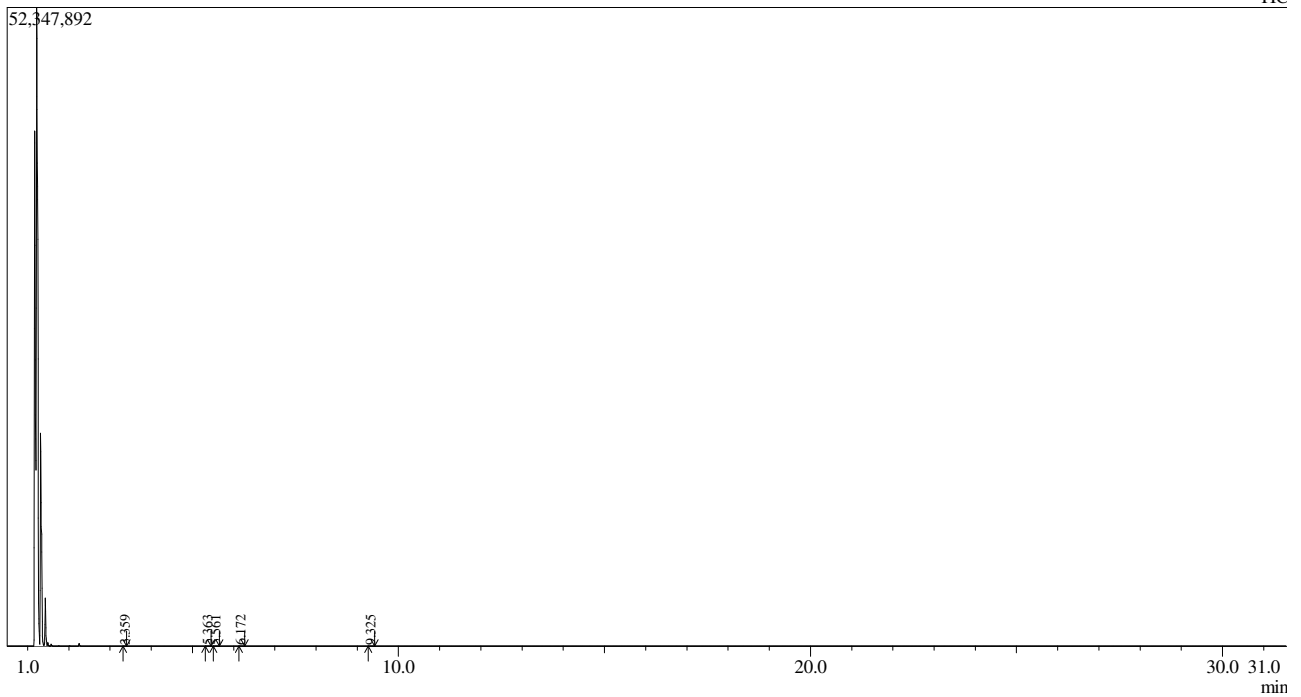




Peak Report TIC									
Peak#	R.Time	I.Time	F.Time	Area	Area%	Height	A/H	Mark	Name
1	3.359	3.315	3.400	16353	17.51	10043	1.63	MI	Toluene
2	5.363	5.310	5.450	17539	18.78	7251	2.42	MI	Ethylbenzene
3	5.561	5.505	5.655	23636	25.31	10403	2.27	MI	o-Xylene
4	6.172	6.125	6.265	21640	23.17	7618	2.84	MI	o-Xylene
5	9.325	9.265	9.425	14230	15.24	4664	3.05	MI	Decane
				93398	100.00	39979			

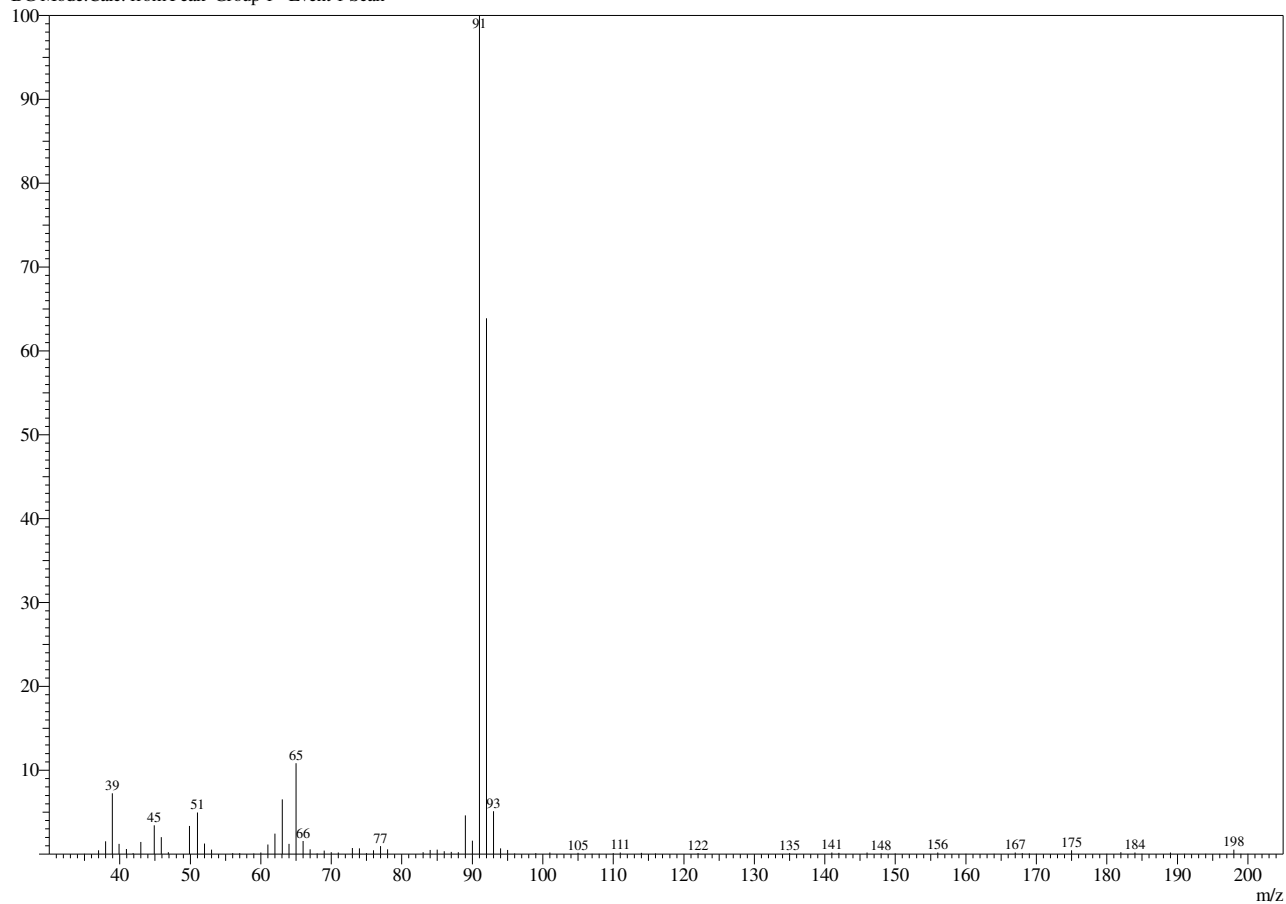
Spectrum

Peak#:1 R.Time:3.359(Scan#:573)

MassPeaks:85

RawMode:Averaged 3.355-3.365(572-574)

BG Mode:Calc. from Peak Group 1 - Event 1 Scan



Mass Table

Peak#:1 R.Time:3.360(Scan#:573)

MassPeaks:85

Group 1 - Event 1 Scan

#	m/z	Rel. Int.	#	m/z	Rel. Int.	#	m/z	Rel. Int.	#	m/z	Rel. Int.
1	35.00	0.02	23	64.00	1.21	45	91.00	100.00	67	156.00	0.20
2	37.00	0.47	24	65.00	10.83	46	92.00	63.90	68	160.00	0.02
3	38.00	1.54	25	66.00	1.56	47	93.00	5.10	69	164.00	0.05
4	38.95	7.23	26	67.00	0.57	48	94.00	0.67	70	167.00	0.17
5	39.90	1.24	27	68.00	0.07	49	95.00	0.50	71	168.00	0.12
6	40.95	0.62	28	69.00	0.42	50	96.00	0.07	72	169.00	0.07
7	41.90	0.10	29	70.00	0.22	51	99.00	0.10	73	170.00	0.02
8	43.00	1.46	30	71.00	0.20	52	101.00	0.17	74	175.00	0.45
9	44.90	3.42	31	73.00	0.72	53	105.00	0.15	75	180.00	0.02
10	45.90	2.03	32	74.00	0.69	54	107.00	0.02	76	182.00	0.22
11	46.90	0.25	33	75.00	0.12	55	108.00	0.05	77	184.00	0.20
12	49.90	3.34	34	76.00	0.47	56	110.00	0.12	78	185.00	0.05
13	51.00	4.96	35	77.00	0.97	57	111.00	0.17	79	189.00	0.17
14	52.00	1.26	36	78.00	0.57	58	114.00	0.15	80	191.00	0.02
15	53.00	0.52	37	83.00	0.22	59	116.00	0.07	81	194.00	0.02
16	56.00	0.10	38	84.00	0.50	60	122.00	0.12	82	196.00	0.05
17	57.00	0.10	39	85.00	0.55	61	135.00	0.10	83	198.00	0.52
18	59.00	0.07	40	86.00	0.35	62	141.00	0.17	84	199.00	0.05
19	60.00	0.17	41	87.00	0.27	63	142.00	0.15	85	200.00	0.02
20	61.00	1.14	42	88.00	0.25	64	146.00	0.17			
21	62.00	2.43	43	89.00	4.61	65	148.00	0.07			
22	63.05	6.54	44	90.00	1.61	66	150.00	0.07			

