

Data Path : Z:\HPCHEM1\BNA M\DATA\BM072617\
 Data File : BM010969.D
 Acq On : 26 Jul 2017 13:52
 Operator : SJ/JU
 Sample : SSTDICC060
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC060

Manual Integrations
 APPROVED

Sohil
 7/27/2017 6:21:27 PM

Quant Time: Jul 26 14:38:58 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM072617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 26 13:34:33 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	247764	20.00	ng	0.00
21) Naphthalene-d8	10.16	136	968498	20.00	ng	0.00
38) Acenaphthene-d10	14.05	164	652791	20.00	ng	0.00
63) Phenanthrene-d10	16.81	188	1604744	20.00	ng	0.00
75) Chrysene-d12	21.03	240	1753460	20.00	ng	0.00
86) Perylene-d12	23.14	264	1499570	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.03	112	1796558	120.11	ng	0.00
7) Phenol-d6	6.59	99	2537390	123.06	ng	0.00
23) Nitrobenzene-d5	8.55	82	3197899	126.92	ng	0.00
41) 2,4,6-Tribromophenol	15.56	330	1280417	128.35	ng	0.00
44) 2-Fluorobiphenyl	12.66	172	6362935	121.26	ng	0.00
78) Terphenyl-d14	19.47	244	10806721	119.20	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.03	88	379575	57.513	ng	# 100
3) Pyridine	3.41	79	1062746	58.910	ng	# 89
4) n-Nitrosodimethylamine	3.33	42	548035	59.440	ng	# 78
6) Aniline	6.75	93	1581696	61.234	ng	# 93
8) 2-Chlorophenol	6.97	128	891968	60.232	ng	# 85
9) Benzaldehyde	6.55	77	667071	46.591	ng	# 81
10) Phenol	6.62	94	1357106	62.249	ng	# 95
11) bis(2-Chloroethyl)ether	6.85	93	1036102	61.707	ng	# 95
12) 1,3-Dichlorobenzene	7.28	146	1085471	59.312	ng	# 85
13) 1,4-Dichlorobenzene	7.43	146	1115130	59.045	ng	# 90
14) 1,2-Dichlorobenzene	7.74	146	1043234	57.969	ng	# 91
15) Benzyl Alcohol	7.64	79	1182934	60.586	ng	# 84
16) 2,2'-oxybis(1-Chloropropan	7.92	45	919766	63.183	ng	# 77
17) 2-Methylphenol	7.85	107	868328	61.585	ng	# 81
18) Hexachloroethane	8.45	117	479177	60.490	ng	# 82
19) n-Nitroso-di-n-propylamine	8.21	70	1031165	63.131	ng	# 93
20) 3+4-Methylphenols	8.18	107	1198899	61.203	ng	# 84
22) Acetophenone	8.22	105	1763742	60.484	ng	# 95
24) Nitrobenzene	8.59	77	1635029	63.317	ng	# 88
25) Isophorone	9.11	82	2619317	63.456	ng	# 87
26) 2-Nitrophenol	9.29	139	545948	63.537	ng	# 30
27) 2,4-Dimethylphenol	9.36	122	922936	60.180	ng	# 74
28) bis(2-Chloroethoxy)methane	9.60	93	1439481	62.209	ng	# 98
29) 2,4-Dichlorophenol	9.82	162	1033733	59.670	ng	# 92
30) 1,2,4-Trichlorobenzene	10.02	180	1262208	59.598	ng	# 99
31) Naphthalene	10.20	128	3006401	61.050	ng	# 99
32) Benzoic acid	9.55	122	750887	62.403	ng	# 60
33) 4-Chloroaniline	10.33	127	1184882	59.721	ng	# 78
34) Hexachlorobutadiene	10.49	225	1011458	61.549	ng	# 97
35) Caprolactam	11.13	113	274931m	59.609	ng	#
36) 4-Chloro-3-methylphenol	11.46	107	1328357	63.390	ng	# 72
37) 2-Methylnaphthalene	11.83	142	2186711	61.822	ng	# 88
39) 1,2,4,5-Tetrachlorobenzene	12.21	216	1731905	60.866	ng	# 100
40) Hexachlorocyclopentadiene	12.19	237	1053616	64.592	ng	# 99

