

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN111820\
 Data File : VN064720.D
 Acq On : 18 Nov 2020 20:36
 Operator : JC/MD
 Sample : L4783-20 40X
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-1

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\82N111820W.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.778	310	323	337	rBV	106900	221752	4.68%	0.788%
2	3.129	415	432	452	rVB5	114358	314070	6.63%	1.117%
3	3.286	469	481	495	rVB	75471	155180	3.28%	0.552%
4	3.438	514	528	540	rBV4	39225	95784	2.02%	0.341%
5	3.524	541	555	582	rVB	259346	632279	13.35%	2.248%
6	4.383	803	822	838	rBV	67501	197332	4.17%	0.702%
7	4.585	870	885	916	rVB6	52774	213676	4.51%	0.760%
8	4.743	919	934	957	rVB4	19204	70654	1.49%	0.251%
9	5.875	1271	1286	1308	rVB3	51237	162449	3.43%	0.578%
10	6.023	1311	1332	1348	rBV6	27893	88889	1.88%	0.316%
11	6.325	1407	1426	1450	rVB4	38019	114977	2.43%	0.409%
12	6.579	1483	1505	1522	rBV4	55698	176934	3.74%	0.629%
13	6.778	1550	1567	1590	rVB	17089	52766	1.11%	0.188%
14	7.347	1725	1744	1761	rVB3	115196	293754	6.20%	1.044%
15	7.550	1786	1807	1817	rBV2	168976	429047	9.06%	1.525%
16	7.624	1818	1830	1849	rVB2	332928	807290	17.05%	2.870%
17	7.994	1926	1945	1967	rBV3	426991	1131762	23.90%	4.023%
18	8.138	1967	1990	2001	rBV5	48439	137295	2.90%	0.488%
19	8.550	2103	2118	2142	rBV	553004	1196710	25.27%	4.254%
20	9.582	2426	2439	2446	rBV	32790	65356	1.38%	0.232%
21	9.630	2448	2454	2469	rVB8	24702	53408	1.13%	0.190%
22	10.058	2574	2587	2596	rBV	847717	1539729	32.52%	5.474%
23	10.122	2596	2607	2641	rVB	2612732	4735150	100.00%	16.834%
24	11.379	2982	2998	3014	rBV	874853	1518642	32.07%	5.399%
25	11.482	3015	3030	3048	rVV	1663664	2707599	57.18%	9.626%
26	11.592	3049	3064	3086	rVB	1381106	2632318	55.59%	9.358%
27	11.920	3149	3166	3181	rBV	653452	1076414	22.73%	3.827%
28	12.222	3246	3260	3273	rBV	270341	431932	9.12%	1.536%
29	12.373	3295	3307	3330	rVB	723964	1220311	25.77%	4.338%
30	12.563	3353	3366	3377	rBV	299444	484384	10.23%	1.722%
31	12.627	3377	3386	3392	rVV	183873	305729	6.46%	1.087%
32	12.659	3392	3396	3403	rVV	137044	204287	4.31%	0.726%
33	12.704	3403	3410	3424	rVB	182811	300731	6.35%	1.069%
34	12.865	3446	3460	3474	rBV	169406	281696	5.95%	1.001%

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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 >
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35	13.013	3493	3506	3519	rBV	591898	935258	19.75%	3.325%
36	13.315	3587	3600	3623	rBV2	914244	1768482	37.35%	6.287%
37	13.514	3643	3662	3672	rBV2	373001	665253	14.05%	2.365%
38	13.707	3710	3722	3734	rVB2	30421	70142	1.48%	0.249%
39	13.862	3760	3770	3780	rBV	68904	105956	2.24%	0.377%
40	13.968	3794	3803	3813	rVB2	37055	58766	1.24%	0.209%
41	14.180	3858	3869	3876	rBV	41091	67259	1.42%	0.239%
42	14.447	3942	3952	3962	rBV2	33002	53394	1.13%	0.190%
43	14.566	3975	3989	4001	rBV3	69861	137477	2.90%	0.489%
44	15.103	4143	4156	4169	rBV	122198	216974	4.58%	0.771%

