

Data Path : Z:\HPCHEM1\BNA F\DATA\BF020216\
 Data File : BF084762.D
 Acq On : 3 Feb 2016 00:13
 Operator : SJ/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

MMDadoda
 2/3/2016 6:25:42 PM

Quant Time: Feb 03 03:02:44 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 01 14:55:11 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	172976	20.00	ng	-0.03
21) Naphthalene-d8	7.98	136	714556	20.00	ng	-0.03
38) Acenaphthene-d10	9.74	164	337044	20.00	ng	-0.02
63) Phenanthrene-d10	11.22	188	632100	20.00	ng	-0.02
75) Chrysene-d12	13.85	240	427906	20.00	ng	-0.02
86) Perylene-d12	15.22	264	351099	20.00	ng	-0.03

System Monitoring Compounds

5) 2-Fluorophenol	5.29	112	931205	83.85	ng	-0.02
7) Phenol-d6	6.35	99	1066237	77.45	ng	-0.02
23) Nitrobenzene-d5	7.28	82	1068517	84.24	ng	-0.02
41) 2,4,6-Tribromophenol	10.53	330	293714	83.54	ng	-0.02
44) 2-Fluorobiphenyl	9.07	172	1790100	91.97	ng	-0.02
78) Terphenyl-d14	12.81	244	1705764	90.33	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.19	88	186858	38.41	ng	# 77
3) Pyridine	2.85	79	564994	44.58	ng	# 77
4) n-Nitrosodimethylamine	2.82	42	262790	40.09	ng	84
6) Aniline	6.36	93	682039	40.49	ng	# 63
8) 2-Chlorophenol	6.49	128	527921	42.00	ng	95
9) Benzaldehyde	6.24	77	323113	39.74	ng	96
10) Phenol	6.36	94	596587	38.68	ng	# 71
11) bis(2-Chloroethyl)ether	6.44	93	485690	40.38	ng	86
12) 1,3-Dichlorobenzene	6.64	146	528998	39.83	ng	# 94
13) 1,4-Dichlorobenzene	6.72	146	545598	41.09	ng	97
14) 1,2-Dichlorobenzene	6.88	146	472155m	38.18	ng	
15) Benzyl Alcohol	6.85	79	383386	40.33	ng	91
16) 2,2'-oxybis(1-Chloropropan	6.99	45	816357	40.35	ng	86
17) 2-Methylphenol	6.98	107	406487	40.89	ng	# 90
18) Hexachloroethane	7.21	117	201464	39.40	ng	92
19) n-Nitroso-di-n-propylamine	7.13	70	332147	38.49	ng	# 76
20) 3+4-Methylphenols	7.13	107	471262	38.19	ng	# 74
22) Acetophenone	7.12	105	713123	37.87	ng	# 98
24) Nitrobenzene	7.29	77	486027	35.78	ng	93
25) Isophorone	7.54	82	983660	39.88	ng	95
26) 2-Nitrophenol	7.61	139	280992	41.94	ng	# 87
27) 2,4-Dimethylphenol	7.65	122	423766	36.37	ng	95
28) bis(2-Chloroethoxy)methane	7.74	93	547724	37.43	ng	99
29) 2,4-Dichlorophenol	7.86	162	405644	40.68	ng	93
30) 1,2,4-Trichlorobenzene	7.93	180	426596	37.88	ng	96
31) Naphthalene	8.01	128	1341270	37.42	ng	99
32) Benzoic acid	7.80	122	313983	42.13	ng	98
33) 4-Chloroaniline	8.06	127	567915	38.31	ng	99
34) Hexachlorobutadiene	8.13	225	275416	42.41	ng	98
35) Caprolactam	8.45	113	134484	43.76	ng	# 70
36) 4-Chloro-3-methylphenol	8.56	107	425079	37.37	ng	93
37) 2-Methylnaphthalene	8.70	142	939435	41.88	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.86	216	396417	37.03	ng	99
40) Hexachlorocyclopentadiene	8.85	237	181624	46.59	ng	95

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42) 2,4,6-Trichlorophenol	8.99	196	275199	39.91	nq	95
43) 2,4,5-Trichlorophenol	9.02	196	281680	40.20	nq	# 88
45) 1,1'-Biphenyl	9.17	154	1159127	38.50	nq	94
46) 2-Chloronaphthalene	9.18	162	836281	37.72	nq	96
47) 2-Nitroaniline	9.29	65	249095	35.48	nq	97
48) Acenaphthylene	9.60	152	1291069	38.38	nq	98
49) Dimethylphthalate	9.48	163	1053314	41.03	nq	# 91
50) 2,6-Dinitrotoluene	9.54	165	249575	44.50	nq	94
51) Acenaphthene	9.78	154	859959	39.29	nq	97
52) 3-Nitroaniline	9.70	138	223776	35.51	nq	99
53) 2,4-Dinitrophenol	9.81	184	106178	43.85	nq	# 82
54) Dibenzofuran	9.95	168	1070269	40.01	nq	95
55) 4-Nitrophenol	9.88	139	176896	38.20	nq	85
56) 2,4-Dinitrotoluene	9.94	165	312756	43.73	nq	95
57) Fluorene	10.28	166	863762	38.67	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.06	232	212397	39.82	nq	92
59) Diethylphthalate	10.17	149	989074	39.46	nq	98
60) 4-Chlorophenyl-phenylether	10.28	204	420924	44.82	nq	97
61) 4-Nitroaniline	10.32	138	242672	39.52	nq	93
62) Azobenzene	10.44	77	1022831	42.44	nq	92
64) 4,6-Dinitro-2-methylphenol	10.35	198	169153	43.36	nq	74
65) n-Nitrosodiphenylamine	10.41	169	842307	41.97	nq	98
66) 4-Bromophenyl-phenylether	10.77	248	297899	42.67	nq	99
67) Hexachlorobenzene	10.83	284	289076	38.00	nq	# 89
68) Atrazine	10.93	200	224798	35.47	nq	98
69) Pentachlorophenol	11.02	266	137636	38.50	nq	97
70) Phenanthrene	11.24	178	1292297	36.86	nq	98
71) Anthracene	11.30	178	1437328	40.69	nq	100
72) Carbazole	11.46	167	1195058	38.80	nq	97
73) Di-n-butylphthalate	11.79	149	1446483	36.88	nq	# 97
74) Fluoranthene	12.43	202	1398689	39.42	nq	93
76) Benzidine	12.56	184	728822	50.62	nq	98
77) Pyrene	12.65	202	1320781	40.31	nq	99
79) Butylbenzylphthalate	13.29	149	635680	42.77	nq	# 80
80) Benzo(a)anthracene	13.84	228	1059804	39.90	nq	99
81) 3,3'-Dichlorobenzidine	13.81	252	411150	43.72	nq	99
82) Chrysene	13.87	228	1011503	39.28	nq	99
83) Bis(2-ethylhexyl)phthalate	13.85	149	793897	42.92	nq	# 96
84) Di-n-octyl phthalate	14.45	149	1199927	39.32	nq	# 89
85) Indeno(1,2,3-cd)pyrene	16.52	276	847486	41.81	nq	99
87) Benzo(b)fluoranthene	14.84	252	833617m	35.34	nq	
88) Benzo(k)fluoranthene	14.88	252	907007m	46.50	nq	
89) Benzo(a)pyrene	15.17	252	795774	39.59	nq	98
90) Dibenzo(a,h)anthracene	16.53	278	706427	41.75	nq	98
91) Benzo(g,h,i)perylene	16.91	276	718786	42.17	nq	99

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

