

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX030521\
 Data File : VX020996.D
 Acq On : 03 Mar 2021 14:09
 Operator : JC/MD
 Sample : M1536-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-1

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022421W.M
 Title : SW846 8260

Signal : TIC: VX020996.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.393	46	52	57	rBV	364698	406927	31.79%	3.614%
2	1.434	57	59	70	rVB	21222	27978	2.19%	0.248%
3	1.681	96	101	104	rBV2	46147	63905	4.99%	0.568%
4	1.717	104	107	114	rVB	104484	131382	10.26%	1.167%
5	1.952	142	147	152	rBV	28952	40283	3.15%	0.358%
6	2.028	157	160	168	rVB	13085	17657	1.38%	0.157%
7	2.175	178	185	193	rBV	51662	89876	7.02%	0.798%
8	2.252	193	198	207	rVB2	10835	19020	1.49%	0.169%
9	2.358	209	216	226	rBV	787260	1246161	97.34%	11.068%
10	2.570	245	252	263	rBV	28180	54633	4.27%	0.485%
11	2.993	316	324	334	rVB3	18858	44844	3.50%	0.398%
12	3.670	430	439	451	rBV	92862	220910	17.26%	1.962%
13	4.564	580	591	600	rBV	71383	200223	15.64%	1.778%
14	5.464	732	744	756	rBV2	146840	407734	31.85%	3.621%
15	5.629	761	772	783	rBV	222557	615896	48.11%	5.470%
16	6.029	829	840	849	rBV	153157	401522	31.36%	3.566%
17	6.158	856	862	874	rVB2	28630	78172	6.11%	0.694%
18	6.299	874	886	900	rBV	339773	933949	72.95%	8.295%
19	6.829	965	976	998	rBV	380807	909829	71.07%	8.081%
20	7.229	1039	1044	1052	rVB3	24802	52878	4.13%	0.470%
21	7.441	1072	1080	1088	rBV4	5770	13355	1.04%	0.119%
22	7.706	1117	1125	1141	rBV	247544	491256	38.37%	4.363%
23	8.023	1172	1179	1188	rBV2	22131	39120	3.06%	0.347%
24	8.470	1248	1255	1263	rBV	31469	55734	4.35%	0.495%
25	8.694	1286	1293	1302	rBV	819811	1280205	100.00%	11.371%
26	9.482	1421	1427	1436	rBV2	38108	64914	5.07%	0.577%
27	9.594	1442	1446	1453	rVB2	20019	29112	2.27%	0.259%
28	9.771	1468	1476	1483	rBV	19944	29190	2.28%	0.259%
29	10.012	1508	1517	1522	rBV	64455	94299	7.37%	0.838%
30	10.094	1525	1531	1542	rVV	822446	1107553	86.51%	9.837%
31	10.682	1627	1631	1637	rVB2	10043	14511	1.13%	0.129%
32	10.888	1660	1666	1671	rBV4	11283	17320	1.35%	0.154%
33	11.053	1689	1694	1700	rBV	10430	14487	1.13%	0.129%
34	11.118	1700	1705	1715	rBV2	674162	910664	71.13%	8.088%
35	11.653	1791	1796	1801	rVB	13037	16167	1.26%	0.144%
36	12.059	1860	1865	1873	rBV	796087	986980	77.10%	8.766%

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Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Peak separation: 5

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Title : SW846 8260

37	12.283	1895	1903	1908	rBV	23778	30819	2.41%	0.274%
38	12.741	1977	1981	1987	rVB4	7486	13209	1.03%	0.117%
39	12.883	1998	2005	2011	rVB2	8718	13108	1.02%	0.116%
40	13.330	2076	2081	2087	rBV	56474	73125	5.71%	0.649%

