

Data Path : Z:\HPCHEM1\BNA F\DATA\BF110715\
 Data File : BF082755.D
 Acq On : 6 Nov 2015 12:00
 Operator : UM/IZ
 Sample : SSTDICC025
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC025

Manual Integrations
 APPROVED

Mmdadoda
 11/9/2015 2:10:02 PM

Quant Time: Nov 06 13:06:53 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 06 12:40:46 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	175199	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	658795	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	362086	20.00	ng	0.00
63) Phenanthrene-d10	11.38	188	665188	20.00	ng	0.00
75) Chrysene-d12	14.02	240	592225	20.00	ng	0.00
86) Perylene-d12	15.46	264	538447	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.45	112	589150	50.65	ng	0.00
7) Phenol-d6	6.48	99	733312	47.45	ng	0.00
23) Nitrobenzene-d5	7.41	82	763197	47.31	ng	0.00
41) 2,4,6-Tribromophenol	10.68	330	225278	56.85	ng	0.00
44) 2-Fluorobiphenyl	9.22	172	1383006	54.23	ng	0.00
78) Terphenyl-d14	12.97	244	1343141	49.63	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.44	88	121010	22.50	ng	96
3) Pyridine	3.16	79	364529	23.12	ng	98
4) n-Nitrosodimethylamine	3.09	42	171746	19.35	ng	86
6) Aniline	6.51	93	539294	25.06	ng	97
8) 2-Chlorophenol	6.64	128	328082	26.07	ng	97
9) Benzaldehyde	6.40	77	230906	21.89	ng	82
10) Phenol	6.49	94	401799	22.95	ng	98
11) bis(2-Chloroethyl)ether	6.59	93	303725	23.51	ng	# 78
12) 1,3-Dichlorobenzene	6.80	146	354609m	24.71	ng	
13) 1,4-Dichlorobenzene	6.86	146	346685	24.36	ng	98
14) 1,2-Dichlorobenzene	7.02	146	367787	27.84	ng	98
15) Benzyl Alcohol	6.99	79	327474	22.72	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.14	45	495063	24.11	ng	96
17) 2-Methylphenol	7.10	107	257353	24.05	ng	98
18) Hexachloroethane	7.37	117	132943	23.93	ng	97
19) n-Nitroso-di-n-propylamine	7.26	70	263617	22.50	ng	93
20) 3+4-Methylphenols	7.26	107	373691	26.87	ng	# 84
22) Acetophenone	7.26	105	513082	26.45	ng	95
24) Nitrobenzene	7.44	77	424636	26.04	ng	99
25) Isophorone	7.68	82	732454	24.68	ng	99
26) 2-Nitrophenol	7.76	139	171773	28.84	ng	# 60
27) 2,4-Dimethylphenol	7.79	122	278665	23.96	ng	98
28) bis(2-Chloroethoxy)methane	7.89	93	429865	26.48	ng	99
29) 2,4-Dichlorophenol	8.00	162	291396	27.49	ng	96
30) 1,2,4-Trichlorobenzene	8.09	180	301088	25.60	ng	# 94
31) Naphthalene	8.16	128	880652	24.70	ng	99
32) Benzoic acid	7.90	122	205479	24.36	ng	97
33) 4-Chloroaniline	8.20	127	381444	25.06	ng	97
34) Hexachlorobutadiene	8.29	225	192619	25.75	ng	98
35) Caprolactam	8.57	113	81168	24.96	ng	88
36) 4-Chloro-3-methylphenol	8.69	107	323608	25.93	ng	# 78
37) 2-Methylnaphthalene	8.85	142	652164	28.23	ng	97
39) 1,2,4,5-Tetrachlorobenzene	9.02	216	316963	26.23	ng	97
40) Hexachlorocyclopentadiene	9.01	237	169562	22.32	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.13	196	220085	25.23	nq	97
43) 2,4,5-Trichlorophenol	9.16	196	205006	23.58	nq	98
45) 1,1'-Biphenyl	9.32	154	766902	25.07	nq	96
46) 2-Chloronaphthalene	9.34	162	600275	25.13	nq	94
47) 2-Nitroaniline	9.42	65	206516	22.25	nq	95
48) Acenaphthylene	9.76	152	1017612	26.64	nq	98
49) Dimethylphthalate	9.62	163	773969	26.17	nq	99
50) 2,6-Dinitrotoluene	9.68	165	175748	28.67	nq	# 79
51) Acenaphthene	9.93	154	646353	26.65	nq	99
52) 3-Nitroaniline	9.84	138	158829	23.29	nq	93
53) 2,4-Dinitrophenol	9.94	184	76158	27.74	nq	# 50
54) Dibenzofuran	10.10	168	883129	26.84	nq	91
55) 4-Nitrophenol	10.00	139	137154	26.73	nq	# 86
56) 2,4-Dinitrotoluene	10.08	165	236834	28.57	nq	# 84
57) Fluorene	10.44	166	728160	27.43	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.21	232	177827	24.21	nq	95
59) Diethylphthalate	10.32	149	726067	24.12	nq	98
60) 4-Chlorophenyl-phenylether	10.43	204	311962	24.18	nq	# 82
61) 4-Nitroaniline	10.45	138	192144	29.53	nq	# 81
62) Azobenzene	10.59	77	699745	20.61	nq	90
64) 4,6-Dinitro-2-methylphenol	10.49	198	132445	33.01	nq	79
65) n-Nitrosodiphenylamine	10.56	169	597956	26.15	nq	99
66) 4-Bromophenyl-phenylether	10.92	248	205477	27.26	nq	# 82
67) Hexachlorobenzene	10.99	284	227163	27.18	nq	# 78
68) Atrazine	11.08	200	212229	28.37	nq	94
69) Pentachlorophenol	11.18	266	143707	26.07	nq	99
70) Phenanthrene	11.40	178	1069662	28.51	nq	97
71) Anthracene	11.45	178	982503	26.01	nq	99
72) Carbazole	11.61	167	891932	27.11	nq	# 96
73) Di-n-butylphthalate	11.95	149	1132772	27.04	nq	# 96
74) Fluoranthene	12.59	202	1052523	27.44	nq	93
76) Benzidine	12.70	184	621463	23.27	nq	99
77) Pyrene	12.81	202	1029520	23.48	nq	98
79) Butylbenzylphthalate	13.44	149	466256	24.27	nq	96
80) Benzo(a)anthracene	14.01	228	997299	26.51	nq	98
81) 3,3'-Dichlorobenzidine	13.96	252	339438	26.19	nq	98
82) Chrysene	14.04	228	963315	28.25	nq	99
83) Bis(2-ethylhexyl)phthalate	14.01	149	673238	27.04	nq	99
84) Di-n-octyl phthalate	14.64	149	1118950	28.81	nq	96
85) Indeno(1,2,3-cd)pyrene	16.85	276	777608	27.63	nq	100
87) Benzo(b)fluoranthene	15.05	252	816947	22.00	nq	# 99
88) Benzo(k)fluoranthene	15.08	252	873848m	30.89	nq	
89) Benzo(a)pyrene	15.40	252	777068	26.58	nq	# 97
90) Dibenzo(a,h)anthracene	16.87	278	656579	24.07	nq	99
91) Benzo(g,h,i)perylene	17.27	276	608847	23.03	nq	98

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

