

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061521\
 Data File : VX022707.D
 Acq On : 15 Jun 2021 18:31
 Operator : JC/MD
 Sample : M2712-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 10 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\82X061021W.M
 Title : SW846 8260

Signal : TIC: VX022707.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.221	19	22	26	rBV	5852	6017	1.26%	0.163%
2	1.374	41	47	52	rBV	135781	132903	27.83%	3.593%
3	1.416	52	54	58	rVB	7502	7875	1.65%	0.213%
4	1.660	89	94	96	rVV2	18752	24998	5.23%	0.676%
5	1.691	96	99	106	rVB	32571	41861	8.77%	1.132%
6	1.916	132	136	142	rBV	11501	15601	3.27%	0.422%
7	2.002	147	150	158	rVB2	3554	5039	1.06%	0.136%
8	2.142	167	173	179	rBV	15470	24923	5.22%	0.674%
9	2.215	181	185	190	rBV2	3615	5659	1.18%	0.153%
10	2.319	196	202	211	rBV	294735	459672	96.26%	12.429%
11	2.971	303	309	318	rBV2	3580	8917	1.87%	0.241%
12	3.617	405	415	426	rVB	22985	51079	10.70%	1.381%
13	4.495	550	559	570	rBV2	22661	61893	12.96%	1.673%
14	5.397	696	707	718	rBV	40807	114633	24.00%	3.100%
15	5.562	723	734	745	rVV	69422	190288	39.85%	5.145%
16	5.964	790	800	810	rBV2	52308	136818	28.65%	3.699%
17	6.099	817	822	832	rVB3	10616	27459	5.75%	0.742%
18	6.245	834	846	859	rBV	178214	477556	100.00%	12.912%
19	6.769	922	932	942	rBV	114822	277591	58.13%	7.506%
20	7.178	994	999	1007	rVB5	6406	13458	2.82%	0.364%
21	7.385	1026	1033	1039	rBV3	2590	5156	1.08%	0.139%
22	7.659	1070	1078	1097	rBV	113321	226442	47.42%	6.123%
23	7.976	1121	1130	1135	rBV3	6315	10692	2.24%	0.289%
24	8.427	1198	1204	1213	rBV	14134	24840	5.20%	0.672%
25	8.653	1234	1241	1248	rBV	250274	385473	80.72%	10.423%
26	9.452	1365	1372	1379	rBV2	19201	31788	6.66%	0.859%
27	9.561	1386	1390	1394	rVB2	7478	10279	2.15%	0.278%
28	9.732	1414	1418	1424	rVB3	6532	8476	1.77%	0.229%
29	9.976	1446	1458	1463	rBV	25074	36777	7.70%	0.994%
30	10.055	1466	1471	1478	rVB	239162	321422	67.31%	8.691%
31	10.378	1519	1524	1532	rBV2	4435	6182	1.29%	0.167%
32	11.018	1623	1629	1635	rBV2	4936	6620	1.39%	0.179%
33	11.085	1635	1640	1646	rBV	185544	237785	49.79%	6.429%
34	12.024	1789	1794	1804	rBV	220771	270378	56.62%	7.311%
35	12.250	1827	1831	1836	rBV2	6639	7736	1.62%	0.209%
36	13.298	1997	2003	2007	rBV3	19055	24145	5.06%	0.653%

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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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061021W.M
Title : SW846 8260

