

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052316\
 Data File : BF087303.D
 Acq On : 23 May 2016 16:35
 Operator : UM/SJ
 Sample : SSTDICC060
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC060

Manual Integrations
 APPROVED

umangi
 5/24/2016 7:39:03 PM

Quant Time: May 24 01:13:48 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 24 00:59:18 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	145802	20.00	ng	0.01
21) Naphthalene-d8	9.64	136	603562	20.00	ng	0.00
38) Acenaphthene-d10	12.48	164	272197	20.00	ng	0.00
63) Phenanthrene-d10	14.88	188	505617	20.00	ng	0.00
75) Chrysene-d12	18.56	240	354606	20.00	ng	0.00
86) Perylene-d12	20.23	264	286781	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.52	112	1059199	121.76	ng	0.00
7) Phenol-d6	7.16	99	1379072	119.34	ng	0.00
23) Nitrobenzene-d5	8.55	82	1432269	117.33	ng	0.01
41) 2,4,6-Tribromophenol	13.78	330	313155	106.82	ng	0.00
44) 2-Fluorobiphenyl	11.44	172	2235251	133.65	ng	0.01
78) Terphenyl-d14	17.22	244	1956521	121.44	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.94	88	254624	57.57	ng	# 73
3) Pyridine	2.55	79	609486	57.64	ng	# 65
4) n-Nitrosodimethylamine	2.54	42	462735	61.35	ng	# 46
6) Aniline	7.12	93	886992	59.93	ng	96
8) 2-Chlorophenol	7.30	128	614800	58.47	ng	94
9) Benzaldehyde	6.94	77	306140	46.61	ng	98
10) Phenol	7.19	94	702952	50.03	ng	81
11) bis(2-Chloroethyl)ether	7.27	93	584854	53.30	ng	90
12) 1,3-Dichlorobenzene	7.52	146	676289	60.56	ng	# 94
13) 1,4-Dichlorobenzene	7.64	146	681411	59.44	ng	94
14) 1,2-Dichlorobenzene	7.88	146	626317	61.33	ng	98
15) Benzyl Alcohol	7.91	79	526898	61.21	ng	89
16) 2,2'-oxybis(1-Chloropropan	8.10	45	1296903	57.35	ng	88
17) 2-Methylphenol	8.12	107	489811	57.03	ng	# 86
18) Hexachloroethane	8.40	117	255753	58.65	ng	# 87
19) n-Nitroso-di-n-propylamine	8.34	70	513022	59.01	ng	# 69
20) 3+4-Methylphenols	8.39	107	616283	54.83	ng	# 59
22) Acetophenone	8.31	105	882386	60.03	ng	# 96
24) Nitrobenzene	8.57	77	726811	55.43	ng	97
25) Isophorone	8.97	82	1290212	58.33	ng	97
26) 2-Nitrophenol	9.07	139	330479	60.06	ng	# 86
27) 2,4-Dimethylphenol	9.21	122	533321	57.03	ng	87
28) bis(2-Chloroethoxy)methane	9.34	93	701264	58.34	ng	98
29) 2,4-Dichlorophenol	9.47	162	492582	59.35	ng	94
30) 1,2,4-Trichlorobenzene	9.58	180	544125	58.89	ng	98
31) Naphthalene	9.69	128	1759593	59.32	ng	100
32) Benzoic acid	9.56	122	281443	52.20	ng	# 77
33) 4-Chloroaniline	9.83	127	656138	54.30	ng	# 94
34) Hexachlorobutadiene	9.90	225	333675	62.95	ng	99
35) Caprolactam	10.50	113	148363m	56.72	ng	
36) 4-Chloro-3-methylphenol	10.68	107	541639	54.52	ng	90
37) 2-Methylnaphthalene	10.81	142	1072035	56.46	ng	96
39) 1,2,4,5-Tetrachlorobenzene	11.08	216	505853	62.09	ng	98
40) Hexachlorocyclopentadiene	11.06	237	200719	79.46	ng	94

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	11.30	196	308044	59.64	ng	96
43) 2,4,5-Trichlorophenol	11.38	196	312624	58.79	ng #	88
45) 1,1'-Biphenyl	11.59	154	1364296	60.09	ng	98
46) 2-Chloronaphthalene	11.60	162	985971	56.39	ng	94
47) 2-Nitroaniline	11.82	65	406563	61.68	ng	85
48) Acenaphthylene	12.25	152	1645021	61.13	ng	99
49) Dimethylphthalate	12.15	163	1293821	61.74	ng	99
50) 2,6-Dinitrotoluene	12.23	165	261436	58.79	ng #	75
51) Acenaphthene	12.54	154	991263	59.09	ng	99
52) 3-Nitroaniline	12.50	138	293858	61.03	ng	96
53) 2,4-Dinitrophenol	12.70	184	133506	60.98	ng	86
54) Dibenzofuran	12.82	168	1298781	59.52	ng	94
55) 4-Nitrophenol	12.87	139	134719	65.36	ng #	66
56) 2,4-Dinitrotoluene	12.89	165	355687	61.41	ng	93
57) Fluorene	13.38	166	1153161	64.48	ng	98
58) 2,3,4,6-Tetrachlorophenol	13.06	232	241471	58.06	ng	98
59) Diethylphthalate	13.29	149	1259696	63.70	ng	97
60) 4-Chlorophenyl-phenylether	13.41	204	555675	67.11	ng	93
61) 4-Nitroaniline	13.51	138	257382	54.84	ng #	68
62) Azobenzene	13.66	77	1347866	59.19	ng	86
64) 4,6-Dinitro-2-methylphenol	13.55	198	203623	61.19	ng #	55
65) n-Nitrosodiphenylamine	13.62	169	940963	59.11	ng	99
66) 4-Bromophenyl-phenylether	14.19	248	324785	61.02	ng	91
67) Hexachlorobenzene	14.26	284	336461	56.55	ng #	74
68) Atrazine	14.54	200	245753	47.45	ng	93
69) Pentachlorophenol	14.62	266	100568	47.00	ng	96
70) Phenanthrene	14.93	178	1549316	56.71	ng	100
71) Anthracene	15.01	178	1585214	56.13	ng	100
72) Carbazole	15.29	167	1353230	53.97	ng	98
73) Di-n-butylphthalate	15.86	149	1806744	61.51	ng #	99
74) Fluoranthene	16.67	202	1609745	58.01	ng	94
76) Benzidine	16.89	184	548246	58.70	ng	98
77) Pyrene	16.97	202	1544112	60.42	ng	99
79) Butylbenzylphthalate	17.89	149	772693	71.19	ng	89
80) Benzo(a)anthracene	18.55	228	1219393	59.07	ng	99
81) 3,3'-Dichlorobenzidine	18.54	252	669704	52.32	ng	99
82) Chrysene	18.59	228	1219178	61.43	ng	99
83) Bis(2-ethylhexyl)phthalate	18.62	149	1053541	77.08	ng #	96
84) Di-n-octyl phthalate	19.39	149	1648308	70.33	ng #	84
85) Indeno(1,2,3-cd)pyrene	21.50	276	1029453	59.55	ng	94
87) Benzo(b)fluoranthene	19.82	252	1158945m	59.17	ng	
88) Benzo(k)fluoranthene	19.85	252	867857m	63.33	ng	
89) Benzo(a)pyrene	20.17	252	948783	59.94	ng	98
90) Dibenzo(a,h)anthracene	21.50	278	858296	63.93	ng	99
91) Benzo(g,h,i)perylene	21.86	276	862464	64.05	ng #	90

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

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