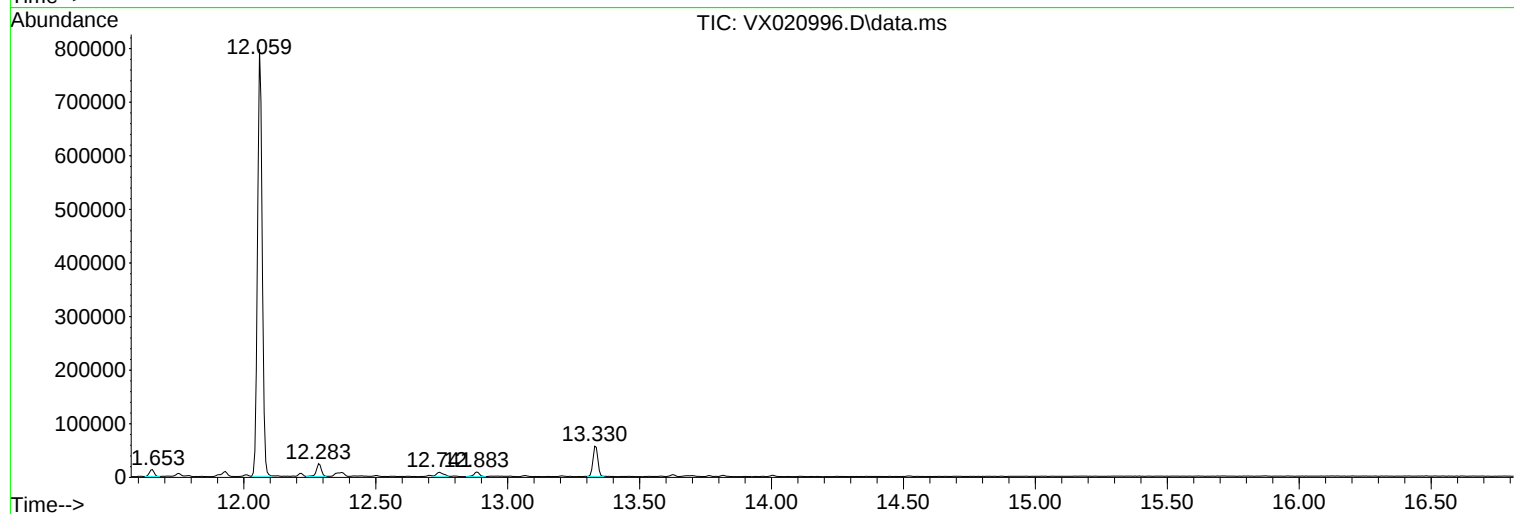
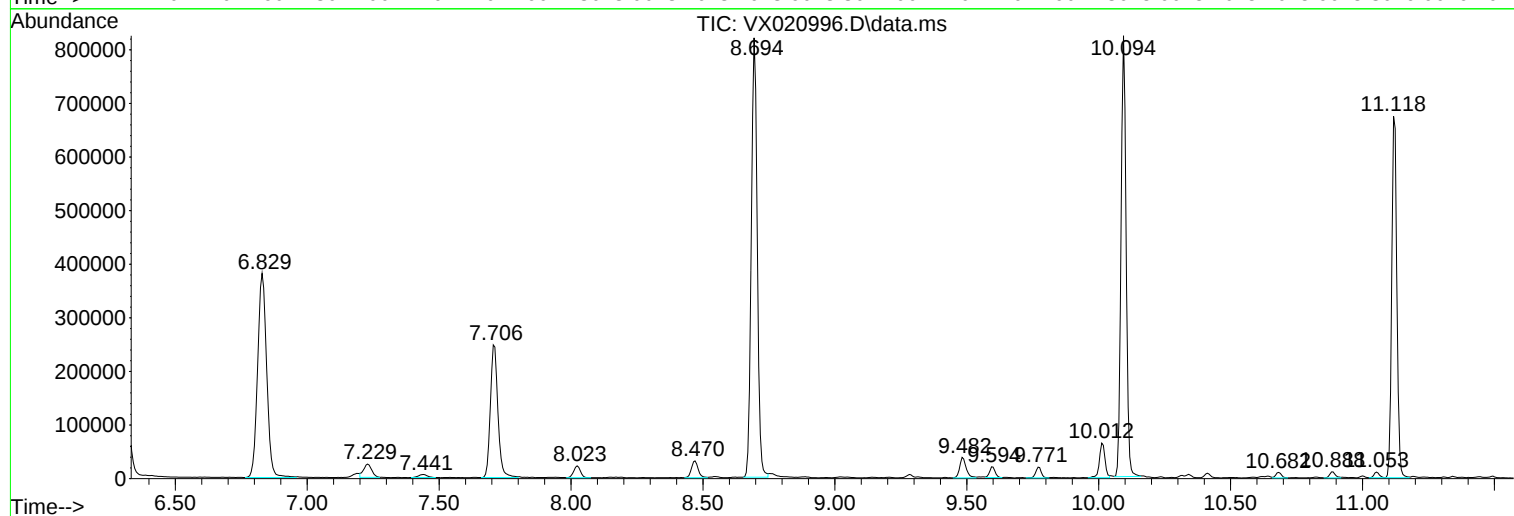