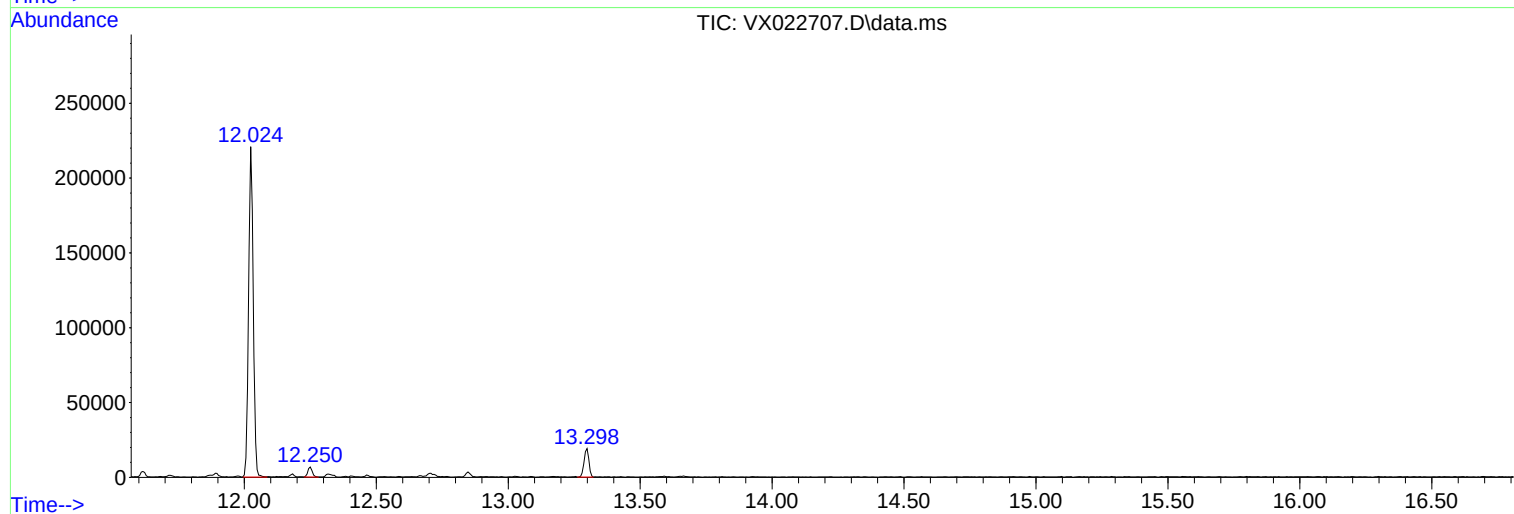
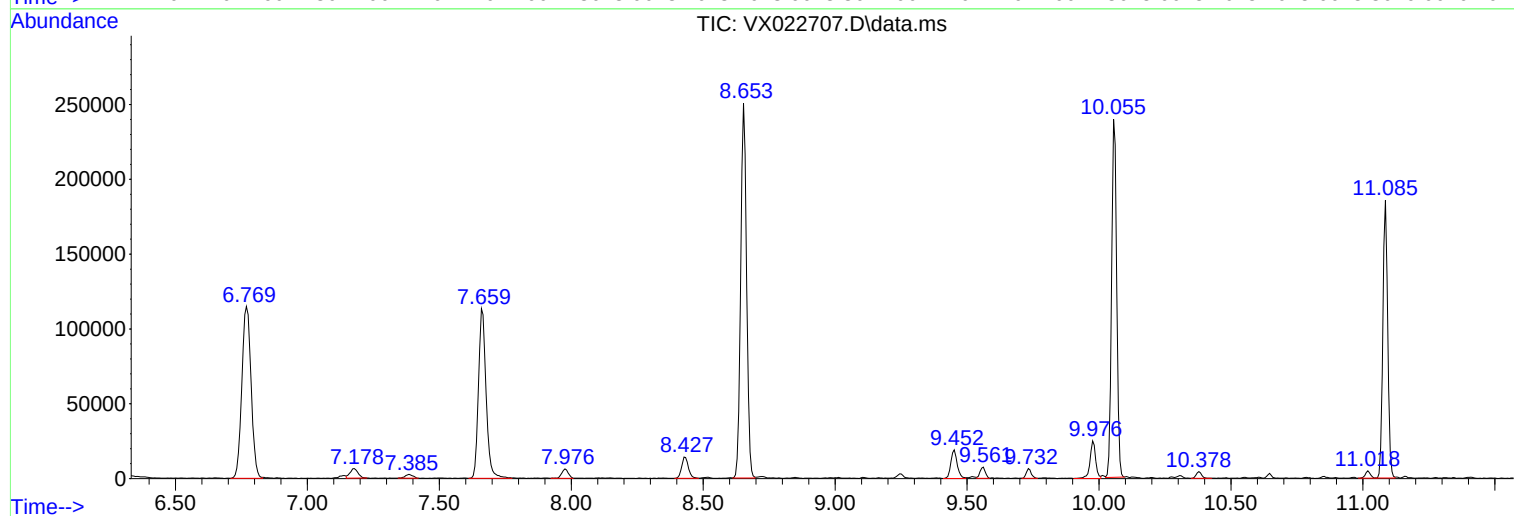
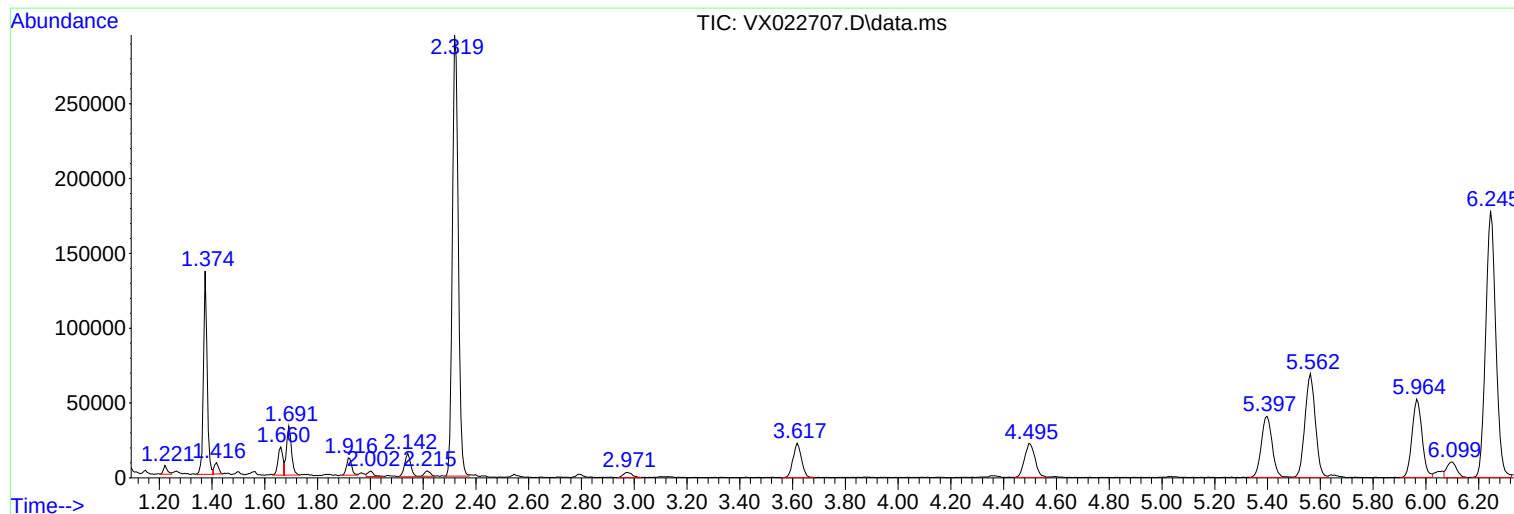
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061021W.M
Quant Title : SW846 8260

