

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111324\
 Data File : VN084835.D
 Acq On : 13 Nov 2024 18:47
 Operator : JC\MD
 Sample : P4791-02 10X
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 HINCHMAN-OILY-WATER

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N103024W.M

Title : SW846 8260

Signal : TIC: VN084835.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.289	50	57	65	rBV3	17098	41187	4.76%	0.728%
2	2.901	146	161	162	rBV3	13354	40100	4.63%	0.709%
3	4.442	410	423	434	rBV2	30978	91408	10.56%	1.616%
4	5.489	594	601	606	rBV3	3642	8997	1.04%	0.159%
5	8.159	1046	1055	1060	rBV	123758	301015	34.76%	5.321%
6	8.218	1060	1065	1076	rVB	216430	487261	56.27%	8.612%
7	8.571	1116	1125	1135	rBV	137780	304792	35.20%	5.387%
8	8.959	1185	1191	1199	rVB3	8084	16058	1.85%	0.284%
9	9.094	1205	1214	1224	rBV	315710	650637	75.14%	11.500%
10	9.183	1224	1229	1236	rVB4	6171	13410	1.55%	0.237%
11	10.441	1438	1443	1451	rVB3	7860	12850	1.48%	0.227%
12	10.565	1456	1464	1472	rBV	472175	865922	100.00%	15.305%
13	11.865	1676	1685	1693	rBV	428495	759826	87.75%	13.430%
14	12.847	1845	1852	1860	rVB	355522	597666	69.02%	10.564%
15	13.600	1971	1980	1985	rBV5	4735	8806	1.02%	0.156%
16	13.665	1985	1991	1994	rVV3	7049	12565	1.45%	0.222%
17	13.700	1994	1997	2002	rVV4	11391	18064	2.09%	0.319%
18	13.788	2002	2012	2019	rVV	469420	785266	90.69%	13.880%
19	13.853	2019	2023	2032	rVB5	46366	87617	10.12%	1.549%
20	14.241	2076	2089	2099	rBV	92544	154774	17.87%	2.736%
21	14.488	2125	2131	2135	rVV2	76725	127951	14.78%	2.262%
22	14.523	2135	2137	2146	rVB4	17444	27278	3.15%	0.482%
23	14.770	2175	2179	2184	rVV4	13455	26352	3.04%	0.466%
24	14.888	2194	2199	2205	rBV4	9304	18419	2.13%	0.326%
25	14.959	2205	2211	2219	rVB	28401	48942	5.65%	0.865%
26	15.106	2230	2236	2245	rVB4	10972	20724	2.39%	0.366%
27	15.282	2261	2266	2269	rBV4	6329	9706	1.12%	0.172%
28	15.323	2269	2273	2285	rVB6	8073	18629	2.15%	0.329%
29	15.429	2285	2291	2294	rBV4	6626	13817	1.60%	0.244%
30	15.512	2300	2305	2313	rVV6	8652	14575	1.68%	0.258%
31	15.600	2313	2320	2322	rVV5	6296	12026	1.39%	0.213%
32	15.635	2322	2326	2333	rVB4	9068	16783	1.94%	0.297%
33	16.012	2384	2390	2396	rVB6	6451	13863	1.60%	0.245%
34	16.706	2501	2508	2516	rBV3	11127	30325	3.50%	0.536%