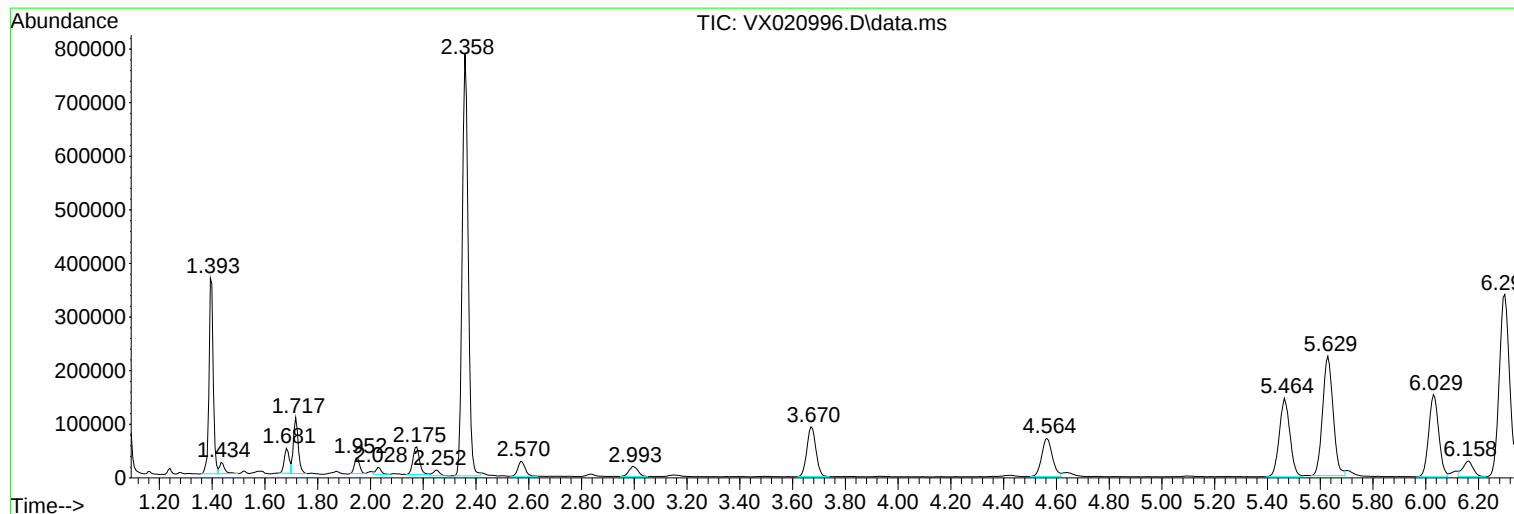
Sum of corrected areas: 11258907