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



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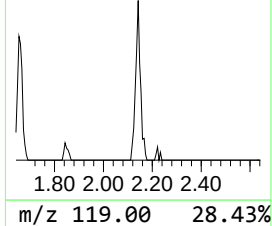
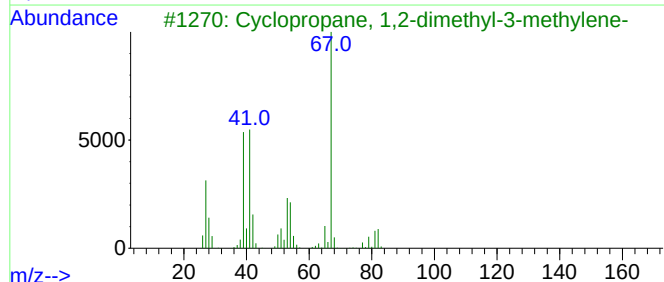
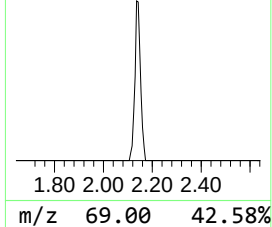
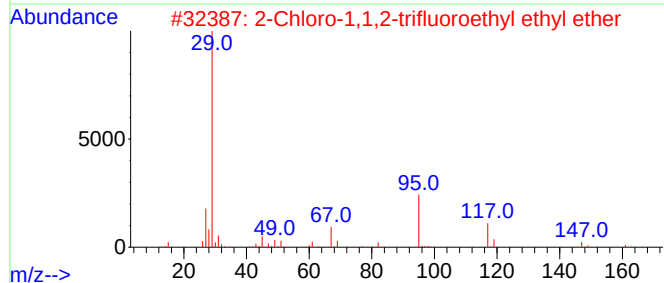
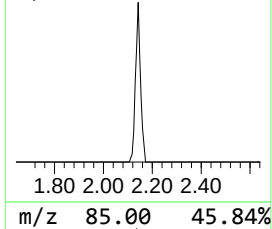
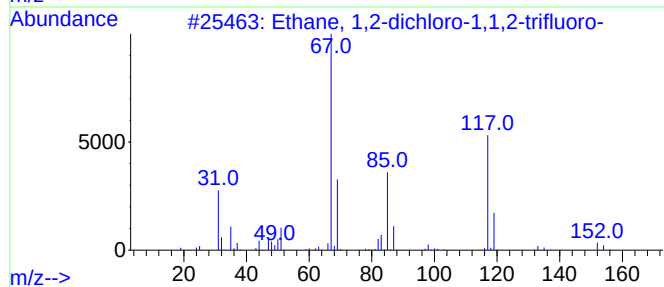
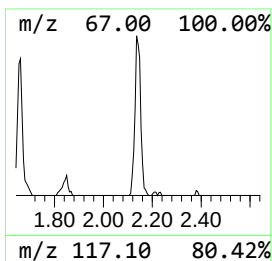
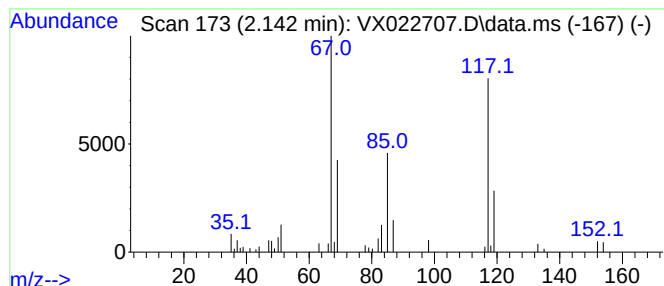
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.142	6.55 ug/l	24923	Pentafluorobenzene	5.562

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			2-Chloro-1,1,2-trifluoroethyl et...	162	C4H6ClF3O	000310-71-4	36
3			Cyclopropane, 1,2-dimethyl-3-met...	82	C6H10	062338-02-7	14
4			1,4-Pentadiene, 3-methyl-	82	C6H10	001115-08-8	14
5			1-Butyne, 3,3-dimethyl-	82	C6H10	000917-92-0	10