Sum of corrected areas: 5657611

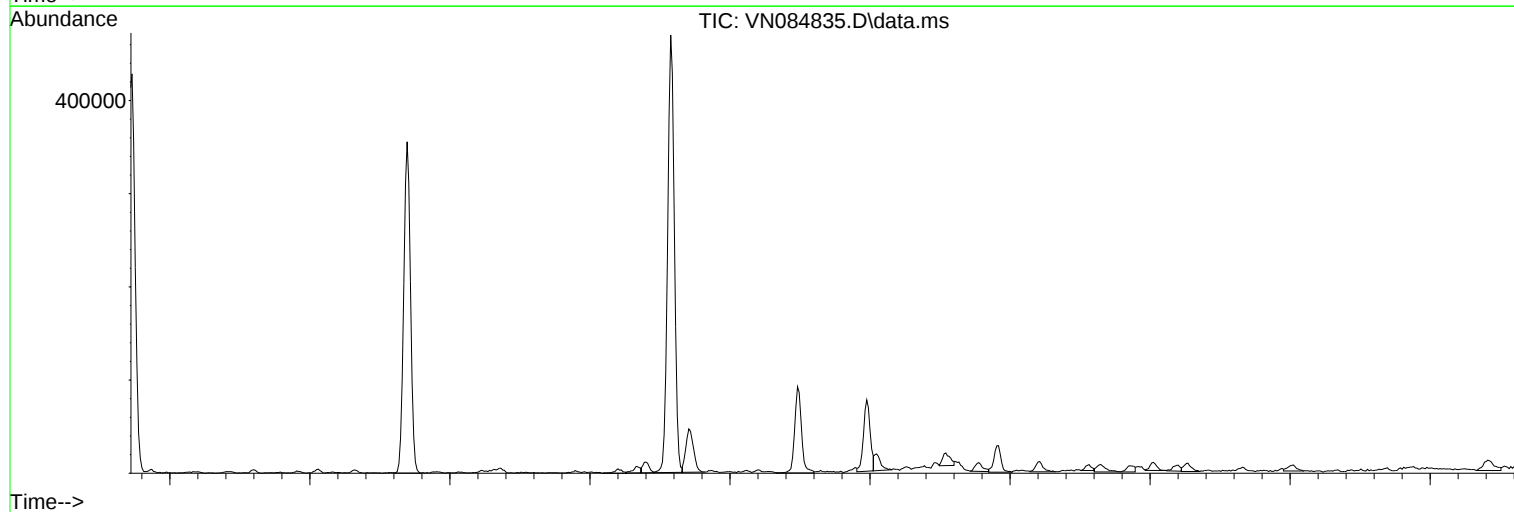
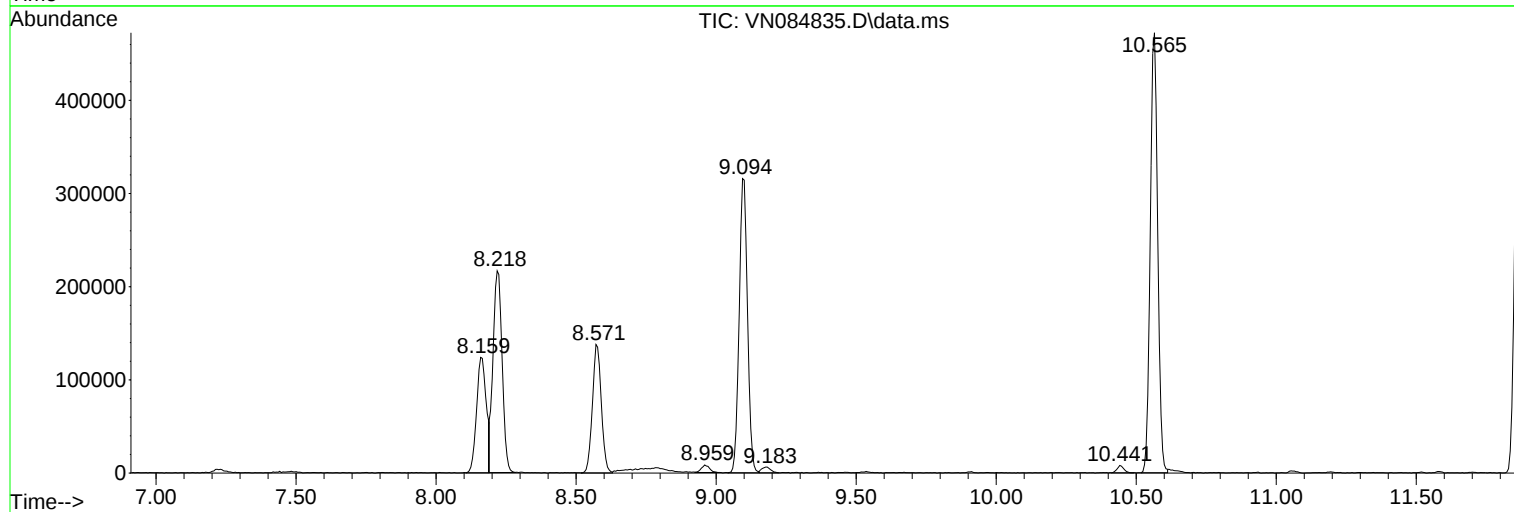
