

Data Path : Z:\HPCHEM1\BNA F\DATA\BF021915\
 Data File : BF077371.D
 Acq On : 19 Feb 2015 13:20
 Operator : TP/IZ
 Sample : SSTDICC060
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC060

Manual Integrations
 APPROVED

apatel
 2/20/2015 2:04:06 PM

Quant Time: Feb 19 15:10:43 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF021915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 19 14:49:45 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.33	152	76018	20.00	ng	0.00
21) Naphthalene-d8	8.91	136	317467	20.00	ng	0.00
38) Acenaphthene-d10	11.09	164	189636	20.00	ng	0.00
63) Phenanthrene-d10	12.93	188	327427	20.00	ng	0.00
75) Chrysene-d12	16.26	240	332663	20.00	ng	0.03
86) Perylene-d12	18.09	264	324157	20.00	ng	0.14

System Monitoring Compounds

5) 2-Fluorophenol	5.62	112	523940	117.47	ng	0.00
7) Phenol-d6	6.89	99	691228	119.17	ng	0.01
23) Nitrobenzene-d5	8.03	82	627127	134.30	ng	0.00
41) 2,4,6-Tribromophenol	12.08	330	219811	132.74	ng	0.01
44) 2-Fluorobiphenyl	10.26	172	1279820	102.94	ng	0.00
78) Terphenyl-d14	14.93	244	1610704	111.21	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.29	88	107882	56.08	ng	98
3) Pyridine	3.02	79	304728	58.76	ng	96
4) n-Nitrosodimethylamine	2.96	42	130304	52.83	ng	97
6) Aniline	6.92	93	475817	58.04	ng	97
8) 2-Chlorophenol	7.06	128	310435	59.43	ng	99
9) Benzaldehyde	6.78	77	190367	53.29	ng	98
10) Phenol	6.90	94	417553	61.27	ng	97
11) bis(2-Chloroethyl)ether	7.02	93	302388	60.09	ng	98
12) 1,3-Dichlorobenzene	7.26	146	338190	57.60	ng	# 88
13) 1,4-Dichlorobenzene	7.36	146	342362	56.55	ng	98
14) 1,2-Dichlorobenzene	7.54	146	331022	56.76	ng	99
15) Benzyl Alcohol	7.52	79	264865	60.26	ng	91
16) 2,2'-oxybis(1-Chloropropan	7.70	45	424069	56.24	ng	100
17) 2-Methylphenol	7.65	107	251386	60.02	ng	98
18) Hexachloroethane	7.96	117	118321	59.98	ng	96
19) n-Nitroso-di-n-propylamine	7.85	70	212822	53.48	ng	96
20) 3+4-Methylphenols	7.85	107	331925	61.17	ng	95
22) Acetophenone	7.85	105	422829	57.13	ng	# 89
24) Nitrobenzene	8.05	77	326739	65.49	ng	96
25) Isophorone	8.35	82	617224	59.68	ng	98
26) 2-Nitrophenol	8.44	139	143083	98.52	ng	99
27) 2,4-Dimethylphenol	8.50	122	285076	60.18	ng	98
28) bis(2-Chloroethoxy)methane	8.62	93	363963	55.35	ng	100
29) 2,4-Dichlorophenol	8.74	162	282762	68.66	ng	98
30) 1,2,4-Trichlorobenzene	8.84	180	287900	56.16	ng	99
31) Naphthalene	8.94	128	920218	57.90	ng	99
32) Benzoic acid	8.61	122	152998m	109.20	ng	
33) 4-Chloroaniline	9.01	127	429845	62.56	ng	98
34) Hexachlorobutadiene	9.09	225	166707	59.20	ng	99
35) Caprolactam	9.45	113	84303	66.32	ng	96
36) 4-Chloro-3-methylphenol	9.61	107	269324	59.74	ng	95
