

Data Path : Z:\HPCHEM1\BNA F\DATA\BF110116\
 Data File : BF090973.D
 Acq On : 2 Nov 2016 1:27
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

Umangi

Quant Time: Nov 02 02:33:49 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102616.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 01 11:53:50 2016
 Response via : Initial Calibration

Internal Standards R.T. QIon Response Conc Units Dev(Min) 11/2/2016 5:42:40 PM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	254122	20.00	ng	-0.01
21) Naphthalene-d8	8.03	136	1036413	20.00	ng	-0.01
38) Acenaphthene-d10	9.78	164	449044	20.00	ng	-0.01
63) Phenanthrene-d10	11.27	188	615331	20.00	ng	0.00
75) Chrysene-d12	13.91	240	582974	20.00	ng	0.00
86) Perylene-d12	15.30	264	404787	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.33	112	1183290	81.41	ng	0.00
7) Phenol-d6	6.40	99	1540627	78.57	ng	0.00
23) Nitrobenzene-d5	7.32	82	1507850	95.24	ng	0.00
41) 2,4,6-Tribromophenol	10.58	330	274774	59.54	ng	0.00
44) 2-Fluorobiphenyl	9.12	172	2327482	91.02	ng	0.00
78) Terphenyl-d14	12.85	244	1713540	74.05	ng	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.26	88	367330	49.43	ng	# 83
3) Pyridine	2.93	79	908688	45.08	ng	# 79
4) n-Nitrosodimethylamine	2.90	42	342588	60.37	ng	# 79
6) Aniline	6.41	93	946957	42.28	ng	# 33
8) 2-Chlorophenol	6.53	128	753072	42.00	ng	# 94
9) Benzaldehyde	6.28	77	490040	48.39	ng	# 67
10) Phenol	6.41	94	758841	35.12	ng	# 1
11) bis(2-Chloroethyl)ether	6.49	93	763500	42.70	ng	# 97
12) 1,3-Dichlorobenzene	6.68	146	727729m	39.68	ng	# 92
13) 1,4-Dichlorobenzene	6.76	146	792316	41.08	ng	# 99
14) 1,2-Dichlorobenzene	6.92	146	670900	36.29	ng	# 78
15) Benzyl Alcohol	6.90	79	574761	44.63	ng	# 45
16) 2,2'-oxybis(1-Chloropropan	7.02	45	1054345	59.26	ng	# 80
17) 2-Methylphenol	7.01	107	548869	40.24	ng	# 90
18) Hexachloroethane	7.25	117	233529	33.27	ng	# 71
19) n-Nitroso-di-n-propylamine	7.17	70	474932	41.89	ng	# 80
20) 3+4-Methylphenols	7.17	107	614235	38.90	ng	# 84
22) Acetophenone	7.16	105	897820	42.11	ng	# 64
24) Nitrobenzene	7.33	77	737466	44.04	ng	# 85
25) Isophorone	7.57	82	1381324	42.99	ng	# 78
26) 2-Nitrophenol	7.65	139	216765	24.16	ng	# 79
27) 2,4-Dimethylphenol	7.70	122	551138	37.95	ng	# 94
28) bis(2-Chloroethoxy)methane	7.79	93	811393	44.41	ng	# 99
29) 2,4-Dichlorophenol	7.90	162	480498	35.73	ng	# 95
30) 1,2,4-Trichlorobenzene	7.97	180	552997	35.58	ng	# 96
31) Naphthalene	8.05	128	1738653	35.89	ng	# 62
32) Benzoic acid	7.84	122	213588	20.25	ng	# 88
33) 4-Chloroaniline	8.11	127	822776	42.03	ng	# 99
34) Hexachlorobutadiene	8.18	225	321322	35.16	ng	# 80
35) Caprolactam	8.50	113	149437	35.80	ng	# 82
36) 4-Chloro-3-methylphenol	8.60	107	580195	39.67	ng	# 90
37) 2-Methylnaphthalene	8.75	142	1184765	37.86	ng	# 97
39) 1,2,4,5-Tetrachlorobenzene	8.91	216	469428	38.11	ng	# 96
40) Hexachlorocyclopentadiene	8.90	237	23200	4.09	ng	