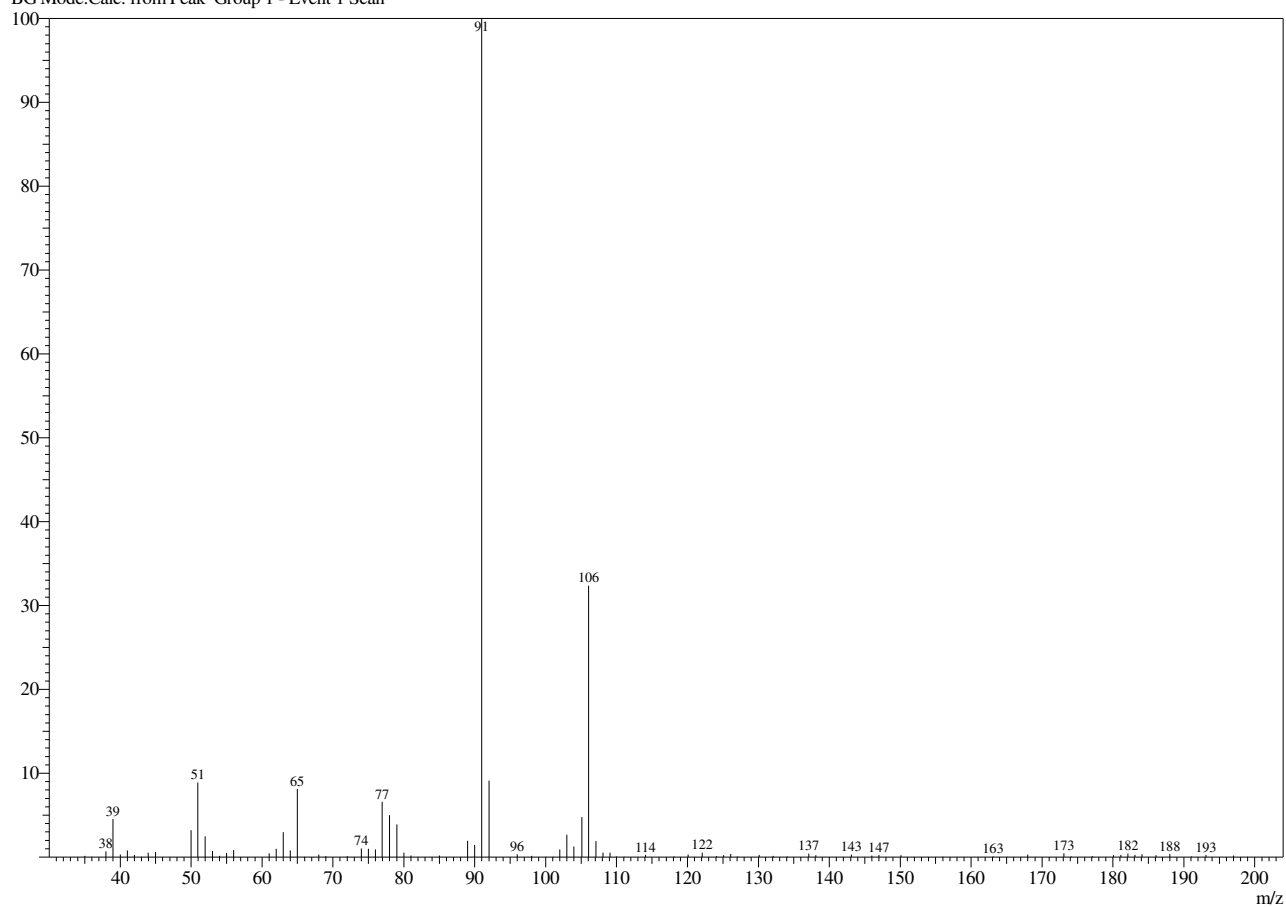
Spectrum

Peak#:2 R.Time:5.363(Scan#:974)

MassPeaks:88

RawMode:Averaged 5.360-5.370(973-975)

BG Mode:Calc. from Peak Group 1 - Event 1 Scan



Mass Table

Peak#:2 R.Time:5.365(Scan#:974)

MassPeaks:88

Group 1 - Event 1 Scan

#	m/z	Rel. Int.	#	m/z	Rel. Int.	#	m/z	Rel. Int.	#	m/z	Rel. Int.
1	35.00	0.15	23	66.00	0.03	45	103.00	2.65	67	143.10	0.31
2	36.00	0.03	24	68.00	0.31	46	104.00	1.25	68	146.00	0.18
3	38.00	0.70	25	69.00	0.12	47	105.10	4.76	69	147.00	0.27
4	39.00	4.55	26	72.00	0.12	48	106.05	32.37	70	150.10	0.24
5	40.00	0.31	27	74.00	1.04	49	107.10	1.89	71	156.00	0.03
6	41.00	0.79	28	75.00	1.01	50	108.10	0.55	72	163.10	0.09
7	42.00	0.27	29	76.00	0.92	51	109.10	0.52	73	165.10	0.03
8	43.00	0.03	30	76.95	6.62	52	110.00	0.06	74	168.00	0.31
9	43.95	0.52	31	78.00	5.00	53	111.10	0.03	75	173.10	0.46
10	45.00	0.61	32	79.00	3.87	54	114.10	0.27	76	174.10	0.06
11	50.00	3.20	33	80.00	0.55	55	118.10	0.06	77	177.10	0.12
12	50.95	8.88	34	81.00	0.18	56	120.10	0.31	78	180.10	0.24
13	52.00	2.47	35	85.00	0.18	57	122.10	0.52	79	181.10	0.27
14	53.00	0.73	36	89.00	1.95	58	124.10	0.12	80	182.10	0.43
15	54.00	0.15	37	90.00	1.43	59	125.10	0.21	81	183.10	0.27
16	55.00	0.49	38	91.00	100.00	60	126.10	0.40	82	184.10	0.34
17	56.00	0.82	39	92.00	9.12	61	127.10	0.09	83	186.10	0.24
18	61.00	0.46	40	96.00	0.34	62	130.10	0.27	84	188.00	0.37
19	62.00	1.01	41	98.00	0.15	63	132.10	0.15	85	191.10	0.09
20	63.00	2.99	42	100.10	0.03	64	134.10	0.06	86	193.10	0.27
21	64.00	0.79	43	101.10	0.03	65	137.10	0.43	87	197.00	0.18
22	65.00	8.11	44	102.00	0.92	66	138.10	0.21	88	199.00	0.03