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022421W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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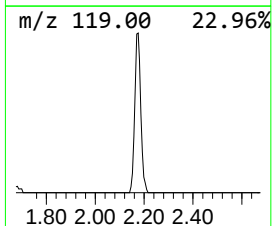
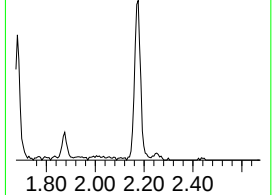
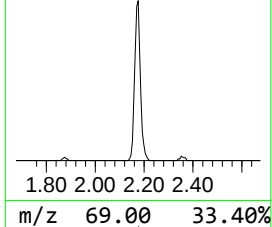
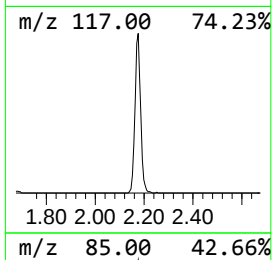
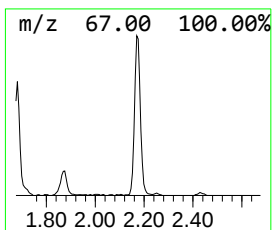
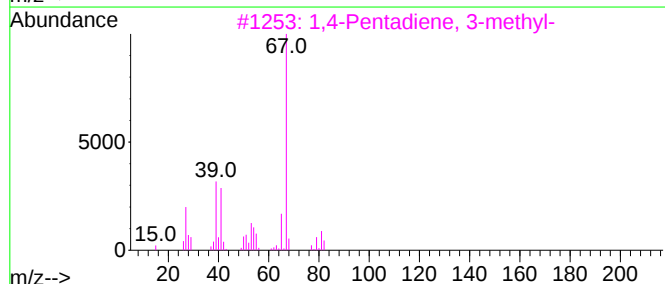
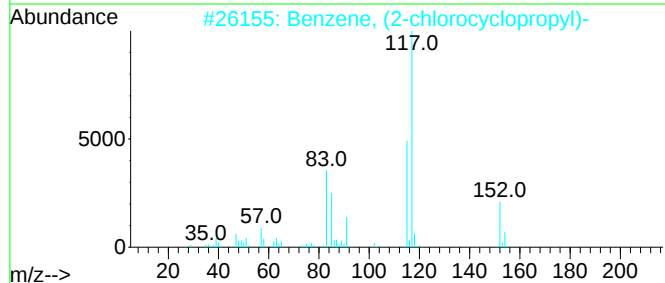
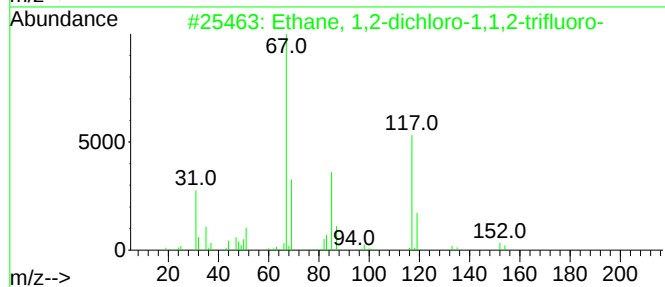
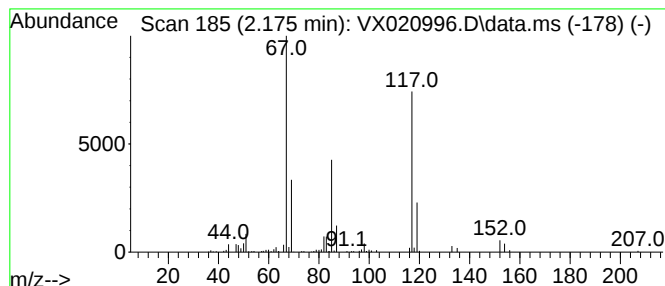
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022421W.M
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TIC Integration Parameters: LSCINT.P