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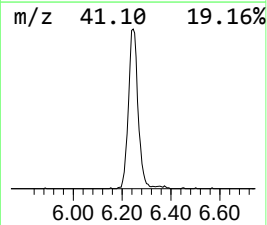
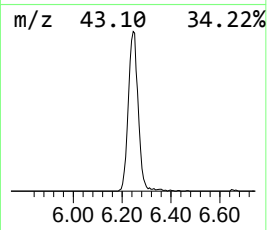
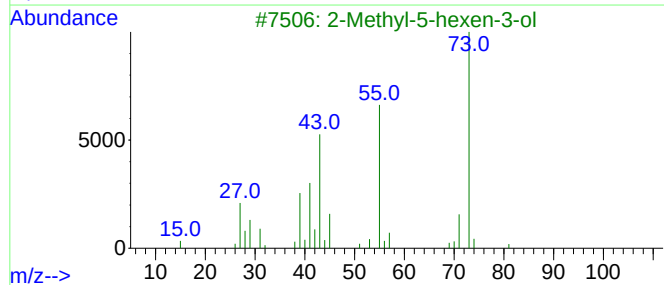
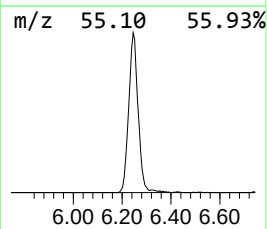
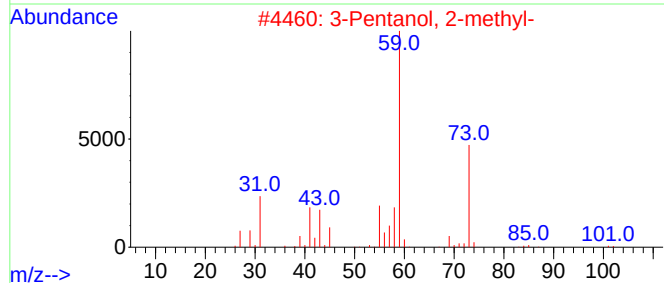
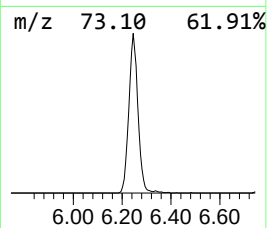
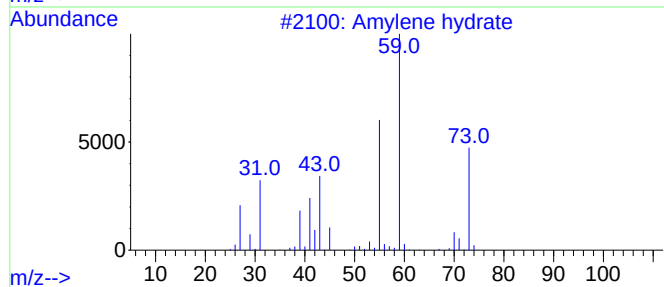
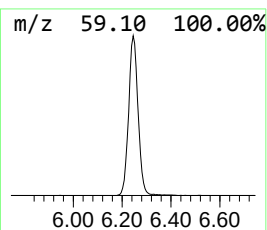
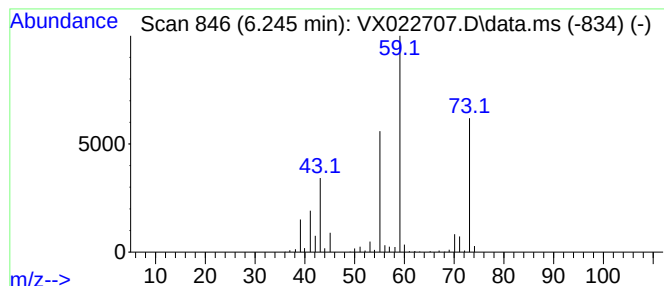
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 Amylene hydrate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.245	86.02 ug/l	477556	1,4-Difluorobenzene	6.769

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			2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	10
4			Silane, trimethyl-	74	C3H10Si	000993-07-7	9
5			Propanoic acid, 2-hydroxy-2-methyl-	104	C4H8O3	000594-61-6	9