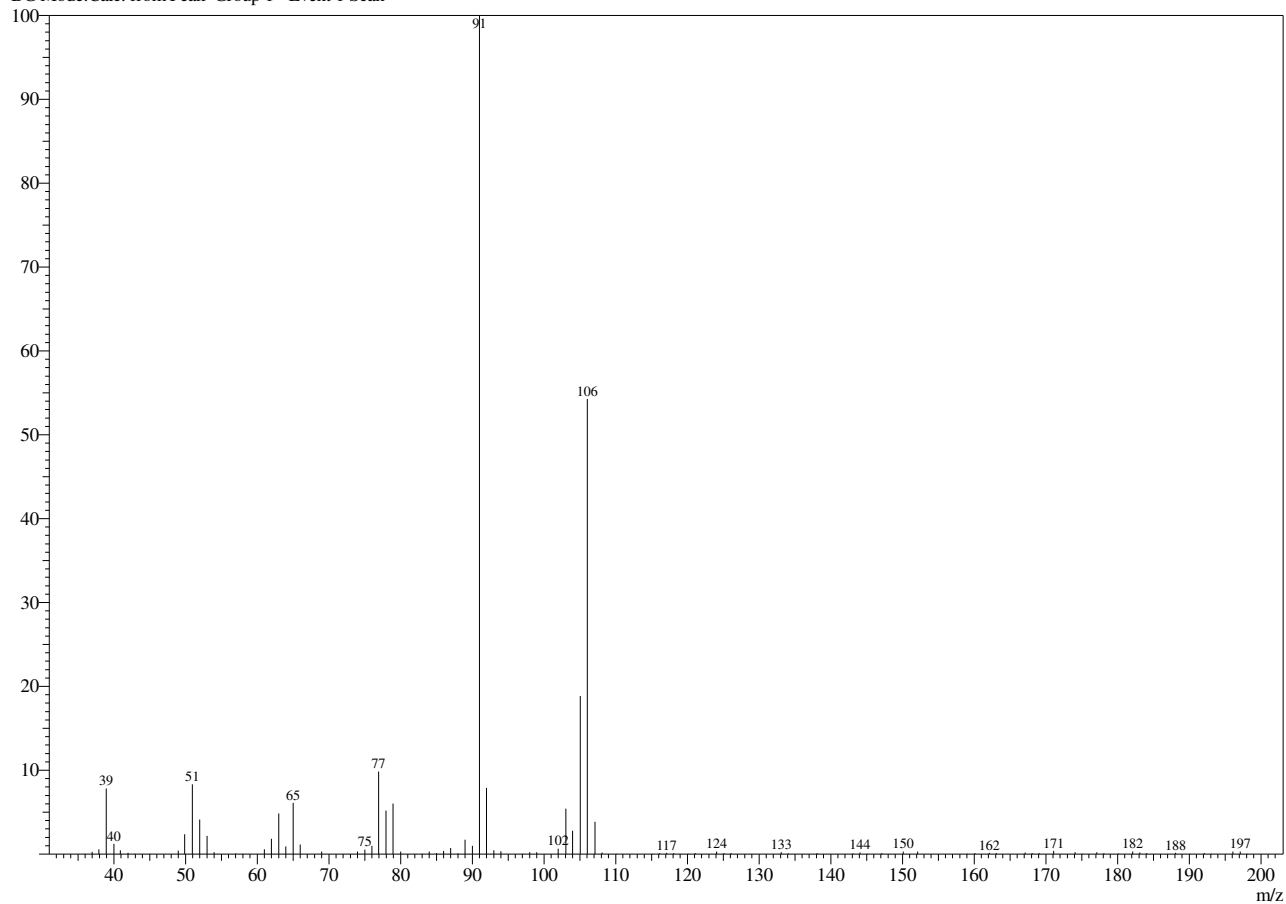
Spectrum

Peak#:3 R.Time:5.561(Scan#:1013)

MassPeaks:95

RawMode:Averaged 5.555-5.565(1012-1014)

BG Mode:Calc. from Peak Group 1 - Event 1 Scan



Mass Table

Peak#:3 R.Time:5.560(Scan#:1013)

MassPeaks:95

Group 1 - Event 1 Scan

#	m/z	Rel. Int.	#	m/z	Rel. Int.	#	m/z	Rel. Int.	#	m/z	Rel. Int.
1	36.00	0.03	25	74.00	0.29	49	104.00	2.80	73	159.10	0.03
2	37.00	0.26	26	75.00	0.53	50	105.05	18.86	74	160.10	0.11
3	37.90	0.55	27	76.00	0.98	51	106.05	54.27	75	162.10	0.16
4	38.95	7.81	28	76.95	9.86	52	107.10	3.85	76	163.10	0.11
5	40.00	1.24	29	77.95	5.20	53	108.10	0.18	77	167.10	0.18
6	40.90	0.45	30	78.95	6.01	54	110.10	0.03	78	168.10	0.11
7	42.00	0.16	31	80.00	0.29	55	115.10	0.05	79	169.10	0.13
8	42.90	0.05	32	83.00	0.05	56	116.10	0.11	80	171.10	0.40
9	45.00	0.03	33	84.00	0.32	57	117.10	0.16	81	174.10	0.21
10	49.00	0.42	34	85.00	0.11	58	118.10	0.11	82	175.10	0.03
11	49.90	2.37	35	86.00	0.37	59	120.10	0.03	83	177.10	0.24
12	50.95	8.31	36	87.00	0.71	60	121.10	0.11	84	178.10	0.11
13	52.00	4.14	37	88.00	0.08	61	124.10	0.32	85	180.10	0.08
14	53.00	2.19	38	89.00	1.71	62	125.10	0.11	86	182.10	0.32
15	54.00	0.21	39	90.00	0.98	63	126.10	0.11	87	183.10	0.18
16	57.00	0.08	40	91.00	100.00	64	133.10	0.21	88	184.10	0.11
17	58.00	0.03	41	92.00	7.91	65	134.10	0.08	89	188.10	0.16
18	61.00	0.55	42	93.00	0.47	66	138.10	0.13	90	190.10	0.03
19	62.00	1.82	43	94.00	0.34	67	143.10	0.08	91	192.10	0.11
20	63.00	4.85	44	97.00	0.05	68	144.10	0.18	92	193.10	0.03
21	64.00	0.92	45	98.00	0.24	69	145.10	0.05	93	196.10	0.29
22	65.00	6.12	46	99.00	0.21	70	147.10	0.11	94	197.10	0.32
23	66.00	1.16	47	102.00	0.66	71	150.10	0.32	95	198.10	0.03
24	69.00	0.32	48	103.05	5.43	72	152.10	0.29			

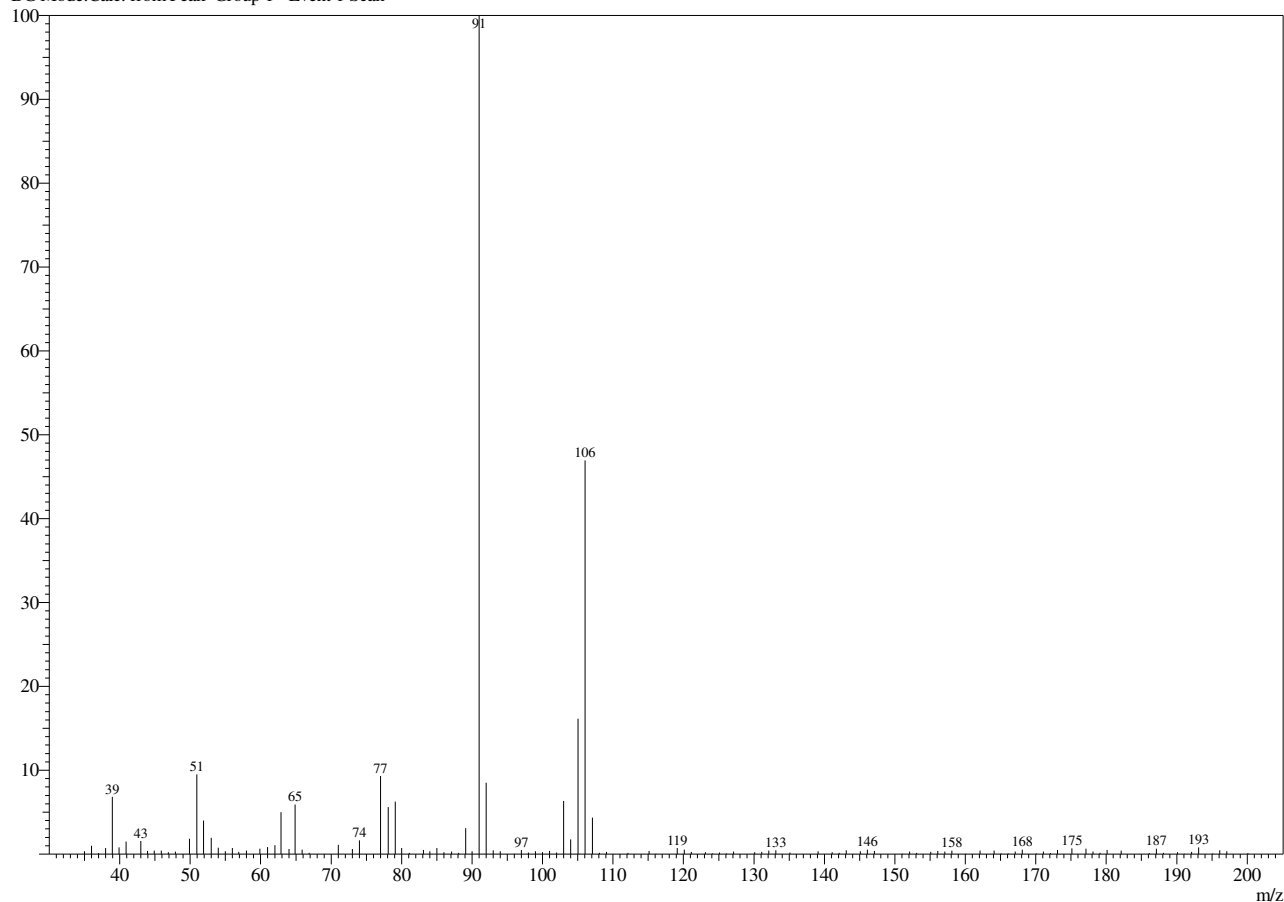
Spectrum

Peak#4 R.Time:6.172(Scan#:1135)

