

Data Path : Z:\HPCHEM1\BNA F\DATA\BF080715\
 Data File : BF080918.D
 Acq On : 7 Aug 2015 10:42
 Operator : TP/UM
 Sample : SSTDICC060
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC060

Manual Integrations
 APPROVED

MMDadoda
 8/10/2015 7:17:15 PM

Quant Time: Aug 08 00:23:42 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF080715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 07 23:54:48 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	56489	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	236974	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	107663	20.00	ng	0.00
63) Phenanthrene-d10	11.33	188	227400	20.00	ng	0.01
75) Chrysene-d12	13.96	240	191001	20.00	ng	0.00
86) Perylene-d12	15.36	264	163985	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	380211	109.97	ng	0.01
7) Phenol-d6	6.44	99	531037	112.09	ng	0.00
23) Nitrobenzene-d5	7.38	82	509088	101.62	ng	0.00
41) 2,4,6-Tribromophenol	10.64	330	124137	104.95	ng	0.00
44) 2-Fluorobiphenyl	9.17	172	787385	102.96	ng	0.00
78) Terphenyl-d14	12.91	244	967214	104.60	ng	0.00

Target Compounds						Qvalue
2) 1,4-Dioxane	2.33	88	92056	56.14	ng	100
3) Pyridine	3.04	79	272889	58.28	ng	98
4) n-Nitrosodimethylamine	3.00	42	144516	60.88	ng	93
6) Aniline	6.48	93	392689	56.42	ng	97
8) 2-Chlorophenol	6.59	128	212685	53.16	ng	85
9) Benzaldehyde	6.35	77	165994	51.71	ng	98
10) Phenol	6.45	94	294722	52.43	ng	81
11) bis(2-Chloroethyl)ether	6.56	93	242301	57.73	ng	88
12) 1,3-Dichlorobenzene	6.75	146	239783m	56.48	ng	
13) 1,4-Dichlorobenzene	6.83	146	258107	58.50	ng	96
14) 1,2-Dichlorobenzene	6.98	146	201408	50.50	ng	96
15) Benzyl Alcohol	6.96	79	208148	53.69	ng	95
16) 2,2'-oxybis(1-Chloropropan	7.09	45	457433	54.17	ng	99
17) 2-Methylphenol	7.07	107	184096	53.91	ng	93
18) Hexachloroethane	7.32	117	94081	55.66	ng	97
19) n-Nitroso-di-n-propylamine	7.24	70	209230	60.42	ng	97
20) 3+4-Methylphenols	7.23	107	244909	57.09	ng	96
22) Acetophenone	7.23	105	321672	55.40	ng	# 91
24) Nitrobenzene	7.40	77	308742	59.85	ng	92
25) Isophorone	7.64	82	524752	54.47	ng	94
26) 2-Nitrophenol	7.71	139	113738	52.40	ng	92
27) 2,4-Dimethylphenol	7.76	122	216033	53.16	ng	95
28) bis(2-Chloroethoxy)methane	7.86	93	322167	55.70	ng	# 96
29) 2,4-Dichlorophenol	7.96	162	185327	55.50	ng	89
30) 1,2,4-Trichlorobenzene	8.04	180	204019	55.90	ng	98
31) Naphthalene	8.12	128	715201	54.75	ng	98
32) Benzoic acid	7.89	122	150211	61.46	ng	92
33) 4-Chloroaniline	8.17	127	261485	49.26	ng	96
34) Hexachlorobutadiene	8.24	225	112229	51.68	ng	96
35) Caprolactam	8.56	113	67919	58.23	ng	# 72
36) 4-Chloro-3-methylphenol	8.66	107	240781	59.09	ng	90
37) 2-Methylnaphthalene	8.81	142	406655	48.91	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.98	216	201158	60.61	ng	97
40) Hexachlorocyclopentadiene	8.97	237	119425	65.94	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.09	196	138190	55.85	nq	96
43) 2,4,5-Trichlorophenol	9.13	196	145799	61.48	nq #	85
45) 1,1'-Biphenyl	9.28	154	521308	56.42	nq	97
46) 2-Chloronaphthalene	9.30	162	398827	55.36	nq	95
47) 2-Nitroaniline	9.39	65	156913	54.32	nq	93
48) Acenaphthylene	9.71	152	653460	52.18	nq	99
49) Dimethylphthalate	9.58	163	498466	55.74	nq	100
50) 2,6-Dinitrotoluene	9.64	165	110222	59.66	nq #	84
51) Acenaphthene	9.88	154	406286	52.97	nq	99
52) 3-Nitroaniline	9.81	138	147984	64.97	nq #	77
53) 2,4-Dinitrophenol	9.90	184	61949	63.51	nq #	68
54) Dibenzofuran	10.05	168	539826	52.05	nq	96
55) 4-Nitrophenol	9.96	139	92450	54.22	nq #	56
56) 2,4-Dinitrotoluene	10.04	165	140405	56.79	nq	91
57) Fluorene	10.40	166	411983	51.35	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.18	232	111237	57.11	nq	93
59) Diethylphthalate	10.28	149	573997	61.66	nq	97
60) 4-Chlorophenyl-phenylether	10.40	204	219278	56.09	nq #	78
61) 4-Nitroaniline	10.42	138	142248	60.69	nq	99
62) Azobenzene	10.56	77	613242	58.91	nq	86
64) 4,6-Dinitro-2-methylphenol	10.45	198	93849	66.59	nq #	74
65) n-Nitrosodiphenylamine	10.51	169	393886	49.85	nq	99
66) 4-Bromophenyl-phenylether	10.88	248	125163	51.97	nq #	89
67) Hexachlorobenzene	10.94	284	153030	57.15	nq #	82
68) Atrazine	11.04	200	130477	56.33	nq	97
69) Pentachlorophenol	11.14	266	92331	59.30	nq	98
70) Phenanthrene	11.36	178	700851	54.43	nq	98
71) Anthracene	11.40	178	650289	51.01	nq	100
72) Carbazole	11.56	167	656949	52.93	nq #	97
73) Di-n-butylphthalate	11.89	149	809320	52.15	nq	100
74) Fluoranthene	12.53	202	696054	50.00	nq	96
76) Benzidine	12.66	184	372470	49.48	nq	98
77) Pyrene	12.76	202	714900	50.67	nq	99
79) Butylbenzylphthalate	13.39	149	418349	61.50	nq	87
80) Benzo(a)anthracene	13.95	228	615965	51.05	nq	99
81) 3,3'-Dichlorobenzidine	13.92	252	222969	50.81	nq #	98
82) Chrysene	13.98	228	581413	50.33	nq	99
83) Bis(2-ethylhexyl)phthalate	13.95	149	526126	55.59	nq	99
84) Di-n-octyl phthalate	14.56	149	860539	53.63	nq	100
85) Indeno(1,2,3-cd)pyrene	16.73	276	607605	55.24	nq	100
87) Benzo(b)fluoranthene	14.97	252	643005	56.39	nq #	98
88) Benzo(k)fluoranthene	15.00	252	590163m	59.37	nq	
89) Benzo(a)pyrene	15.31	252	569711	58.13	nq #	97
90) Dibenzo(a,h)anthracene	16.75	278	509041	58.71	nq	98
91) Benzo(g,h,i)perylene	17.14	276	492355	58.38	nq	98

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