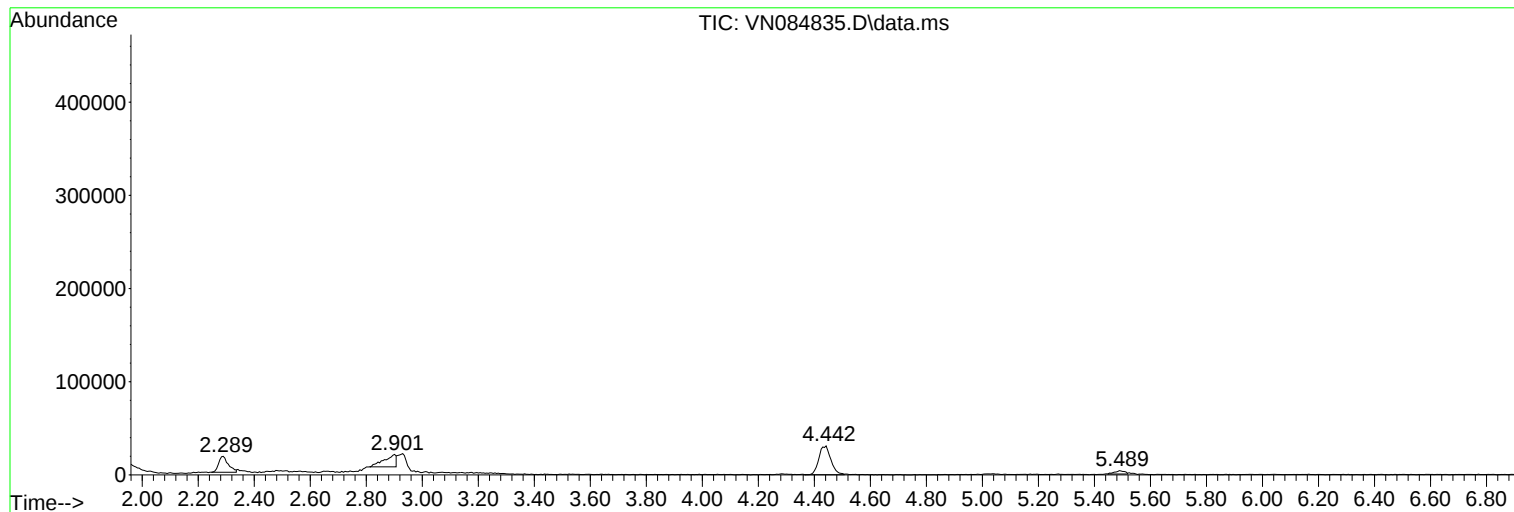
Instrument :
MSVOA_N
ClientSampleId :
HINCHMAN-OILY-WATER