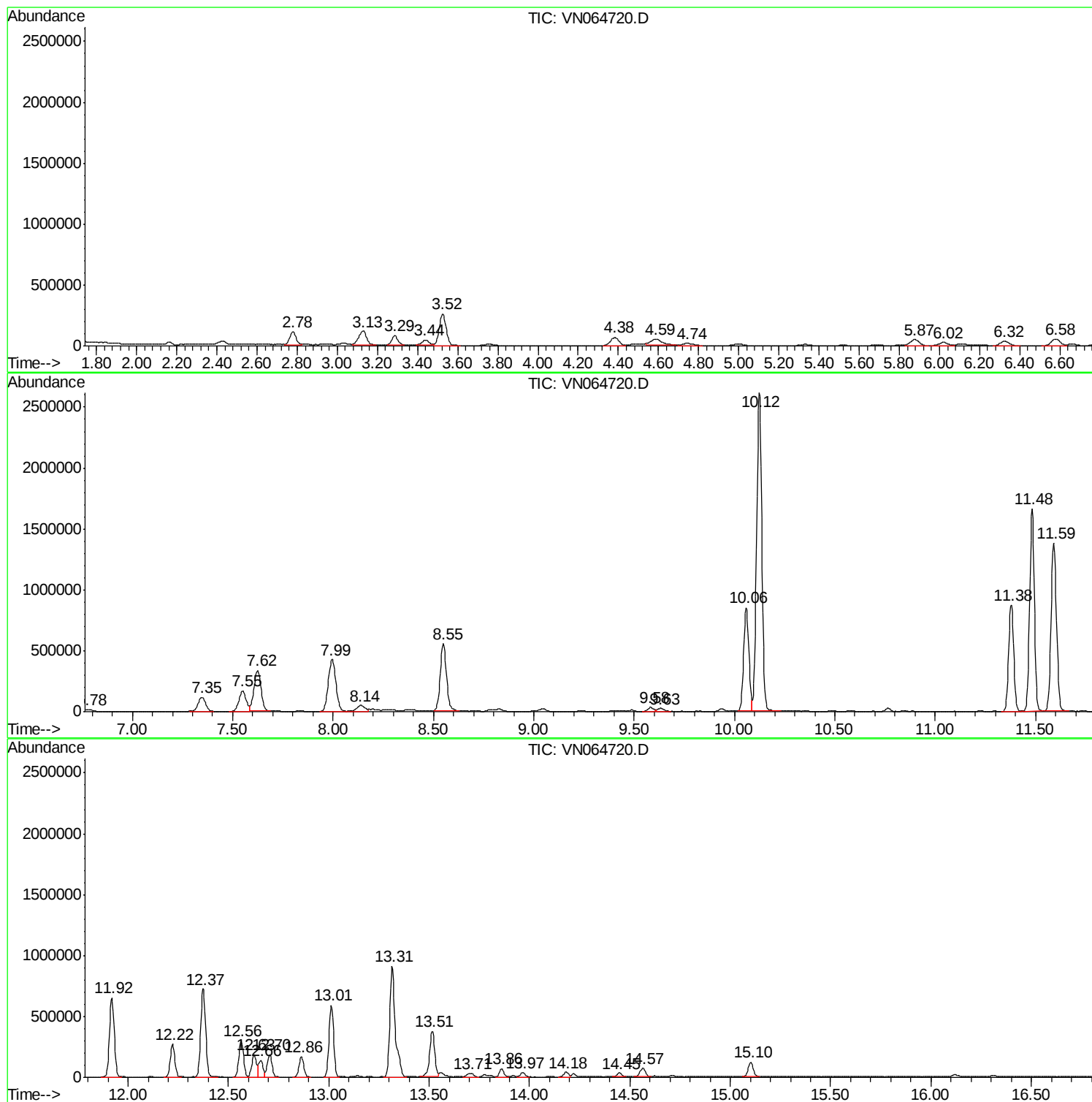
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N111820W.M
Quant Title : SW846 8260