Data Path : Z:\HPCHEM1\BNA M\DATA\BM072617\
 Data File : BM010969.D
 Acq On : 26 Jul 2017 13:52
 Operator : SJ/JU
 Sample : SSTDICC060
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC060

Manual Integrations
 APPROVED

Sohil
 7/27/2017 6:21:27 PM

Quant Time: Jul 26 14:38:58 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM072617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 26 13:34:33 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	12.46	196	1054270	60.172	nq	99
43) 2,4,5-Trichlorophenol	12.53	196	1072360	61.992	nq #	88
45) 1,1'-Biphenyl	12.87	154	3461189	60.860	nq	97
46) 2-Chloronaphthalene	12.91	162	2518055	60.384	nq #	94
47) 2-Nitroaniline	13.13	65	925507	67.515	nq #	69
48) Acenaphthylene	13.77	152	3824474	61.664	nq	99
49) Dimethylphthalate	13.53	163	3315684	59.704	nq #	99
50) 2,6-Dinitrotoluene	13.64	165	692093	63.535	nq #	62
51) Acenaphthene	14.12	154	2353602	62.104	nq	99
52) 3-Nitroaniline	13.97	138	595904	60.176	nq #	40
53) 2,4-Dinitrophenol	14.18	184	504493	66.997	nq #	87
54) Dibenzofuran	14.46	168	3734151	60.913	nq #	90
55) 4-Nitrophenol	14.30	139	535103	62.218	nq #	1
56) 2,4-Dinitrotoluene	14.44	165	984597	65.139	nq	95
57) Fluorene	15.11	166	3270108	62.053	nq	99
58) 2,3,4,6-Tetrachlorophenol	14.69	232	987115	61.414	nq #	100
59) Diethylphthalate	14.91	149	3425547	62.457	nq	99
60) 4-Chlorophenyl-phenylether	15.12	204	1961038	60.579	nq	97
61) 4-Nitroaniline	15.15	138	650712	62.868	nq #	1
62) Azobenzene	15.41	77	3659230	65.703	nq	88
64) 4,6-Dinitro-2-methylphenol	15.20	198	702273	66.366	nq	96
65) n-Nitrosodiphenylamine	15.33	169	2853771	62.149	nq	98
66) 4-Bromophenyl-phenylether	16.01	248	1281866	63.030	nq #	87
67) Hexachlorobenzene	16.12	284	1420764	61.094	nq #	86
68) Atrazine	16.30	200	1132719	59.566	nq	96
69) Pentachlorophenol	16.46	266	933479	62.616	nq	98
70) Phenanthrene	16.85	178	5156030	61.761	nq	100
71) Anthracene	16.94	178	5171844	62.410	nq	100
72) Carbazole	17.22	167	4512559	61.830	nq	99
73) Di-n-butylphthalate	17.80	149	5671208	65.679	nq #	97
74) Fluoranthene	18.89	202	6388880	61.792	nq	98
76) Benzidine	19.09	184	2497923	53.235	nq	99
77) Pyrene	19.25	202	6378165	62.746	nq	99
79) Butylbenzylphthalate	20.19	149	2437094	68.480	nq #	72
80) Benzo(a)anthracene	21.02	228	6061817	59.662	nq	99
81) 3,3'-Dichlorobenzidine	20.96	252	2463570	61.666	nq #	97
82) Chrysene	21.07	228	5831355	60.264	nq	100
83) Bis(2-ethylhexyl)phthalate	20.97	149	3575864	68.999	nq	99
84) Di-n-octyl phthalate	21.83	149	5687270	66.978	nq	100
85) Indeno(1,2,3-cd)pyrene	25.23	276	6013682	54.198	nq #	100
87) Benzo(b)fluoranthene	22.52	252	5867067	61.594	nq #	92
88) Benzo(k)fluoranthene	22.56	252	5740030	63.090	nq #	96
89) Benzo(a)pyrene	23.05	252	5330716	61.483	nq #	95
90) Dibenzo(a,h)anthracene	25.24	278	4901944	58.252	nq #	91
91) Benzo(g,h,i)perylene	25.86	276	4808935	58.134	nq #	85

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

