34) Hexachlorobutadiene	11.65	225	43793	4.891	ng	98
35) Caprolactam	12.18	113	16953	2.541	ng	97
36) 4-Chloro-3-methylphenol	12.54	107	61515	3.014	ng	99
37) 2-Methylnaphthalene	12.95	142	194614	4.509	ng	97
39) 1,2,4,5-Tetrachlorobenzene	13.30	216	92865	4.853	ng	# 95
40) Hexachlorocyclopentadiene	13.28	237	23872	3.735	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.52	196	32509	2.874	nq	99
43) 2,4,5-Trichlorophenol	13.59	196	38755	2.909	nq	# 92
45) 1,1'-Biphenyl	13.93	154	297454	5.067	nq	98
46) 2-Chloronaphthalene	13.98	162	211734	4.805	nq	96
47) 2-Nitroaniline	14.16	65	29909	2.512	nq	99
48) Acenaphthylene	14.81	152	303723	4.104	nq	97
49) Dimethylphthalate	14.52	163	248303	4.229	nq	99
50) 2,6-Dinitrotoluene	14.64	165	26665	2.295	nq	# 86
51) Acenaphthene	15.15	154	216474	4.636	nq	98
52) 3-Nitroaniline	14.97	138	35074	2.445	nq	92
53) 2,4-Dinitrophenol	15.16	184	6662	7.301	nq	# 1
54) Dibenzofuran	15.47	168	328874	4.908	nq	92
55) 4-Nitrophenol	15.25	139	22196	2.054	nq	# 83
56) 2,4-Dinitrotoluene	15.42	165	32422	2.072	nq	# 82
57) Fluorene	16.12	166	269722	4.780	nq	97
58) 2,3,4,6-Tetrachlorophenol	15.69	232	29442	2.714	nq	# 70
59) Diethylphthalate	15.87	149	240938	3.927	nq	94
60) 4-Chlorophenyl-phenylether	16.10	204	122279	4.903	nq	88
61) 4-Nitroaniline	16.12	138	38321	2.511	nq	96
62) Azobenzene	16.40	77	238550	3.891	nq	92
64) 4,6-Dinitro-2-methylphenol	16.17	198	11984	2.099	nq	88
65) n-Nitrosodiphenylamine	16.32	169	214243	4.209	nq	98
66) 4-Bromophenyl-phenylether	17.00	248	64878	4.451	nq	# 86
67) Hexachlorobenzene	17.11	284	83096	4.949	nq	# 83
68) Atrazine	17.24	200	55697	3.590	nq	92
69) Pentachlorophenol	17.45	266	26682	2.653	nq	97
70) Phenanthrene	17.85	178	469802	5.013	nq	98
71) Anthracene	17.94	178	405203	4.364	nq	97
72) Carbazole	18.20	167	411361	4.282	nq	96
73) Di-n-butylphthalate	18.74	149	321400	3.013	nq	# 97
74) Fluoranthene	19.83	202	477598	4.488	nq	94
76) Benzidine	20.00	184	142401	2.854	nq	96
77) Pyrene	20.19	202	543202	4.368	nq	99
79) Butylbenzylphthalate	21.05	149	99811	1.815	nq	# 88
80) Benzo(a)anthracene	21.92	228	563235	4.609	nq	97
81) 3,3'-Dichlorobenzidine	21.84	252	129698	3.167	nq	# 98
82) Chrysene	21.97	228	606095	4.994	nq	97
83) Bis(2-ethylhexyl)phthalate	21.83	149	182893	2.230	nq	# 91
84) Di-n-octyl phthalate	22.79	149	251275	1.893	nq	98
85) Indeno(1,2,3-cd)pyrene	27.00	276	674429	4.308	nq	# 86
87) Benzo(b)fluoranthene	23.67	252	555715	4.396	nq	# 96
88) Benzo(k)fluoranthene	23.72	252	579589	4.573	nq	# 96
89) Benzo(a)pyrene	24.32	252	512818	4.176	nq	# 95
90) Dibenzo(a,h)anthracene	27.02	278	567639	4.416	nq	# 91
91) Benzo(g,h,i)perylene	27.79	276	559860	4.285	nq	# 86

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