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



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Instrument :
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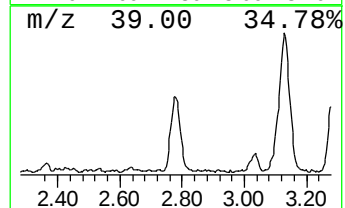
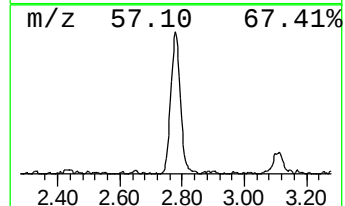
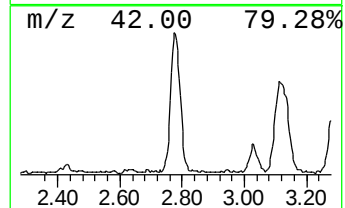
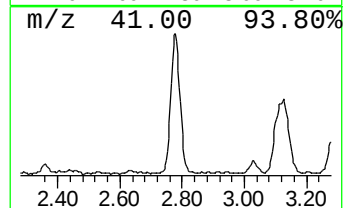
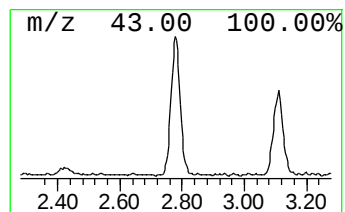
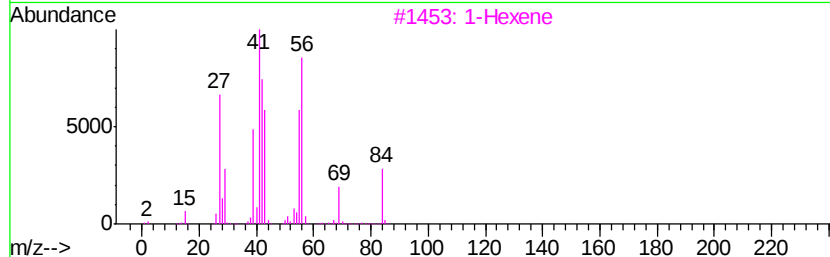
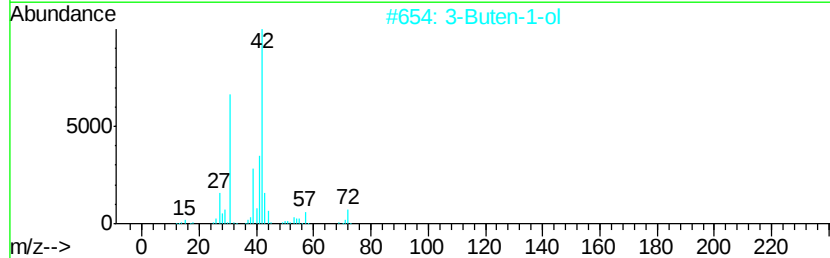
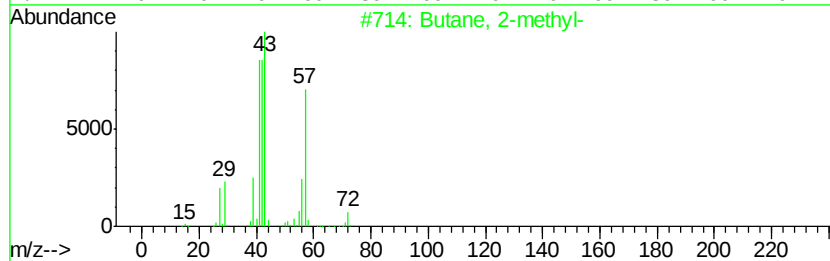
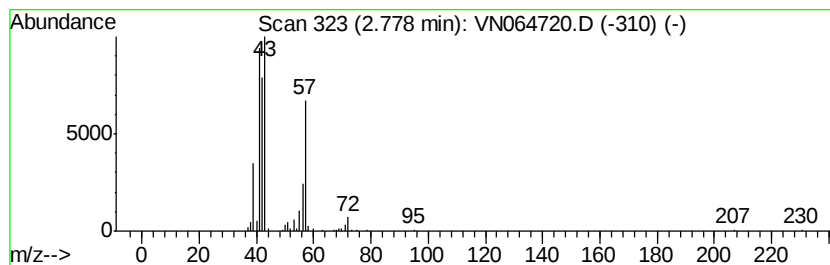
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N111820W.M
Quant Title : SW846 8260

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

Peak Number 1 Butane, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.78	13.73 ug/l	221752	Pentafluorobenzene	7.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methyl-	72	C5H12	000078-78-4	91
2		3-Buten-1-ol	72	C4H8O	000627-27-0	38
3		1-Hexene	84	C6H12	000592-41-6	36
4		Pentane	72	C5H12	000109-66-0	33
5		2(3H)-Furanone, dihydro-4-methyl-	100	C5H8O2	001679-49-8	28