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N103024W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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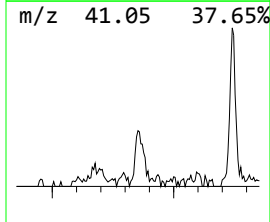
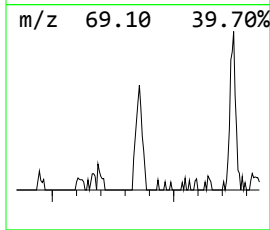
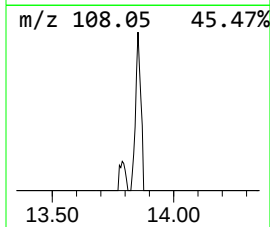
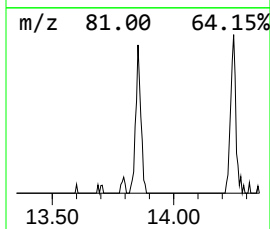
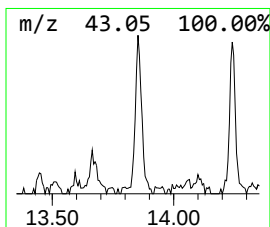
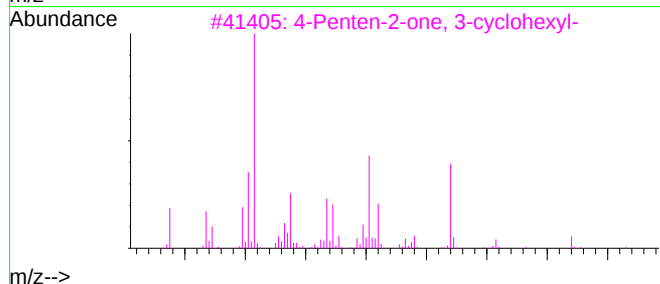
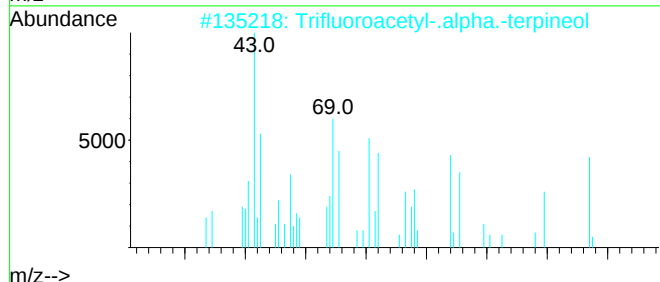
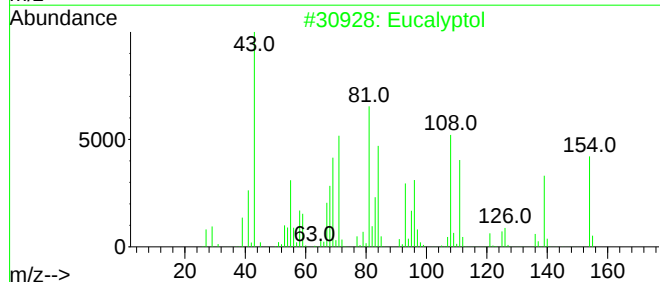
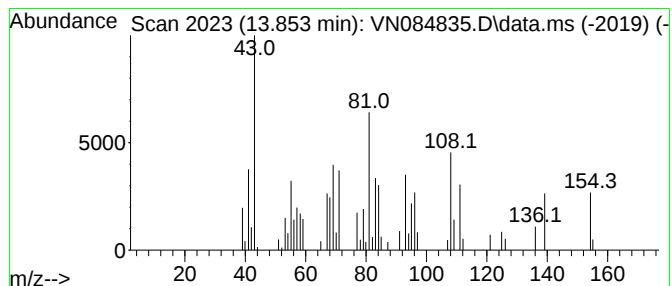
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N103024W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Eucalyptol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.853	5.58 ug/l	87617	1,4-Dichlorobenzene-d4	13.788