MassPeaks:122

RawMode:Averaged 6.165-6.175(1134-1136)

BG Mode:Calc. from Peak Group 1 - Event 1 Scan



Mass Table

Peak#4 R.Time:6.170(Scan#:1135)

MassPeaks:122

Group 1 - Event 1 Scan

#	m/z	Rel. Int.	#	m/z	Rel. Int.	#	m/z	Rel. Int.	#	m/z	Rel. Int.
1	35.00	0.34	32	73.00	0.61	63	108.10	0.15	94	158.10	0.42
2	36.00	1.00	33	74.00	1.65	64	109.10	0.27	95	160.10	0.04
3	37.00	0.11	34	77.00	9.30	65	112.10	0.15	96	162.10	0.42
4	38.00	0.73	35	78.10	5.63	66	115.10	0.38	97	164.10	0.42
5	38.95	6.85	36	79.10	6.28	67	116.10	0.04	98	165.10	0.08
6	39.90	0.80	37	80.00	0.73	68	119.10	0.73	99	166.10	0.08
7	40.90	1.49	38	81.10	0.15	69	120.10	0.54	100	167.10	0.31
8	43.00	1.57	39	83.10	0.50	70	121.10	0.27	101	168.10	0.54
9	43.95	0.38	40	84.00	0.34	71	123.10	0.23	102	170.10	0.08
10	44.90	0.42	41	85.00	0.73	72	124.10	0.08	103	171.10	0.31
11	45.90	0.42	42	86.00	0.23	73	125.10	0.19	104	172.10	0.11
12	46.90	0.23	43	87.10	0.31	74	126.10	0.08	105	173.10	0.50
13	47.90	0.31	44	88.10	0.15	75	127.10	0.31	106	175.10	0.69
14	49.90	1.84	45	89.10	3.10	76	130.10	0.23	107	177.10	0.65
15	50.95	9.49	46	90.00	0.34	77	131.10	0.27	108	178.10	0.31
16	51.90	4.02	47	91.00	100.00	78	132.10	0.42	109	179.10	0.15
17	53.00	1.95	48	92.00	8.50	79	133.10	0.46	110	180.10	0.50
18	54.00	0.77	49	93.00	0.46	80	135.10	0.19	111	182.10	0.42
19	55.00	0.34	50	94.00	0.34	81	139.10	0.34	112	186.10	0.11
20	56.00	0.73	51	96.00	0.04	82	141.10	0.23	113	187.10	0.65
21	56.90	0.27	52	97.00	0.50	83	142.10	0.15	114	188.10	0.15
22	58.00	0.42	53	98.00	0.19	84	143.10	0.46	115	190.10	0.23
23	59.90	0.65	54	99.00	0.34	85	145.10	0.34	116	191.10	0.04
24	61.00	0.84	55	100.00	0.23	86	146.10	0.54	117	192.10	0.27
25	62.00	1.07	56	101.00	0.38	87	147.10	0.38	118	193.10	0.80
26	62.90	5.01	57	102.00	0.19	88	151.10	0.04	119	195.10	0.08
27	64.00	0.61	58	103.00	6.35	89	152.10	0.31	120	196.10	0.46
28	64.90	5.93	59	104.00	1.76	90	153.10	0.15	121	197.10	0.34
29	65.90	0.54	60	105.05	16.15	91	155.10	0.27	122	200.10	0.11
30	66.90	0.15	61	106.05	46.96	92	156.10	0.38			
31	71.00	1.11	62	107.10	4.36	93	157.10	0.31			

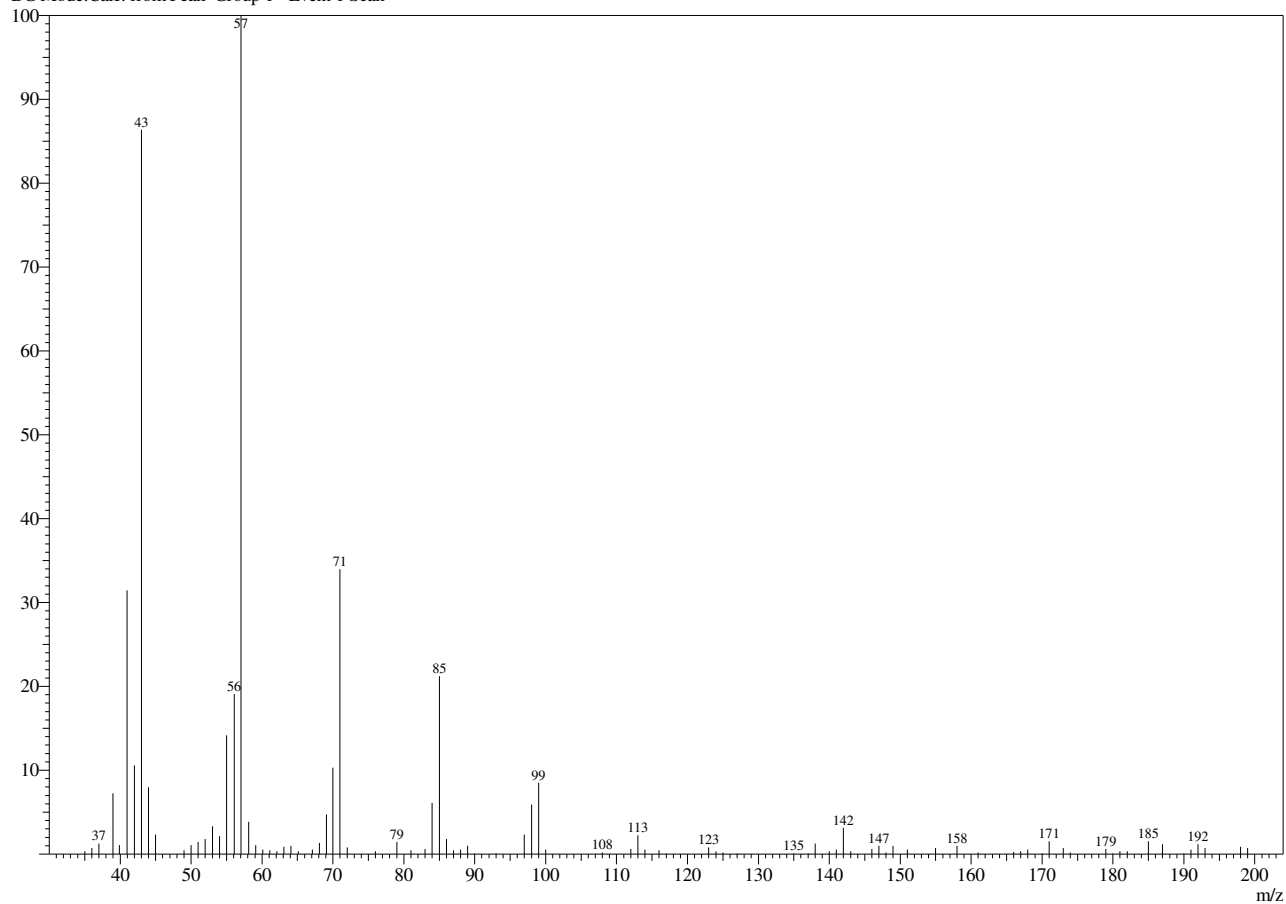
Spectrum

Peak#:5 R.Time:9.325(Scan#:1766)

MassPeaks:93

RawMode:Averaged 9.320-9.330(1765-1767)

BG Mode:Calc. from Peak Group 1 - Event 1 Scan



Mass Table

Peak#:5 R.Time:9.325(Scan#:1766)

MassPeaks:93

Group 1 - Event 1 Scan

#	m/z	Rel. Int.	#	m/z	Rel. Int.	#	m/z	Rel. Int.	#	m/z	Rel. Int.
1	35.00	0.36	25	63.10	0.90	49	100.00	0.54	73	158.00	0.99
2	36.00	0.72	26	64.10	0.99	50	108.00	0.18	74	161.00	0.18
3	37.00	1.25	27	65.10	0.36	51	112.00	0.63	75	166.00	0.27
4	39.00	7.26	28	67.10	0.54	52	113.00	2.24	76	167.00	0.36
5	39.90	1.08	29	68.10	1.34	53	114.00	0.54	77	168.00	0.54
6	40.95	31.45	30	69.10	4.75	54	116.00	0.45	78	170.00	0.09
7	42.00	10.57	31	70.00	10.30	55	117.00	0.09	79	171.00	1.52
8	43.00	86.38	32	71.00	33.96	56	123.00	0.81	80	173.00	0.72
9	44.00	7.97	33	72.00	0.81	57	124.00	0.36	81	174.00	0.18
10	45.00	2.33	34	76.00	0.36	58	125.00	0.18	82	179.00	0.63
11	49.00	0.45	35	79.00	1.43	59	127.00	0.09	83	180.00	0.09
12	50.00	1.08	36	81.00	0.45	60	135.00	0.09	84	181.00	0.36
13	51.00	1.43	37	83.00	0.63	61	137.00	0.09	85	182.00	0.36
14	52.00	1.79	38	84.00	6.09	62	138.00	1.25	86	185.00	1.52
15	53.00	3.32	39	85.00	21.24	63	140.00	0.36	87	187.00	1.16
16	54.00	2.15	40	86.00	1.79	64	141.00	0.54	88	190.00	0.09
17	55.00	14.16	41	87.00	0.45	65	142.00	3.14	89	191.00	0.54
18	56.10	19.09	42	88.00	0.54	66	143.00	0.36	90	192.00	1.16
19	57.05	100.00	43	89.00	0.99	67	146.00	0.63	91	193.00	0.72
20	58.10	3.85	44	90.00	0.09	68	147.00	0.99	92	198.00	0.90
21	59.10	1.08	45	93.00	0.09	69	148.00	0.18	93	199.00	0.72
22	60.10	0.54	46	97.00	2.33	70	149.00	0.99			
23	61.10	0.45	47	98.00	5.91	71	151.00	0.54			
24	62.10	0.36	48	99.00	8.51	72	155.00	0.72			