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Instrument :
MSVOA_N
ClientSampled :
MW-1

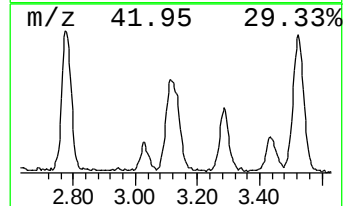
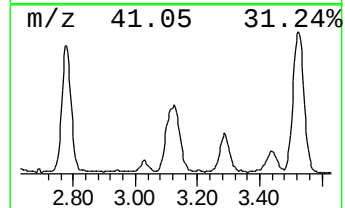
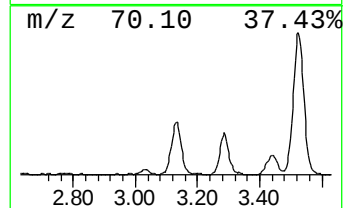
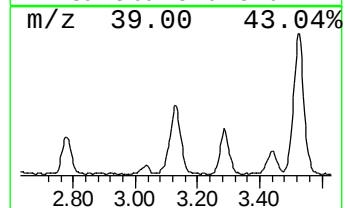
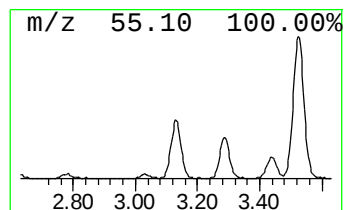
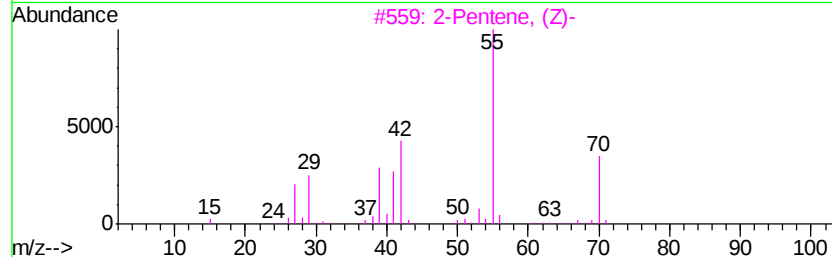
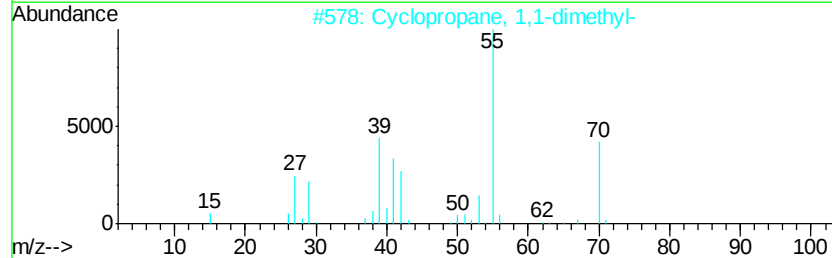
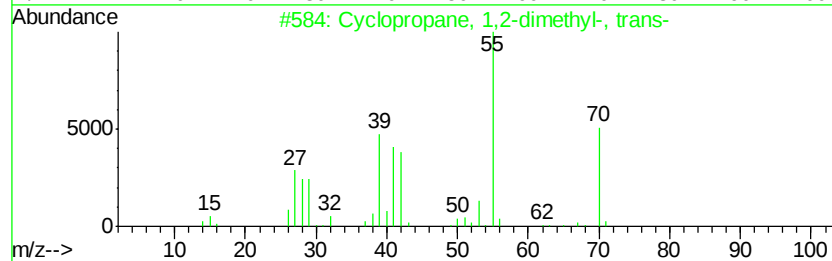
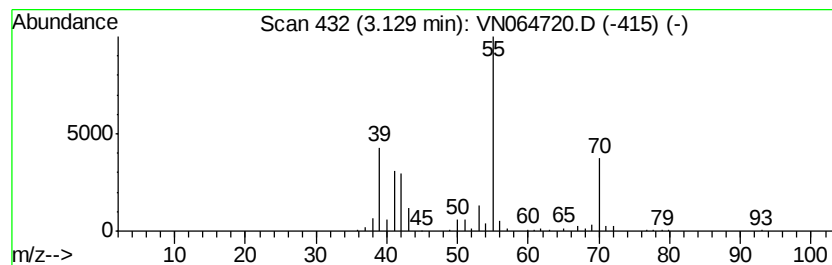
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N111820W.M
Quant Title : SW846 8260