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)	11/2/2016 5:42:40 PM
42) 2,4,6-Trichlorophenol	9.02	196	288690	36.30	nq		98
43) 2,4,5-Trichlorophenol	9.07	196	314405	38.38	nq	#	91
45) 1,1'-Biphenyl	9.21	154	1304042	40.32	nq		90
46) 2-Chloronaphthalene	9.23	162	997485	40.54	nq	#	88
47) 2-Nitroaniline	9.33	65	403989	58.19	nq	#	68
48) Acenaphthylene	9.64	152	1466132	39.36	nq		95
49) Dimethylphthalate	9.52	163	1162366	38.74	nq	#	97
50) 2,6-Dinitrotoluene	9.57	165	198580	29.97	nq	#	15
51) Acenaphthene	9.82	154	972132	37.89	nq		89
52) 3-Nitroaniline	9.74	138	299371	42.44	nq	#	50
54) Dibenzofuran	9.98	168	1226562	37.04	nq	#	81
55) 4-Nitrophenol	9.92	139	127849	29.72	nq	#	25
56) 2,4-Dinitrotoluene	9.98	165	255361	31.02	nq	#	76
57) Fluorene	10.33	166	908587	33.50	nq		92
58) 2,3,4,6-Tetrachlorophenol	10.11	232	187944m	25.69	nq		
59) Diethylphthalate	10.21	149	1223574	41.06	nq		99
60) 4-Chlorophenyl-phenylether	10.33	204	530544	40.07	nq		95
61) 4-Nitroaniline	10.36	138	304832	43.63	nq	#	47
62) Azobenzene	10.49	77	1352267	46.15	nq		88
65) n-Nitrosodiphenylamine	10.44	169	836449	44.38	nq		91
66) 4-Bromophenyl-phenylether	10.82	248	268441	39.84	nq	#	85
67) Hexachlorobenzene	10.88	284	272682	34.19	nq	#	67
68) Atrazine	10.98	200	110564	18.96	nq		82
69) Pentachlorophenol	11.07	266	103699	24.24	nq		98
70) Phenanthrene	11.29	178	1186920	37.88	nq		95
71) Anthracene	11.35	178	1380159	41.83	nq		97
72) Carbazole	11.49	167	1126403	38.66	nq	#	93
73) Di-n-butylphthalate	11.84	149	1681187	44.76	nq	#	98
74) Fluoranthene	12.48	202	1280634	37.32	nq		92
76) Benzidine	12.60	184	655601	51.50	nq	#	97
77) Pyrene	12.71	202	1311132m	35.54	nq		
79) Butylbenzylphthalate	13.32	149	626014	37.40	nq	#	71
80) Benzo(a)anthracene	13.89	228	1178479	38.09	nq		96
81) 3,3'-Dichlorobenzidine	13.86	252	511347	42.57	nq	#	95
82) Chrysene	13.93	228	1132954m	36.20	nq		
83) Bis(2-ethylhexyl)phthalate	13.89	149	919997	42.20	nq		98
84) Di-n-octyl phthalate	14.50	149	1478127	40.38	nq	#	94
85) Indeno(1,2,3-cd)pyrene	16.65	276	889419	32.07	nq	#	95
87) Benzo(b)fluoranthene	14.91	252	1182858m	43.46	nq		
88) Benzo(k)fluoranthene	14.93	252	881328m	40.97	nq		
89) Benzo(a)pyrene	15.24	252	923739	39.50	nq	#	91
90) Dibenzo(a,h)anthracene	16.66	278	784273	39.61	nq	#	81
91) Benzo(g,h,i)perylene	17.05	276	696885	37.37	nq	#	79

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

