

Data Path : Z:\HPCHEM1\BNA M\DATA\BM051117\
 Data File : BM009986.D
 Acq On : 11 May 2017 16:06
 Operator : SJ/MA
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

mohammad
 5/12/2017 4:30:00 PM

Quant Time: May 11 19:02:45 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM051117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 11 18:53:14 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	59517	20.00	ng	0.00
21) Naphthalene-d8	10.56	136	253203	20.00	ng	0.00
38) Acenaphthene-d10	14.42	164	167297	20.00	ng	0.00
63) Phenanthrene-d10	17.17	188	425991	20.00	ng	0.00
75) Chrysene-d12	21.36	240	575958	20.00	ng	0.00
86) Perylene-d12	23.66	264	458278	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.36	112	75810	20.28	ng	0.00
7) Phenol-d6	6.95	99	112197	19.58	ng	0.00
23) Nitrobenzene-d5	8.93	82	131660	18.82	ng	0.00
41) 2,4,6-Tribromophenol	15.92	330	42714	18.69	ng	0.00
44) 2-Fluorobiphenyl	13.03	172	244768	19.06	ng	0.00
78) Terphenyl-d14	19.80	244	488364	17.39	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.28	88	13528	9.42	ng	# 100
3) Pyridine	3.69	79	47070	9.53	ng	# 93
4) n-Nitrosodimethylamine	3.60	42	27150	8.43	ng	# 54
6) Aniline	7.10	93	74320	9.81	ng	# 87
8) 2-Chlorophenol	7.33	128	37965	9.12	ng	# 64
9) Benzaldehyde	6.93	77	46238	10.16	ng	90
10) Phenol	6.97	94	62175	9.91	ng	91
11) bis(2-Chloroethyl)ether	7.20	93	50625	10.14	ng	90
12) 1,3-Dichlorobenzene	7.65	146	41430	8.98	ng	# 80
13) 1,4-Dichlorobenzene	7.80	146	44188	9.29	ng	# 94
14) 1,2-Dichlorobenzene	8.11	146	42317	9.26	ng	# 92
15) Benzyl Alcohol	8.02	79	44524	8.39	ng	# 80
16) 2,2'-oxybis(1-Chloropropan	8.29	45	77417	8.39	ng	94
17) 2-Methylphenol	8.22	107	40270	9.53	ng	# 89
18) Hexachloroethane	8.83	117	18526	8.79	ng	98
19) n-Nitroso-di-n-propylamine	8.57	70	47539	8.75	ng	# 88
20) 3+4-Methylphenols	8.55	107	52120	8.98	ng	92
22) Acetophenone	8.60	105	76719	9.68	ng	# 88
24) Nitrobenzene	8.98	77	73136	10.07	ng	# 88
25) Isophorone	9.49	82	112633	9.13	ng	# 84
26) 2-Nitrophenol	9.69	139	19622	8.35	ng	# 27
27) 2,4-Dimethylphenol	9.75	122	39657	9.38	ng	# 82
28) bis(2-Chloroethoxy)methane	9.97	93	70778	9.85	ng	94
29) 2,4-Dichlorophenol	10.23	162	38190	9.42	ng	95
30) 1,2,4-Trichlorobenzene	10.42	180	44863	9.47	ng	96
31) Naphthalene	10.61	128	135171	9.71	ng	99
32) Benzoic acid	9.87	122	20908m	7.53	ng	
33) 4-Chloroaniline	10.75	127	55269	9.41	ng	# 76
34) Hexachlorobutadiene	10.87	225	31056	9.42	ng	95
35) Caprolactam	11.53	113	13925	9.01	ng	# 91
36) 4-Chloro-3-methylphenol	11.87	107	50763	9.22	ng	86
37) 2-Methylnaphthalene	12.22	142	95857	9.75	ng	# 87
39) 1,2,4,5-Tetrachlorobenzene	12.60	216	56229	9.34	ng	# 100
40) Hexachlorocyclopentadiene	12.56	237	1288	1.05	ng	94

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	12.86	196	36419	9.36	nq	97
43) 2,4,5-Trichlorophenol	12.94	196	38157	9.12	nq	# 87
45) 1,1'-Biphenyl	13.25	154	136513	9.58	nq	90
46) 2-Chloronaphthalene	13.29	162	105756	9.39	nq	97
47) 2-Nitroaniline	13.52	65	46208	8.50	nq	# 68
48) Acenaphthylene	14.14	152	165252	9.35	nq	99
49) Dimethylphthalate	13.87	163	144668	9.08	nq	98
50) 2,6-Dinitrotoluene	14.01	165	28761	9.25	nq	# 60
51) Acenaphthene	14.48	154	108935	9.99	nq	95
52) 3-Nitroaniline	14.36	138	29047	8.64	nq	# 48
53) 2,4-Dinitrophenol	14.60	184	5325	3.72	nq	94
54) Dibenzofuran	14.82	168	162356	9.74	nq	96
55) 4-Nitrophenol	14.70	139	17542	7.76	nq	# 11
56) 2,4-Dinitrotoluene	14.82	165	38904	8.61	nq	# 65
57) Fluorene	15.47	166	136241	9.26	nq	98
58) 2,3,4,6-Tetrachlorophenol	15.06	232	31761	9.91	nq	# 100
59) Diethylphthalate	15.24	149	149032	8.75	nq	97
60) 4-Chlorophenyl-phenylether	15.46	204	75943	9.73	nq	99
61) 4-Nitroaniline	15.53	138	29890	8.87	nq	# 15
62) Azobenzene	15.76	77	181066	9.55	nq	85
64) 4,6-Dinitro-2-methylphenol	15.59	198	17280	6.53	nq	97
65) n-Nitrosodiphenylamine	15.68	169	127872	9.47	nq	99
66) 4-Bromophenyl-phenylether	16.36	248	46793	9.39	nq	# 92
67) Hexachlorobenzene	16.47	284	52956	9.33	nq	# 86
68) Atrazine	16.63	200	44764	8.66	nq	94
69) Pentachlorophenol	16.83	266	17362	8.08	nq	90
70) Phenanthrene	17.22	178	231600	9.39	nq	98
71) Anthracene	17.31	178	233952	9.41	nq	99
72) Carbazole	17.59	167	214886	9.53	nq	94
73) Di-n-butylphthalate	18.13	149	262285	8.10	nq	# 97
74) Fluoranthene	19.23	202	304419	9.78	nq	97
76) Benzidine	19.43	184	133863	7.70	nq	97
77) Pyrene	19.60	202	314844	9.15	nq	99
79) Butylbenzylphthalate	20.49	149	124085	7.48	nq	# 66
80) Benzo(a)anthracene	21.34	228	322007	9.28	nq	99
81) 3,3'-Dichlorobenzidine	21.29	252	115605	8.39	nq	98
82) Chrysene	21.40	228	307816	9.44	nq	98
83) Bis(2-ethylhexyl)phthalate	21.25	149	191104	7.49	nq	96
84) Di-n-octyl phthalate	22.14	149	329083	7.43	nq	# 86
85) Indeno(1,2,3-cd)pyrene	25.99	276	249308	6.78	nq	# 100
87) Benzo(b)fluoranthene	22.96	252	293778	10.06	nq	# 93
88) Benzo(k)fluoranthene	23.01	252	276860	9.75	nq	98
89) Benzo(a)pyrene	23.56	252	255843	9.34	nq	98
90) Dibenzo(a,h)anthracene	26.00	278	222669	8.10	nq	97
91) Benzo(g,h,i)perylene	26.71	276	209396	8.12	nq	# 91

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