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

Peak Number 2 Cyclopropane, 1,2-dimethyl-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.13	19.45 ug/l	314070	Pentafluorobenzene	7.62

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopropane, 1,2-dimethyl-, trans-	70	C5H10	002402-06-4	87
2		Cyclopropane, 1,1-dimethyl-	70	C5H10	001630-94-0	83
3		2-Pentene, (Z)-	70	C5H10	000627-20-3	81
4		2-Butene, 2-methyl-	70	C5H10	000513-35-9	64
5		2-Pentene, (E)-	70	C5H10	000646-04-8	60



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ALS Vial : 23 Sample Multiplier: 1

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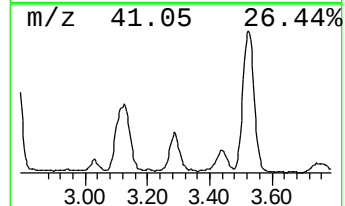
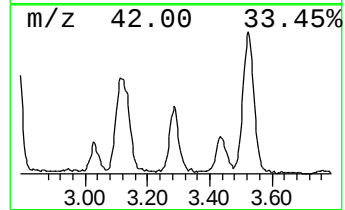
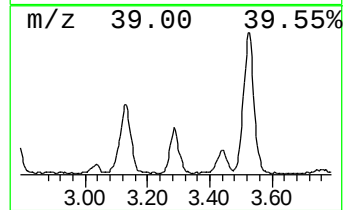
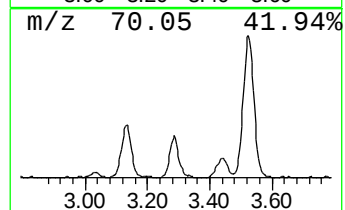
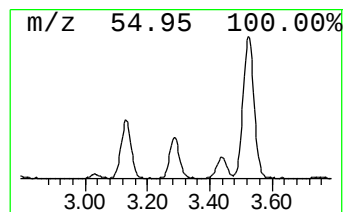
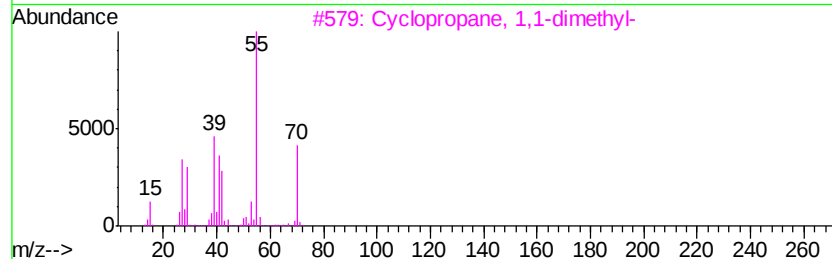
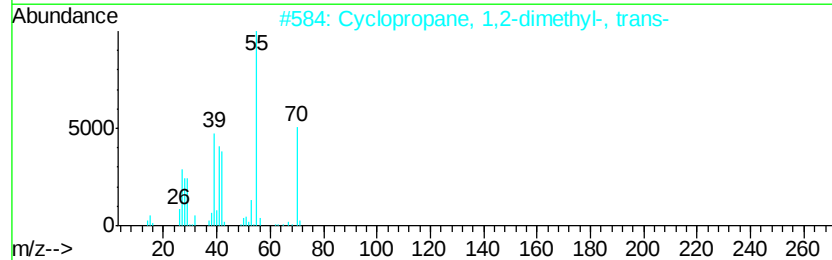
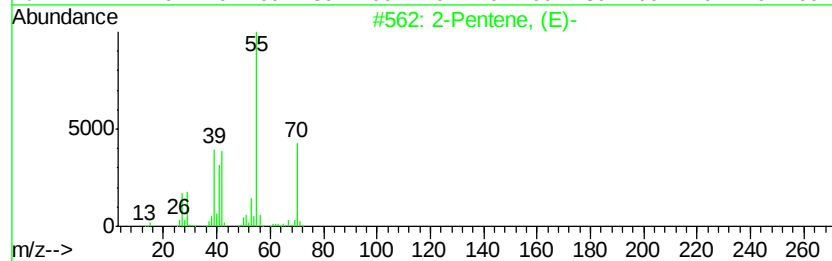
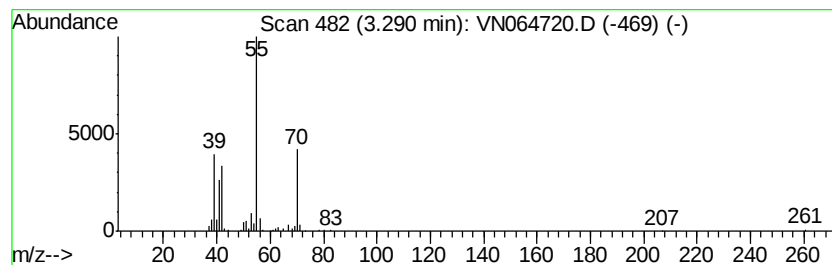
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentene, (E)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.29	9.61 ug/l	155180	Pentafluorobenzene	7.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentene, (E)-	70	C5H10	000646-04-8	90
2		Cyclopropane, 1,2-dimethyl-, trans-	70	C5H10	002402-06-4	87
3		Cyclopropane, 1,1-dimethyl-	70	C5H10	001630-94-0	86
4		2-Butene, 2-methyl-	70	C5H10	000513-35-9	78
5		2-Pentene, (Z)-	70	C5H10	000627-20-3	74