Library

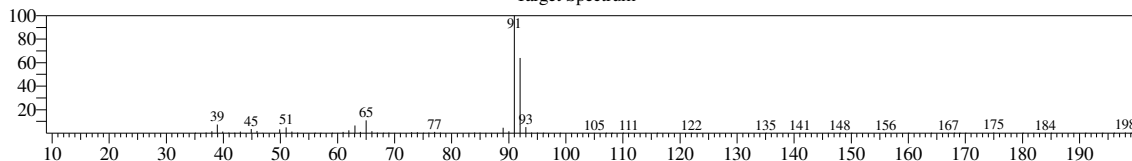
<< Target >>

Line#:1 R.Time:3.360(Scan#:573) MassPeaks:85

RawMode:Averaged 3.355-3.365(572-574) BasePeak:91.00(4036)

BG Mode:Calc. from Peak Group 1 - Event 1 Scan

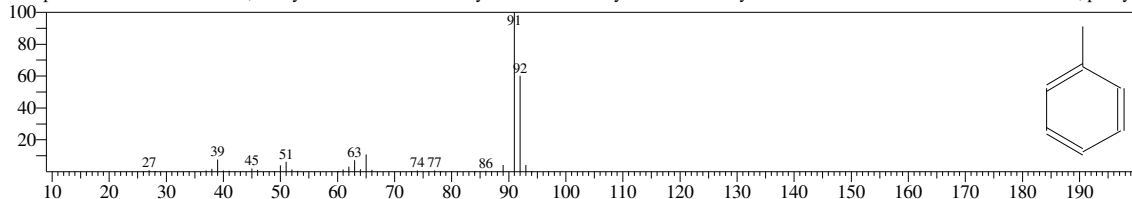
Target Spectrum



Hit#:1 Entry:1275 Library:NIST14s.lib

SI:97 Formula:C7H8 CAS:108-88-3 MolWeight:92 RetIndex:794

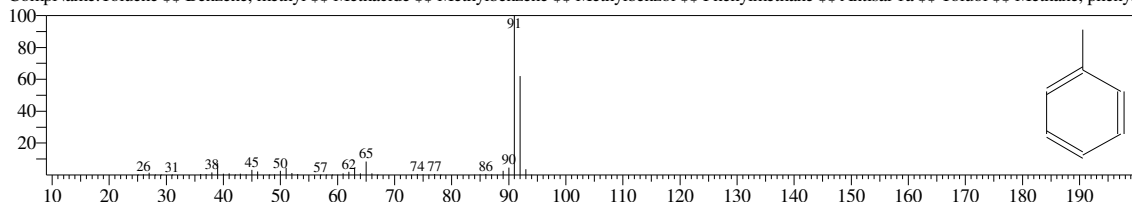
CompName:Toluene \$\$ Benzene, methyl \$\$ Methacide \$\$ Methylbenzene \$\$ Methylbenzol \$\$ Phenylmethane \$\$ Antisal 1a \$\$ Toluol \$\$ Methane, phenyl-



Hit#:2 Entry:1273 Library:NIST14s.lib

SI:97 Formula:C7H8 CAS:108-88-3 MolWeight:92 RetIndex:794

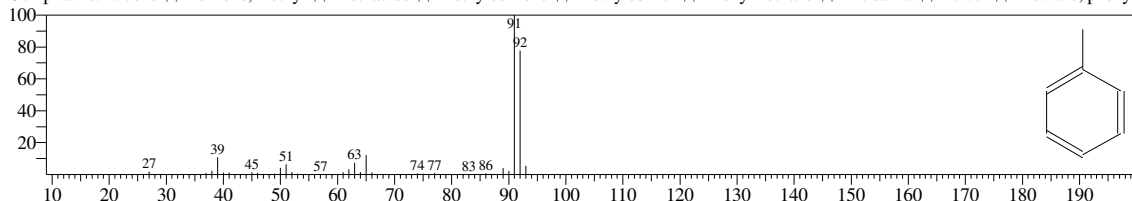
CompName:Toluene \$\$ Benzene, methyl \$\$ Methacide \$\$ Methylbenzene \$\$ Methylbenzol \$\$ Phenylmethane \$\$ Antisal 1a \$\$ Toluol \$\$ Methane, phenyl-



Hit#:3 Entry:1224 Library:NIST14s.lib

SI:95 Formula:C7H8 CAS:108-88-3 MolWeight:92 RetIndex:794

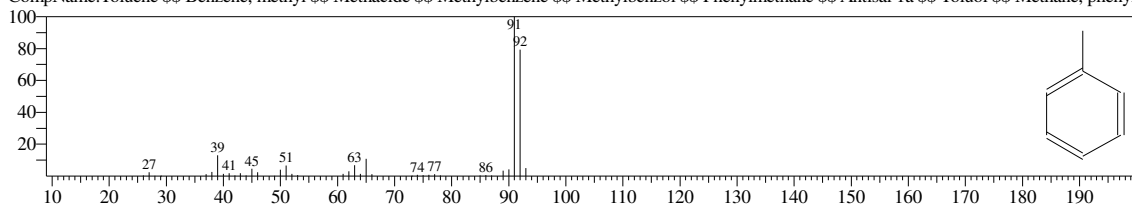
CompName:Toluene \$\$ Benzene, methyl \$\$ Methacide \$\$ Methylbenzene \$\$ Methylbenzol \$\$ Phenylmethane \$\$ Antisal 1a \$\$ Toluol \$\$ Methane, phenyl-



Hit#:4 Entry:1267 Library:NIST14s.lib

SI:95 Formula:C7H8 CAS:108-88-3 MolWeight:92 RetIndex:794

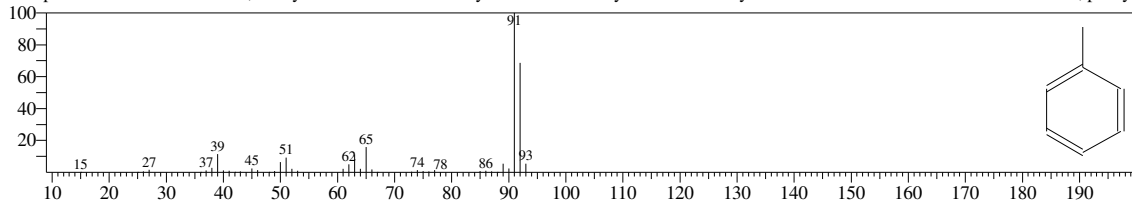
CompName:Toluene \$\$ Benzene, methyl \$\$ Methacide \$\$ Methylbenzene \$\$ Methylbenzol \$\$ Phenylmethane \$\$ Antisal 1a \$\$ Toluol \$\$ Methane, phenyl-



Hit#:5 Entry:1274 Library:NIST14s.lib

SI:95 Formula:C7H8 CAS:108-88-3 MolWeight:92 RetIndex:794

CompName:Toluene \$\$ Benzene, methyl \$\$ Methacide \$\$ Methylbenzene \$\$ Methylbenzol \$\$ Phenylmethane \$\$ Antisal 1a \$\$ Toluol \$\$ Methane, phenyl-



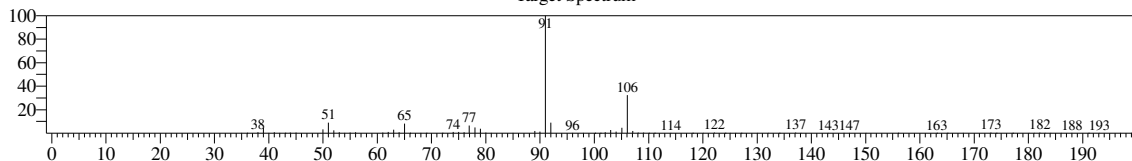
<< Target >>

Line#:2 R.Time:5.365(Scan#:974) MassPeaks:88

RawMode:Averaged 5.360-5.370(973-975) BasePeak:91.00(3278)