37) 2-Methylnaphthalene	9.80	142	675784	58.87	ng	99
39) 1,2,4,5-Tetrachlorobenzene	10.01	216	313913	57.23	ng	100
40) Hexachlorocyclopentadiene	10.00	237	185968	67.68	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	10.14	196	207945	66.65	nq	99
43) 2,4,5-Trichlorophenol	10.19	196	229619	68.42	nq	97
45) 1,1'-Biphenyl	10.38	154	811660	53.93	nq	98
46) 2-Chloronaphthalene	10.41	162	633165	56.41	nq	94
47) 2-Nitroaniline	10.53	65	168131	75.28	nq	# 61
48) Acenaphthylene	10.92	152	1006501	54.79	nq	99
49) Dimethylphthalate	10.77	163	767421	58.68	nq	# 98
50) 2,6-Dinitrotoluene	10.84	165	168080	75.22	nq	# 83
51) Acenaphthene	11.13	154	565106	51.62	nq	100
52) 3-Nitroaniline	11.04	138	175905	66.38	nq	92
53) 2,4-Dinitrophenol	11.17	184	55521	135.86	nq	93
54) Dibenzofuran	11.34	168	900877	54.10	nq	98
55) 4-Nitrophenol	11.24	139	158257	85.63	nq	85
56) 2,4-Dinitrotoluene	11.33	165	209233	77.69	nq	89
57) Fluorene	11.77	166	717047	54.74	nq	99
58) 2,3,4,6-Tetrachlorophenol	11.49	232	186141	73.77	nq	99
59) Diethylphthalate	11.65	149	772336	57.29	nq	96
60) 4-Chlorophenyl-phenylether	11.78	204	344974	54.37	nq	94
61) 4-Nitroaniline	11.80	138	188074	73.76	nq	87
62) Azobenzene	11.97	77	691634	51.97	nq	96
64) 4,6-Dinitro-2-methylphenol	11.84	198	86006	131.90	nq	82
65) n-Nitrosodiphenylamine	11.93	169	647649	60.78	nq	99
66) 4-Bromophenyl-phenylether	12.39	248	224162	61.52	nq	93
67) Hexachlorobenzene	12.45	284	238707	58.26	nq	# 76
68) Atrazine	12.60	200	223480	69.50	nq	94
69) Pentachlorophenol	12.69	266	139908	87.81	nq	98
70) Phenanthrene	12.97	178	1096369	58.10	nq	98
71) Anthracene	13.03	178	1088280	57.99	nq	99
72) Carbazole	13.23	167	985503	58.91	nq	99
73) Di-n-butylphthalate	13.69	149	1296454	63.36	nq	# 98
74) Fluoranthene	14.44	202	1152914	59.54	nq	96
76) Benzidine	14.63	184	629185	62.02	nq	99
77) Pyrene	14.73	202	1214661	59.80	nq	99
79) Butylbenzylphthalate	15.56	149	551738	73.65	nq	96
80) Benzo(a)anthracene	16.25	228	1168397	60.61	nq	98
81) 3,3'-Dichlorobenzidine	16.23	252	436072	67.65	nq	99
82) Chrysene	16.29	228	1047923	57.59	nq	100
83) Bis(2-ethylhexyl)phthalate	16.31	149	841057	65.92	nq	# 98
84) Di-n-octyl phthalate	17.14	149	1394414m	69.89	nq	
85) Indeno(1,2,3-cd)pyrene	19.61	276	1228442m	59.44	nq	
87) Benzo(b)fluoranthene	17.62	252	1110585m	53.51	nq	
88) Benzo(k)fluoranthene	17.65	252	1067002	63.00	nq	# 96
89) Benzo(a)pyrene	18.02	252	1030772	57.05	nq	# 98
90) Dibenzo(a,h)anthracene	19.64	278	997310	55.29	nq	97
91) Benzo(g,h,i)perylene	20.05	276	1011911	56.07	nq	97

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