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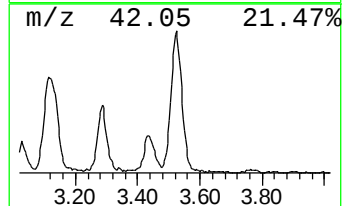
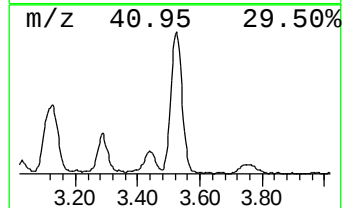
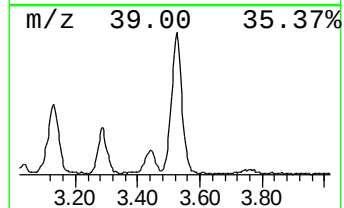
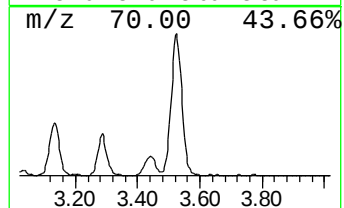
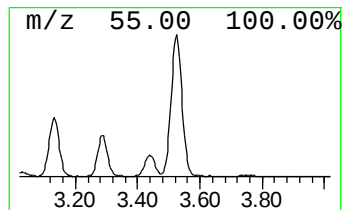
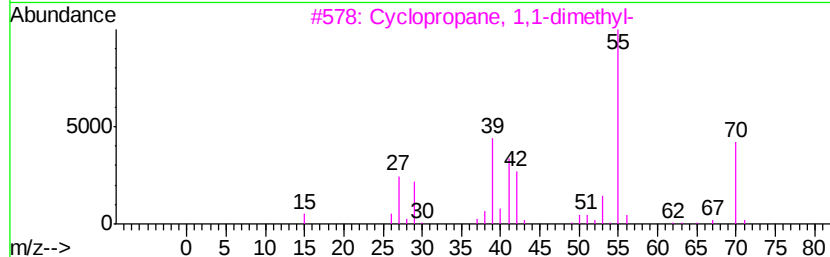
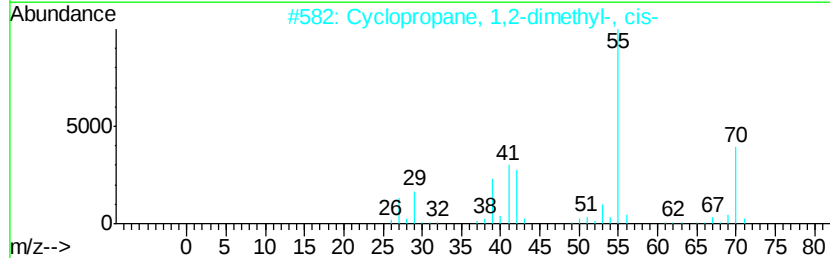
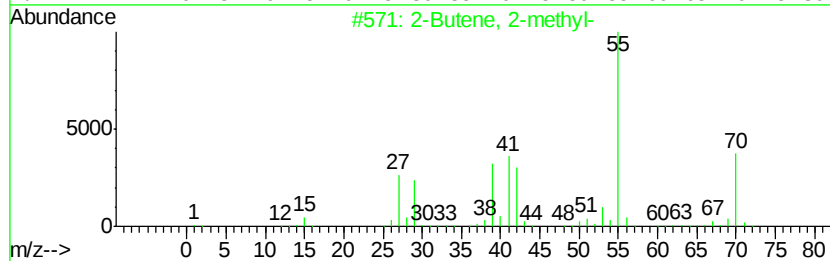
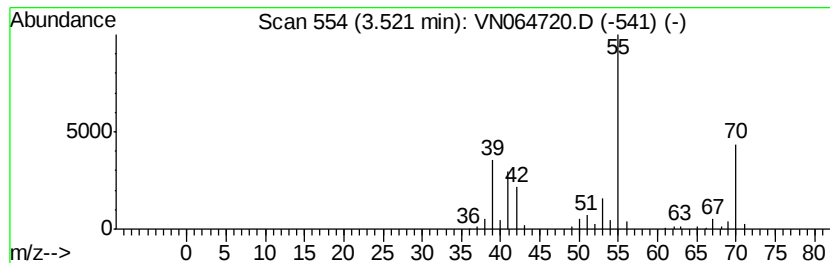
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 2-Butene, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.52	39.16 ug/l	632279	Pentafluorobenzene	7.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Butene, 2-methyl-	70	C5H10	000513-35-9	90
2		Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	86
3		Cyclopropane, 1,1-dimethyl-	70	C5H10	001630-94-0	86
4		Cyclopropane, 1,2-dimethyl-, trans-	70	C5H10	002402-06-4	80
5		1-Butene, 3-methyl-	70	C5H10	000563-45-1	80