BG Mode:Calc. from Peak Group 1 - Event 1 Scan

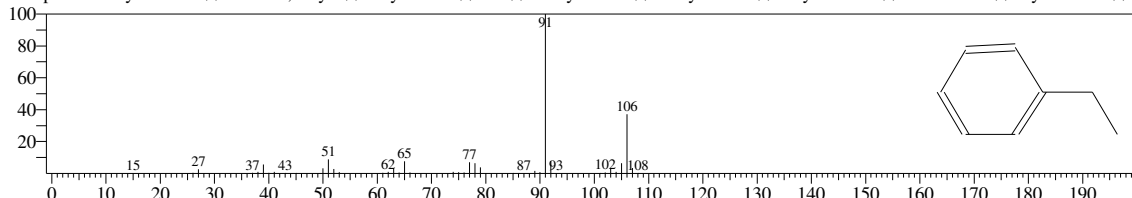
Target Spectrum



Hit#:1 Entry:2463 Library:NIST14s.lib

SI:96 Formula:C8H10 CAS:100-41-4 MolWeight:106 RetIndex:893

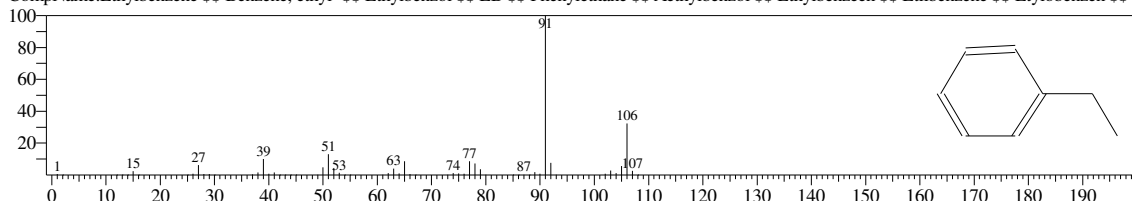
CompName:Ethylbenzene \$\$ Benzene, ethyl- \$\$ Ethylbenzol \$\$ EB \$\$ Phenylethane \$\$ Aethylbenzol \$\$ Ethylbenzeen \$\$ Etilbenzene \$\$ Etylobenzen \$\$ N



Hit#:2 Entry:2461 Library:NIST14s.lib

SI:95 Formula:C8H10 CAS:100-41-4 MolWeight:106 RetIndex:893

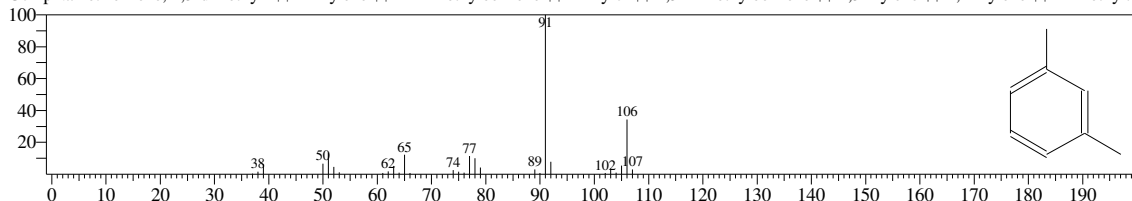
CompName:Ethylbenzene \$\$ Benzene, ethyl- \$\$ Ethylbenzol \$\$ EB \$\$ Phenylethane \$\$ Aethylbenzol \$\$ Ethylbenzeen \$\$ Etilbenzene \$\$ Etylobenzen \$\$ N



Hit#:3 Entry:2464 Library:NIST14s.lib

SI:93 Formula:C8H10 CAS:108-38-3 MolWeight:106 RetIndex:907

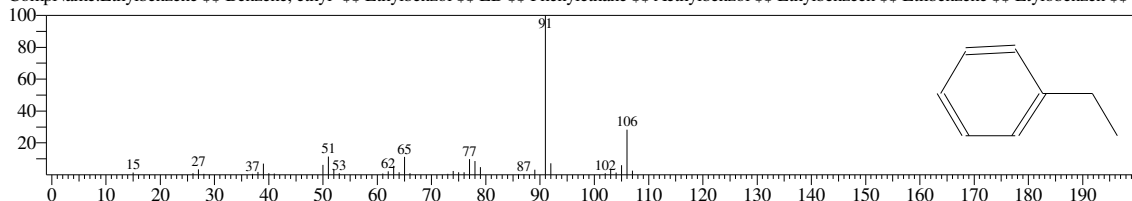
CompName:Benzene, 1,3-dimethyl- \$\$ m-Xylene \$\$ m-Dimethylbenzene \$\$ m-Xylol \$\$ 1,3-Dimethylbenzene \$\$ 1,3-Xylene \$\$ 2,4-Xylene \$\$ m-Methyltol



Hit#:4 Entry:2723 Library:NIST14s.lib

SI:93 Formula:C8H10 CAS:100-41-4 MolWeight:106 RetIndex:893

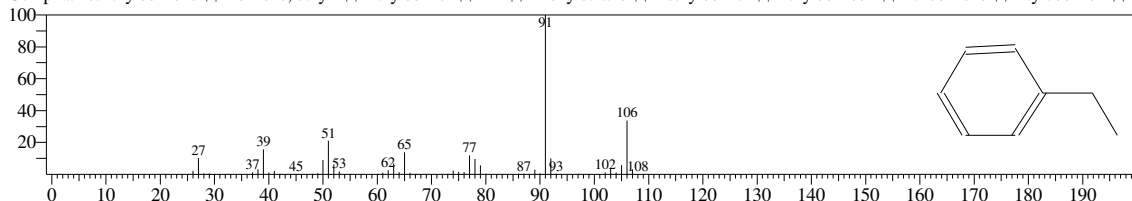
CompName:Ethylbenzene \$\$ Benzene, ethyl- \$\$ Ethylbenzol \$\$ EB \$\$ Phenylethane \$\$ Aethylbenzol \$\$ Ethylbenzeen \$\$ Etilbenzene \$\$ Etylobenzen \$\$ N



Hit#:5 Entry:2462 Library:NIST14s.lib

SI:91 Formula:C8H10 CAS:100-41-4 MolWeight:106 RetIndex:893

CompName:Ethylbenzene \$\$ Benzene, ethyl- \$\$ Ethylbenzol \$\$ EB \$\$ Phenylethane \$\$ Aethylbenzol \$\$ Ethylbenzeen \$\$ Etilbenzene \$\$ Etylobenzen \$\$ N



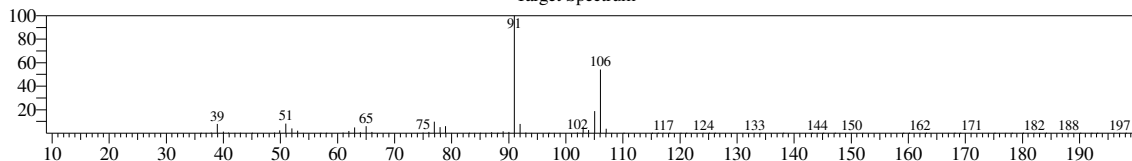
<< Target >>

Line#:3 R.Time:5.560(Scan#:1013) MassPeaks:95

RawMode:Averaged 5.555-5.565(1012-1014) BasePeak:91.00(3792)

BG Mode:Calc. from Peak Group 1 - Event 1 Scan

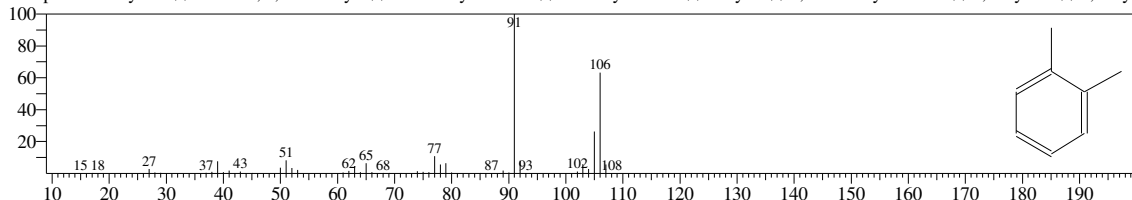
Target Spectrum



Hit#:1 Entry:2478 Library:NIST14s.lib

SI:96 Formula:C₈H₁₀ CAS:95-47-6 MolWeight:106 RetIndex:907

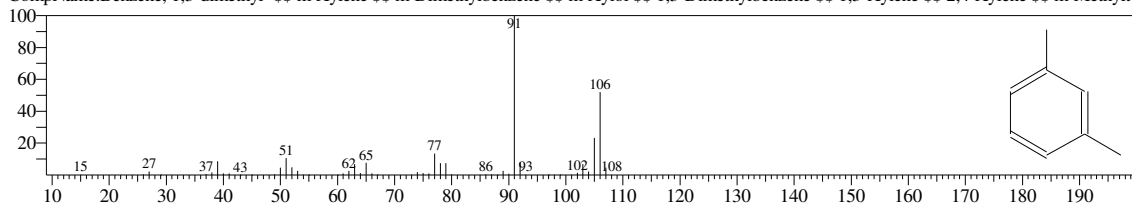
CompName:o-Xylene \$\$ Benzene, 1,2-dimethyl- \$\$ o-Dimethylbenzene \$\$ o-Methyltoluene \$\$ o-Xylol \$\$ 1,2-Dimethylbenzene \$\$ 1,2-Xylene \$\$ 3,4-Xyle