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eucalyptol	154	C10H18O	000470-82-6	94
2	Trifluoroacetyl-.alpha.-terpineol	250	C12H17F3O2	1000058-17-6	42
3	4-Penten-2-one, 3-cyclohexyl-	166	C11H18O	055702-54-0	22
4	3-Ethylidene- heptan-2,6-dione	154	C9H14O2	1000223-47-9	22
5	5-Hepten-2-one, 6-methyl-	126	C8H14O	000110-93-0	22



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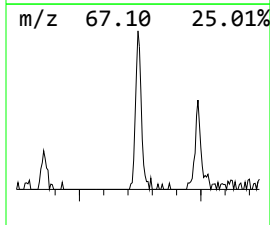
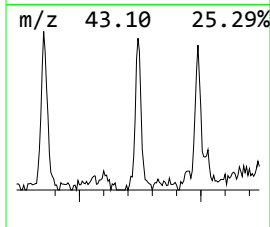
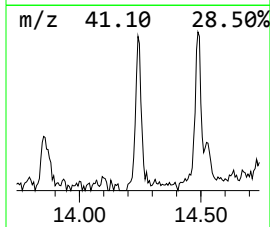
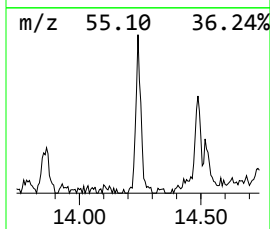
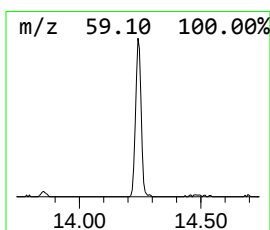
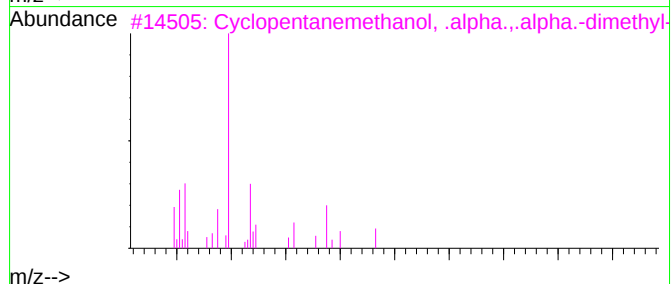
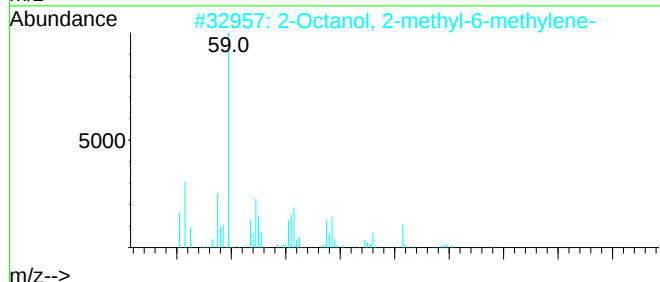
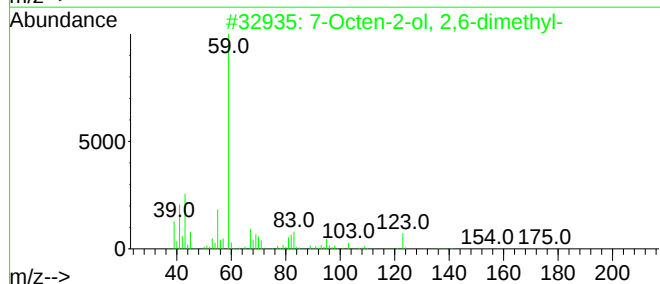
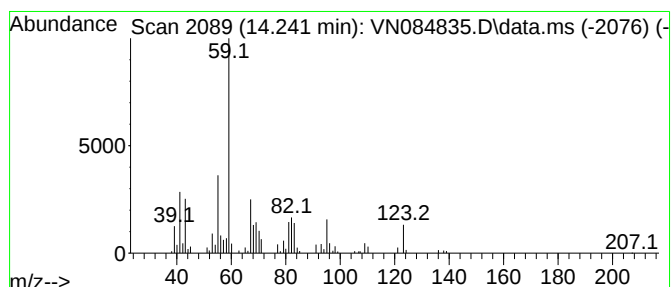
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TIC Integration Parameters: LSCINT.P

Peak Number 2 7-Octen-2-ol, 2,6-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.241	9.85 ug/l	154774	1,4-Dichlorobenzene-d4	13.788

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-Octen-2-ol, 2,6-dimethyl-	156	C10H20O	018479-58-8	90
2	2-Octanol, 2-methyl-6-methylene-	156	C10H20O	018479-59-9	72
3	Cyclopentanemethanol, .alpha.,.a...	128	C8H16O	001462-06-2	53
4	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	000498-81-7	42
5	3-(2-Hydroxy-2-methyl-propyl)-cy...	168	C10H16O2	1000185-60-2	42