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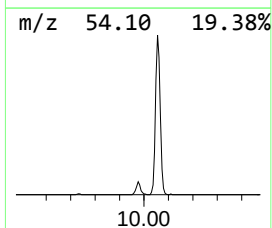
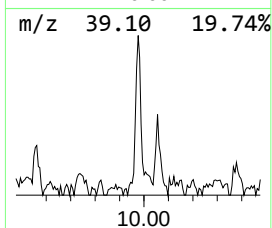
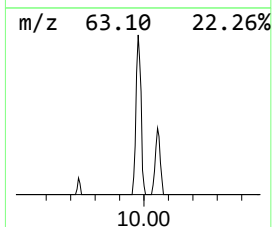
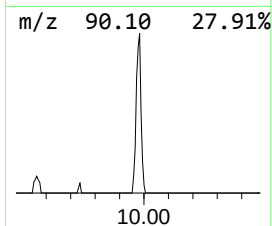
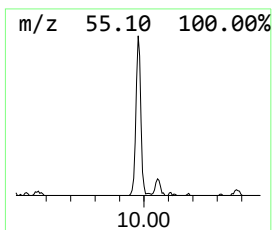
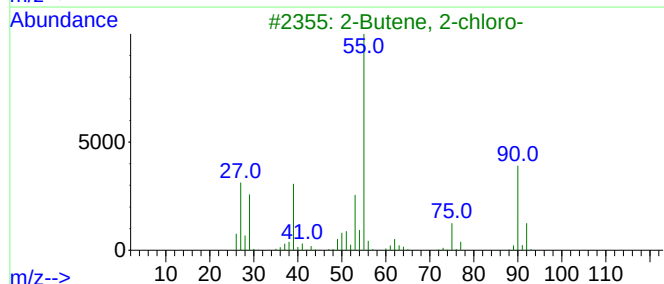
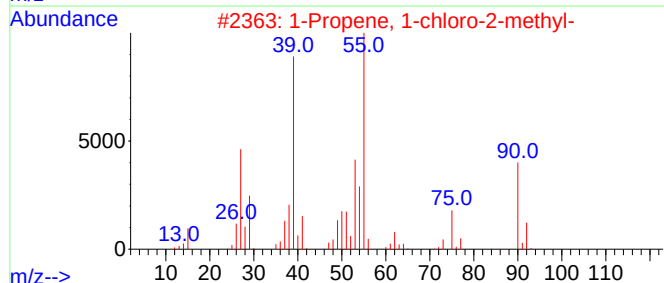
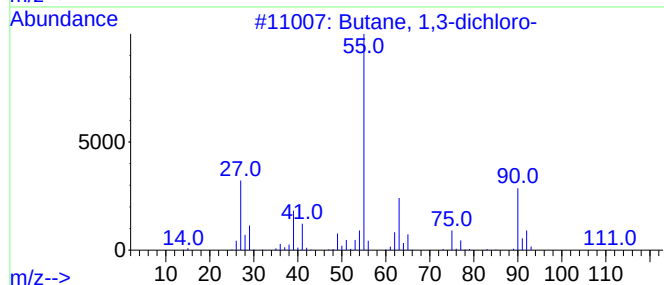
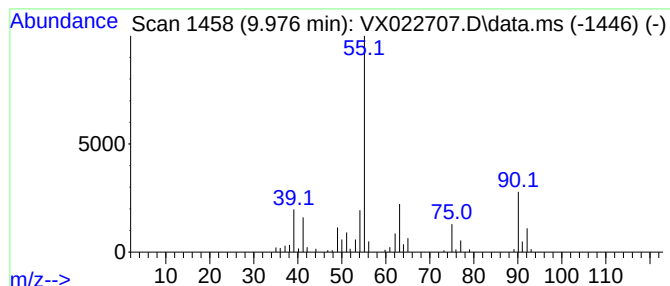
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Butane, 1,3-dichloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.976	5.72 ug/l	36777	Chlorobenzene-d5	10.055

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 1,3-dichloro-	126	C4H8Cl2	001190-22-3	80
2			1-Propene, 1-chloro-2-methyl-	90	C4H7Cl	000513-37-1	64
3			2-Butene, 2-chloro-	90	C4H7Cl	004461-41-0	59
4			1-Propene, 3-chloro-2-methyl-	90	C4H7Cl	000563-47-3	58
5			2-Butene, 1-chloro-	90	C4H7Cl	000591-97-9	58



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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethane, 1,2-dic...	2.142	6.5	ug/l	24923	1	5.562	190288	50.0
Amylene hydrate	6.245	86.0	ug/l	477556	2	6.769	277591	50.0
Butane, 1,3-dic...	9.976	5.7	ug/l	36777	3	10.055	321422	50.0