Hit#:2 Entry:2725 Library:NIST14.lib

SI:96 Formula:C₈H₁₀ CAS:108-38-3 MolWeight:106 RetIndex:907

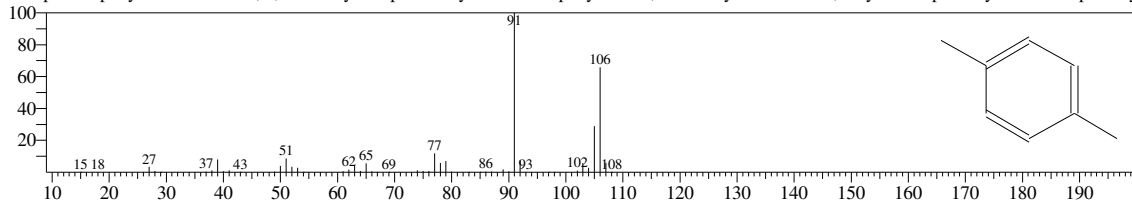
CompName:Benzen, 1,3-dimethyl- \$\$ m-Xylene \$\$ m-Dimethylbenzene \$\$ m-Xylol \$\$ 1,3-Dimethylbenzene \$\$ 1,3-Xylene \$\$ 2,4-Xylene \$\$ m-Methyltol



Hit#:3 Entry:2726 Library:NIST14.lib

SI:96 Formula:C₈H₁₀ CAS:106-42-3 MolWeight:106 RetIndex:907

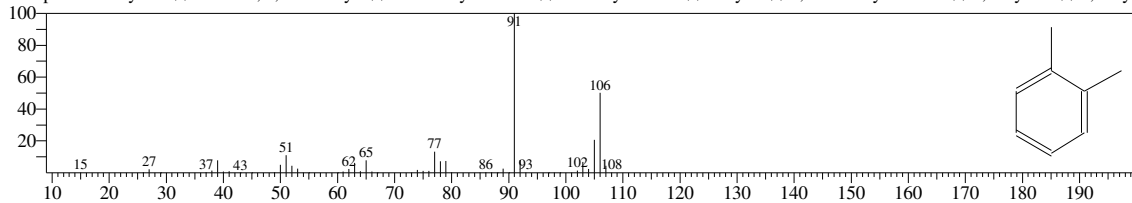
CompName:p-Xylene \$\$ Benzen, 1,4-dimethyl- \$\$ p-Dimethylbenzene \$\$ p-Xylol \$\$ 1,4-Dimethylbenzene \$\$ 1,4-Xylene \$\$ p-Methyltoluene \$\$ para-Xyl



Hit#:4 Entry:2727 Library:NIST14s.lib

SI:96 Formula:C₈H₁₀ CAS:95-47-6 MolWeight:106 RetIndex:907

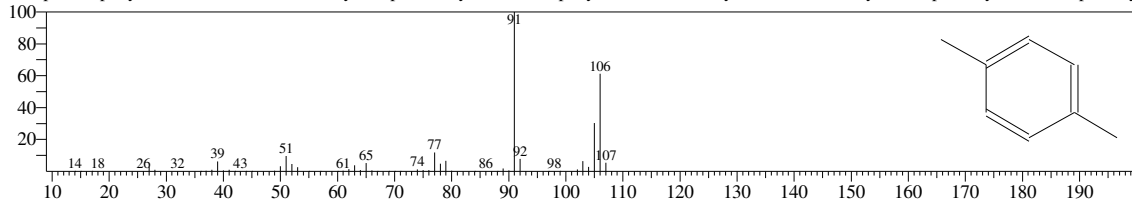
CompName:o-Xylene \$\$ Benzene, 1,2-dimethyl- \$\$ o-Dimethylbenzene \$\$ o-Methyltoluene \$\$ o-Xylol \$\$ 1,2-Dimethylbenzene \$\$ 1,2-Xylene \$\$ 3,4-Xyle



Hit#:5 Entry:2470 Library:NIST14s.lib

SI:96 Formula:C₈H₁₀ CAS:106-42-3 MolWeight:106 RetIndex:907

CompName:p-Xylene \$\$ Benzen, 1,4-dimethyl- \$\$ p-Dimethylbenzene \$\$ p-Xylol \$\$ 1,4-Dimethylbenzene \$\$ 1,4-Xylene \$\$ p-Methyltoluene \$\$ para-Xyl



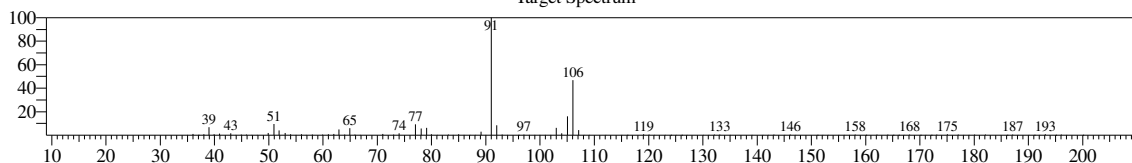
<< Target >>

Line#4 R.Time:6.170(Scan#:1135) MassPeaks:122

RawMode:Averaged 6.165-6.175(1134-1136) BasePeak:91.00(2613)

BG Mode:Calc. from Peak Group 1 - Event 1 Scan

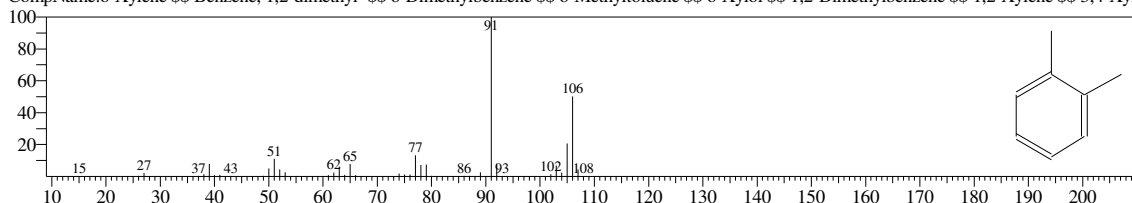
Target Spectrum



Hit#:1 Entry:2727 Library:NIST14.lib

SI:94 Formula:C₈H₁₀ CAS:95-47-6 MolWeight:106 RetIndex:907

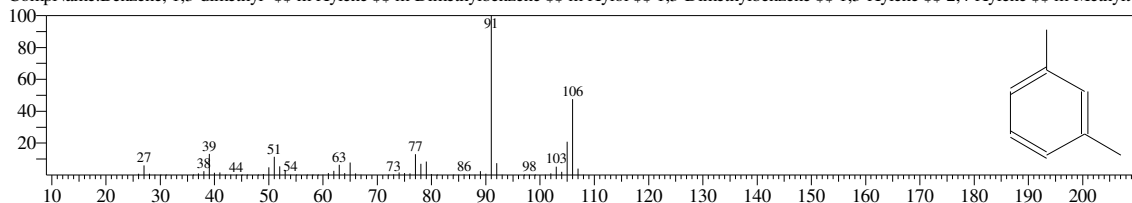
CompName:o-Xylene \$\$ Benzene, 1,2-dimethyl- \$\$ o-Dimethylbenzene \$\$ o-Methyltoluene \$\$ o-Xylol \$\$ 1,2-Dimethylbenzene \$\$ 1,2-Xylene \$\$ 3,4-Xyle



Hit#:2 Entry:2466 Library:NIST14s.lib

SI:93 Formula:C₈H₁₀ CAS:108-38-3 MolWeight:106 RetIndex:907

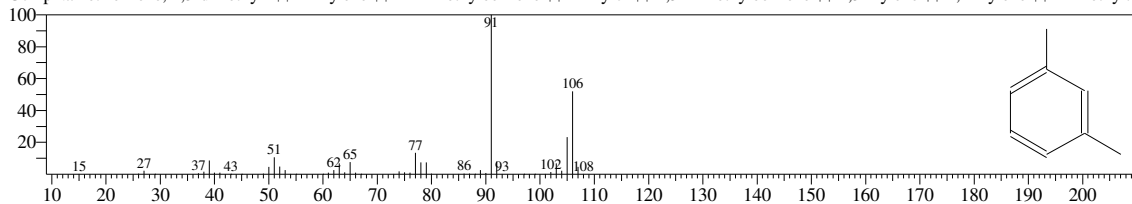
CompName:Benzen, 1,3-dimethyl- \$\$ m-Xylene \$\$ m-Dimethylbenzene \$\$ m-Xylol \$\$ 1,3-Dimethylbenzene \$\$ 1,3-Xylene \$\$ 2,4-Xylene \$\$ m-Methyltol