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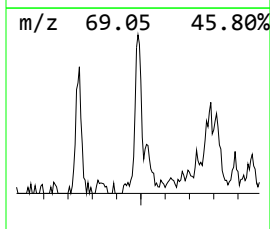
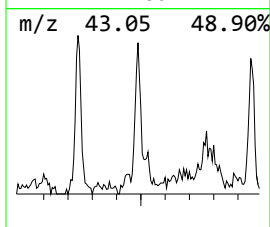
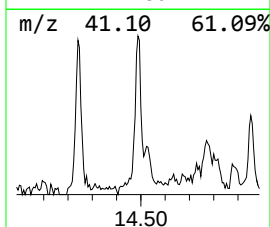
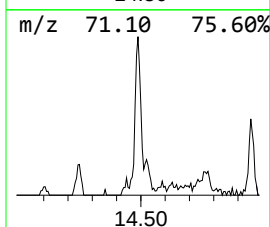
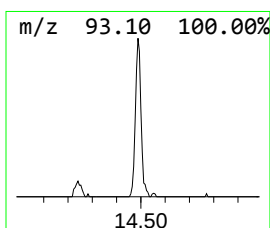
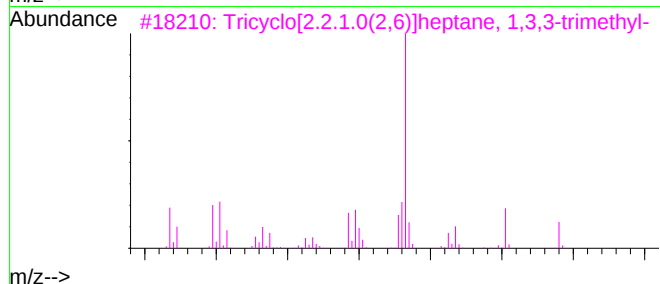
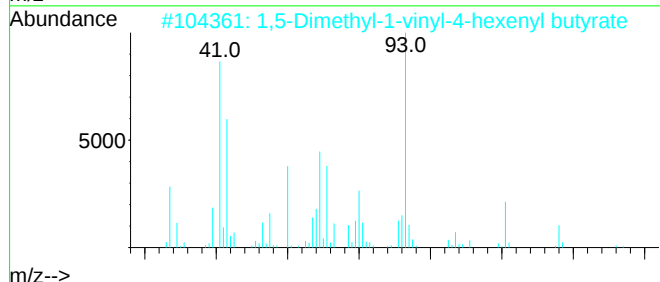
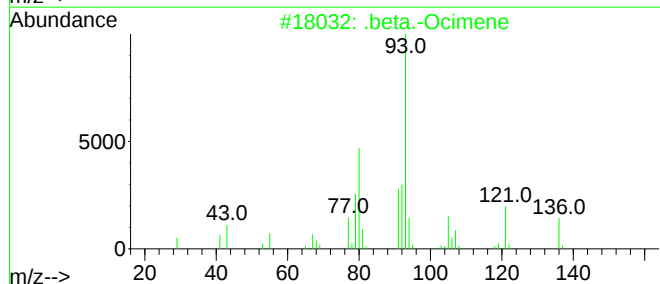
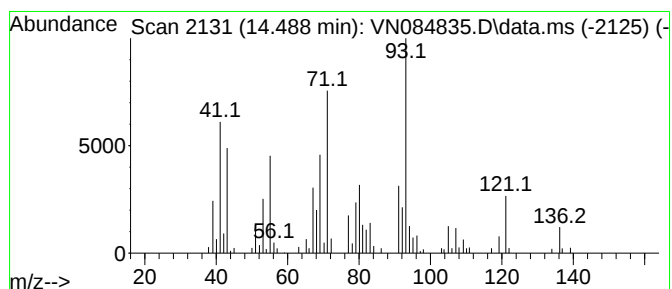
Instrument :
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ClientSampleId :
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 .beta.-Ocimene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.488	8.15 ug/l	127951	1,4-Dichlorobenzene-d4	13.788	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.beta.-Ocimene	136	C10H16	013877-91-3	70
2	1,5-Dimethyl-1-vinyl-4-hexenyl b...	224	C14H24O2	000078-36-4	58
3	Tricyclo[2.2.1.0(2,6)]heptane, 1...	136	C10H16	000488-97-1	55
4	Hexanoic acid, 1-ethenyl-1,5-dim...	252	C16H28O2	007779-23-9	52
5	Butanoic acid, 3-methyl-, 1-ethe...	238	C15H26O2	001118-27-0	52



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Eucalyptol	13.853	5.6	ug/l	87617	4	13.788	785266	50.0
7-Octen-2-ol, 2...	14.241	9.8	ug/l	154774	4	13.788	785266	50.0
.beta.-Ocimene	14.488	8.2	ug/l	127951	4	13.788	785266	50.0