Peak Number 1 Ethane, 1,2-dichloro-1,1,2-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.175	7.30 ug/l	89876	Pentafluorobenzene	5.629

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethane, 1,2-dichloro-1,1,2-trifl...	152	C2HCl2F3	000354-23-4	91
2			Benzene, (2-chlorocyclopropyl)-	152	C9H9Cl	1000327-31-4	12
3			1,4-Pentadiene, 3-methyl-	82	C6H10	001115-08-8	9
4			Methyl 2-butynoate	98	C5H6O2	023326-27-4	9
5			Cyclopropane, 1,2-dimethyl-3-met...	82	C6H10	062338-02-7	9



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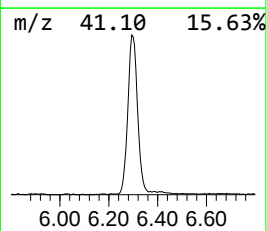
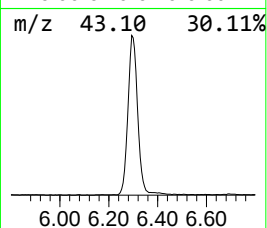
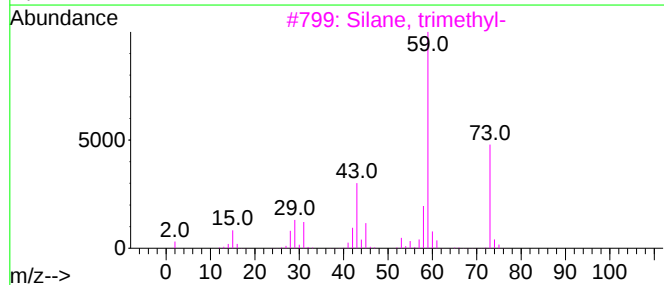
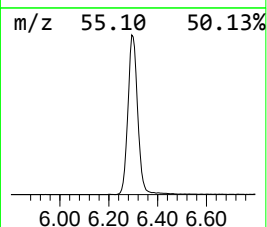
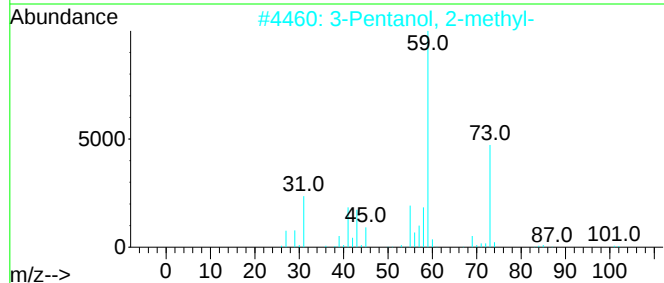
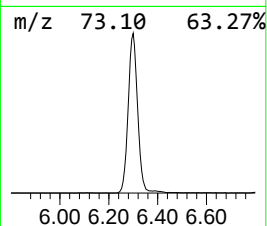
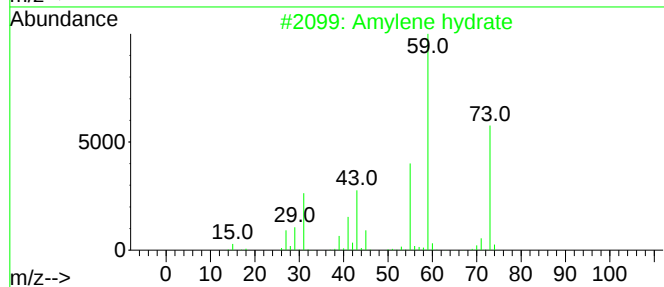
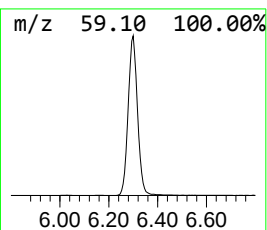
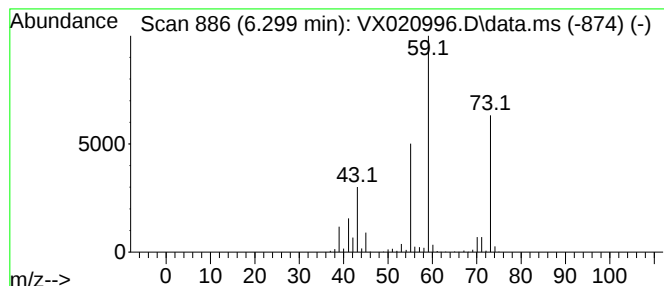
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Peak Number 2 Amylene hydrate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.299	51.33 ug/l	933949	1,4-Difluorobenzene	6.829

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Amylene hydrate	88	C5H12O	000075-85-4	83
2			3-Pentanol, 2-methyl-	102	C6H14O	000565-67-3	39
3			Silane, trimethyl-	74	C3H10Si	000993-07-7	10
4			Propane, 2-methoxy-	74	C4H10O	000598-53-8	9
5			3-Hexanol	102	C6H14O	000623-37-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethane, 1,2-dic...	2.175	7.3	ug/l	89876	1	5.629	615896	50.0
Amylene hydrate	6.299	51.3	ug/l	933949	2	6.829	909829	50.0