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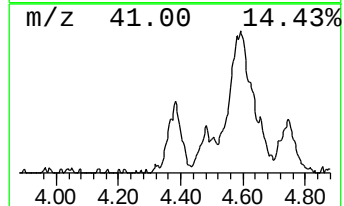
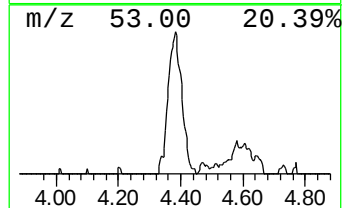
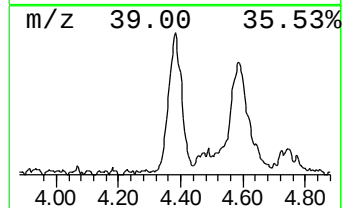
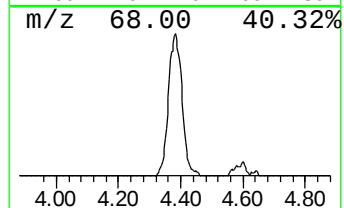
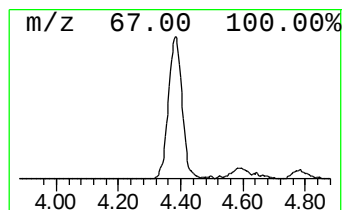
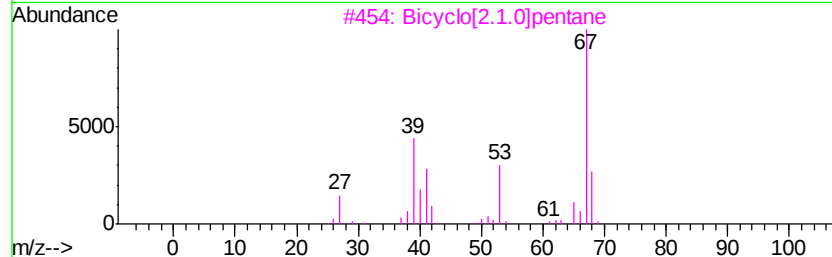
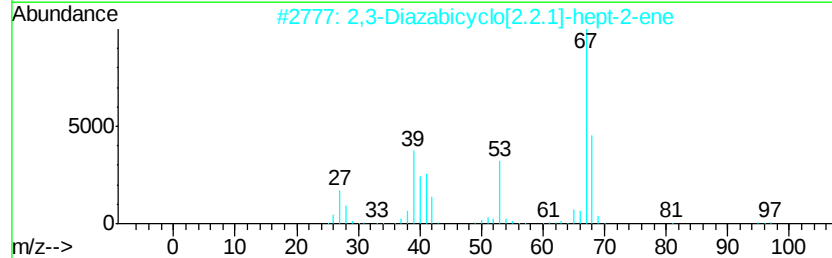
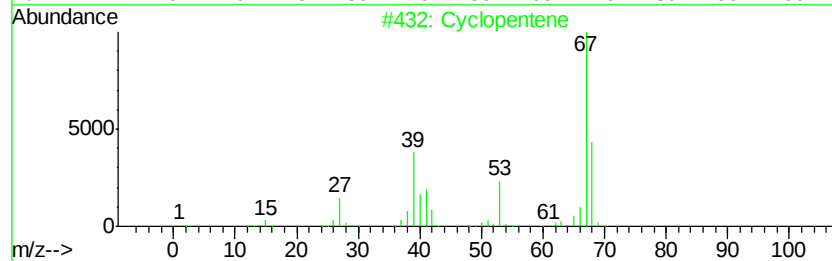
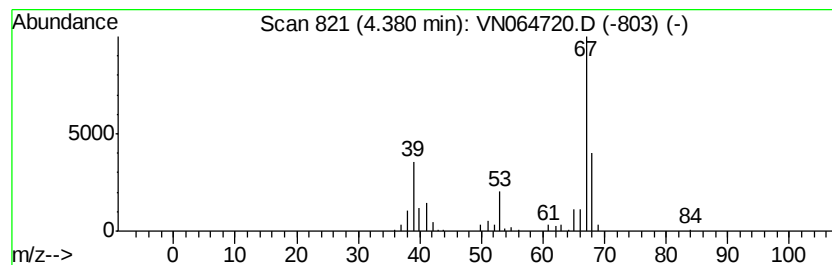
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Quant Title : SW846 8260