Hit#:3 Entry:2725 Library:NIST14.lib

SI:93 Formula:C₈H₁₀ CAS:108-38-3 MolWeight:106 RetIndex:907

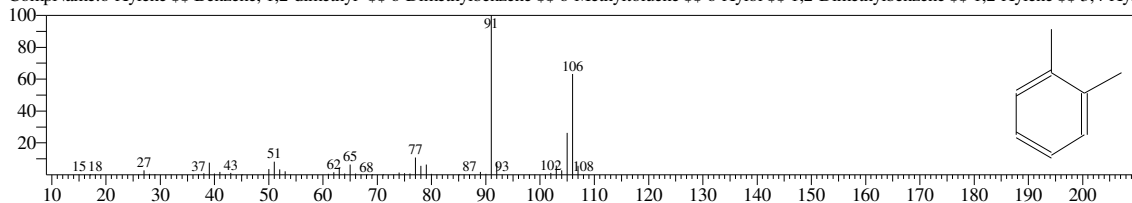
CompName:Benzen, 1,3-dimethyl- \$\$ m-Xylene \$\$ m-Dimethylbenzene \$\$ m-Xylol \$\$ 1,3-Dimethylbenzene \$\$ 1,3-Xylene \$\$ 2,4-Xylene \$\$ m-Methyltol



Hit#:4 Entry:2478 Library:NIST14s.lib

SI:93 Formula:C₈H₁₀ CAS:95-47-6 MolWeight:106 RetIndex:907

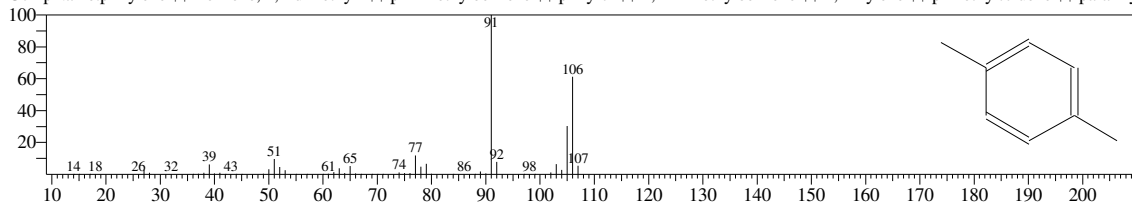
CompName:o-Xylene \$\$ Benzene, 1,2-dimethyl- \$\$ o-Dimethylbenzene \$\$ o-Methyltoluene \$\$ o-Xylol \$\$ 1,2-Dimethylbenzene \$\$ 1,2-Xylene \$\$ 3,4-Xyle



Hit#:5 Entry:2470 Library:NIST14s.lib

SI:93 Formula:C₈H₁₀ CAS:106-42-3 MolWeight:106 RetIndex:907

CompName:p-Xylene \$\$ Benzene, 1,4-dimethyl- \$\$ p-Dimethylbenzene \$\$ p-Xylol \$\$ 1,4-Dimethylbenzene \$\$ 1,4-Xylene \$\$ p-Methyltoluene \$\$ para-Xyl



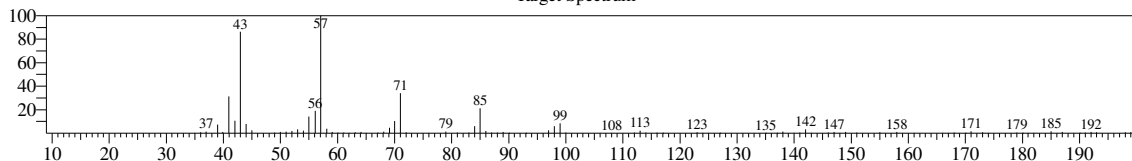
<< Target >>

Line#:5 R.Time:9.325(Scan#:1766) MassPeaks:93

RawMode:Averaged 9.320-9.330(1765-1767) BasePeak:57.05(1116)

BG Mode:Calc. from Peak Group 1 - Event 1 Scan

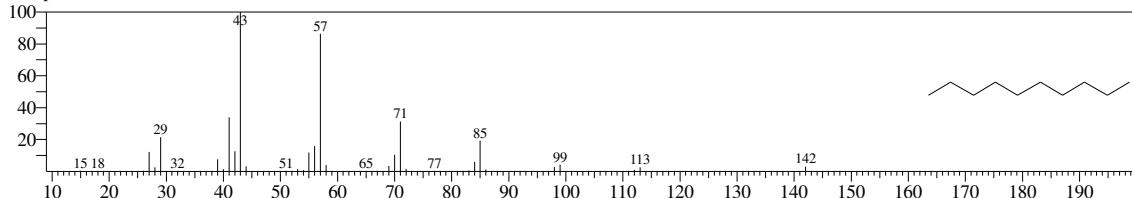
Target Spectrum



Hit#:1 Entry:7907 Library:NIST14s.lib

SI:92 Formula:C10H22 CAS:124-18-5 MolWeight:142 RetIndex:1015

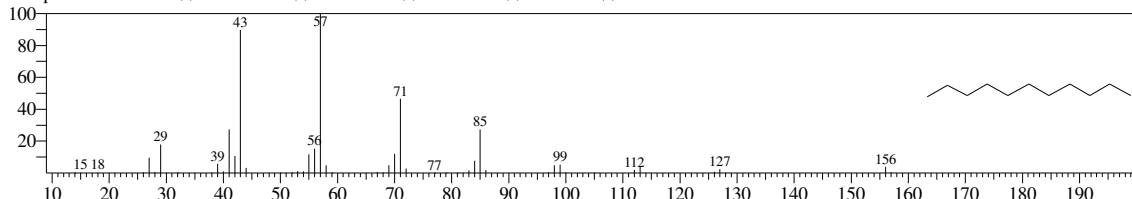
CompName:Decane \$\$ n-Decane \$\$ n-C10H22 \$\$ UN 2247 \$\$



Hit#:2 Entry:19287 Library:NIST14s.lib

SI:91 Formula:C11H24 CAS:1120-21-4 MolWeight:156 RetIndex:1115

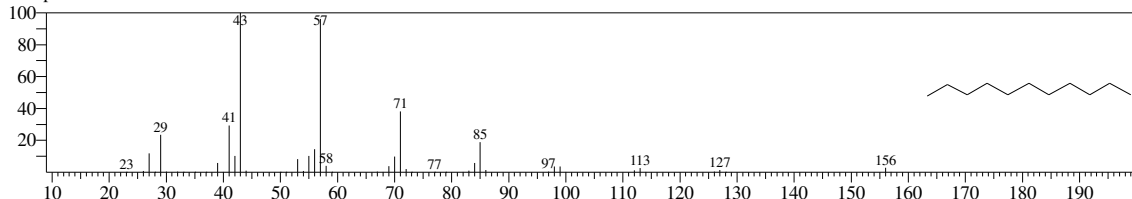
CompName:Undecane \$\$ n-Undecane \$\$ Hendecane \$\$ n-C11H24 \$\$ UN 2330 \$\$



Hit#:3 Entry:10781 Library:NIST14s.lib

SI:91 Formula:C11H24 CAS:1120-21-4 MolWeight:156 RetIndex:1115

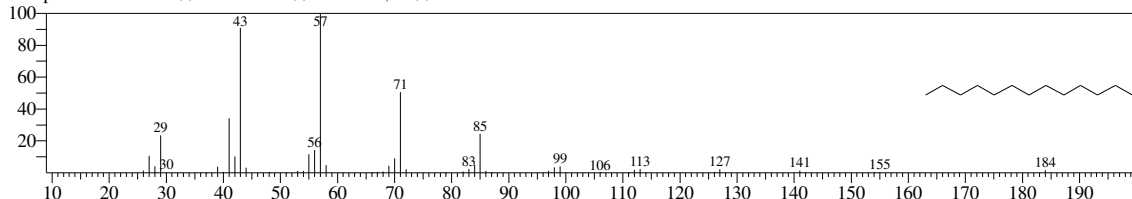
CompName:Undecane \$\$ n-Undecane \$\$ Hendecane \$\$ n-C11H24 \$\$ UN 2330 \$\$



Hit#:4 Entry:36063 Library:NIST14s.lib

SI:91 Formula:C13H28 CAS:629-50-5 MolWeight:184 RetIndex:1313

CompName:Tridecane \$\$ n-Tridecane \$\$ Tridecane, n- \$\$



Hit#:5 Entry:10790 Library:NIST14s.lib

SI:91 Formula:C11H24 CAS:1120-21-4 MolWeight:156 RetIndex:1115

CompName:Undecane \$\$ n-Undecane \$\$ Hendecane \$\$ n-C11H24 \$\$ UN 2330 \$\$

