

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF073015\
 Data File : BF080721.D
 Acq On : 31 Jul 2015 6:04
 Operator : TP/UM
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

MMDadoda
 7/31/2015 11:23:12 AM

Quant Time: Jul 31 06:59:52 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF073015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 31 01:45:22 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	198401	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	747286	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	364344	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	632424	20.00	ng	0.00
75) Chrysene-d12	14.00	240	379122	20.00	ng	0.00
86) Perylene-d12	15.41	264	339466	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.45	112	889235	76.89	ng	0.00
7) Phenol-d6	6.49	99	1230420	78.15	ng	0.00
23) Nitrobenzene-d5	7.42	82	1228684	79.62	ng	0.00
41) 2,4,6-Tribromophenol	10.68	330	320666	84.82	ng	0.00
44) 2-Fluorobiphenyl	9.22	172	1792347	73.39	ng	0.00
78) Terphenyl-d14	12.96	244	1673854	83.50	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.50	88	213454	37.43	ng	99
3) Pyridine	3.22	79	576531	36.43	ng	98
4) n-Nitrosodimethylamine	3.17	42	357378	39.51	ng	99
6) Aniline	6.52	93	858715	37.89	ng	99
8) 2-Chlorophenol	6.64	128	479726	37.01	ng	100
9) Benzaldehyde	6.40	77	346348	34.62	ng	87
10) Phenol	6.50	94	751971	41.27	ng	96
11) bis(2-Chloroethyl)ether	6.59	93	475319	36.52	ng	98
12) 1,3-Dichlorobenzene	6.80	146	558759	38.72	ng	99
13) 1,4-Dichlorobenzene	6.88	146	590930	40.14	ng	99
14) 1,2-Dichlorobenzene	7.02	146	480704	35.72	ng	98
15) Benzyl Alcohol	7.00	79	446483	34.14	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.14	45	1050661	37.78	ng	98
17) 2-Methylphenol	7.10	107	419173	39.39	ng	99
18) Hexachloroethane	7.37	117	209971	38.21	ng	99
19) n-Nitroso-di-n-propylamine	7.28	70	395443	37.30	ng	99
20) 3+4-Methylphenols	7.26	107	501774	38.22	ng	97
22) Acetophenone	7.26	105	679527	37.67	ng	97
24) Nitrobenzene	7.44	77	541233	35.01	ng	100
25) Isophorone	7.69	82	1069601	38.98	ng	# 92
26) 2-Nitrophenol	7.76	139	273873m	43.64	ng	
27) 2,4-Dimethylphenol	7.79	122	450659m	38.23	ng	
28) bis(2-Chloroethoxy)methane	7.89	93	648568	39.89	ng	100
29) 2,4-Dichlorophenol	8.00	162	423754	40.44	ng	97
30) 1,2,4-Trichlorobenzene	8.08	180	434999	38.07	ng	98
31) Naphthalene	8.17	128	1400820	37.55	ng	98
32) Benzoic acid	7.92	122	348776m	42.23	ng	
33) 4-Chloroaniline	8.21	127	600199	38.13	ng	99
34) Hexachlorobutadiene	8.28	225	287836	39.01	ng	98
35) Caprolactam	8.59	113	127407	40.60	ng	# 89
36) 4-Chloro-3-methylphenol	8.69	107	481172	42.59	ng	88
37) 2-Methylnaphthalene	8.85	142	946528	40.42	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.01	216	395105	34.53	ng	99
40) Hexachlorocyclopentadiene	9.01	237	268416	38.13	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.13	196	288819m	36.29	ng	
43) 2,4,5-Trichlorophenol	9.16	196	312238m	39.55	ng	
45) 1,1'-Biphenyl	9.32	154	1043643	37.36	ng	98
46) 2-Chloronaphthalene	9.34	162	850489	37.27	ng	99
47) 2-Nitroaniline	9.44	65	377903	42.57	ng	91
48) Acenaphthylene	9.76	152	1438254	39.64	ng	100
49) Dimethylphthalate	9.62	163	923582	35.97	ng	99
50) 2,6-Dinitrotoluene	9.68	165	198518	36.43	ng	91
51) Acenaphthene	9.93	154	880802	39.79	ng	100
52) 3-Nitroaniline	9.85	138	251707	40.03	ng	# 81
53) 2,4-Dinitrophenol	9.94	184	112013	38.60	ng	88
54) Dibenzofuran	10.10	168	1158139	39.12	ng	100
55) 4-Nitrophenol	10.00	139	182739	39.45	ng	83
56) 2,4-Dinitrotoluene	10.08	165	278158	38.93	ng	98
57) Fluorene	10.44	166	901482	39.23	ng	98
58) 2,3,4,6-Tetrachlorophenol	10.21	232	256375	43.06	ng	98
59) Diethylphthalate	10.32	149	935060	37.78	ng	99
60) 4-Chlorophenyl-phenylether	10.43	204	381087	38.10	ng	98
61) 4-Nitroaniline	10.45	138	198624	34.99	ng	96
62) Azobenzene	10.59	77	1078019	40.21	ng	99
64) 4,6-Dinitro-2-methylphenol	10.49	198	161713	42.46	ng	# 35
65) n-Nitrosodiphenylamine	10.56	169	811396	39.14	ng	97
66) 4-Bromophenyl-phenylether	10.92	248	281408	43.11	ng	98
67) Hexachlorobenzene	10.98	284	274090	36.31	ng	98
68) Atrazine	11.08	200	221628	44.84	ng	90
69) Pentachlorophenol	11.17	266	187116	40.97	ng	97
70) Phenanthrene	11.39	178	1179823	37.03	ng	99
71) Anthracene	11.45	178	1277382	40.80	ng	100
72) Carbazole	11.60	167	971867	35.77	ng	99
73) Di-n-butylphthalate	11.94	149	1402656	41.63	ng	100
74) Fluoranthene	12.58	202	1267299	41.14	ng	98
76) Benzidine	12.70	184	571174	41.38	ng	99
77) Pyrene	12.81	202	1231877	41.90	ng	99
79) Butylbenzylphthalate	13.43	149	470662	38.82	ng	98
80) Benzo(a)anthracene	13.99	228	882808	38.72	ng	99
81) 3,3'-Dichlorobenzidine	13.95	252	307638	39.53	ng	97
82) Chrysene	14.03	228	883447	39.92	ng	99
83) Bis(2-ethylhexyl)phthalate	14.00	149	610702	41.72	ng	100
84) Di-n-octyl phthalate	14.60	149	1017630	41.46	ng	100
85) Indeno(1,2,3-cd)pyrene	16.80	276	806586	40.68	ng	# 87
87) Benzo(b)fluoranthene	15.01	252	808521m	39.87	ng	
88) Benzo(k)fluoranthene	15.04	252	633076m	37.94	ng	
89) Benzo(a)pyrene	15.36	252	660621	39.68	ng	98
90) Dibenzo(a,h)anthracene	16.81	278	674809	45.28	ng	# 97
91) Benzo(g,h,i)perylene	17.22	276	678409	46.26	ng	# 93

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

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