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

Peak Number 5 Cyclopentene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.38	12.22 ug/l	197332	Pentafluorobenzene	7.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentene	68	C5H8	000142-29-0	91
2		2,3-Diazabicyclo[2.2.1]-hept-2-ene	96	C5H8N2	002721-32-6	72
3		Bicyclo[2.1.0]pentane	68	C5H8	000185-94-4	59
4		1,3-Pentadiene, (Z)-	68	C5H8	001574-41-0	50
5		1,3-Butadiene, 2-methyl-	68	C5H8	000078-79-5	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN111820\
Data File : VN064720.D
Acq On : 18 Nov 2020 20:36
Operator : JC/MD
Sample : L4783-20 40X
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1

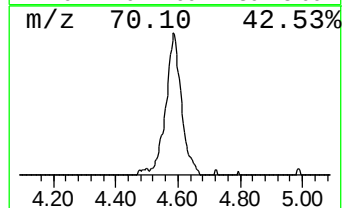
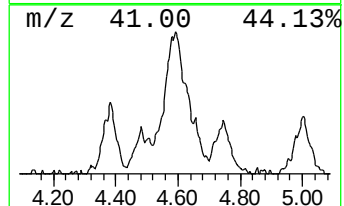
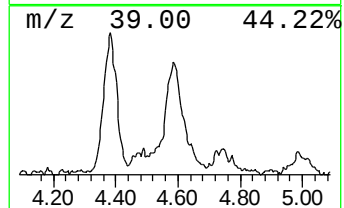
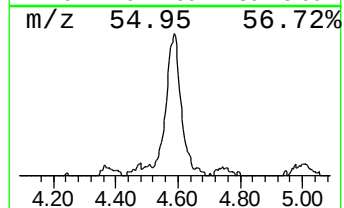
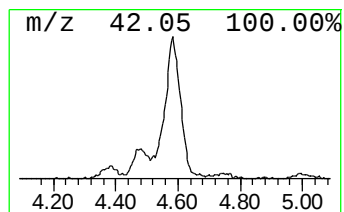
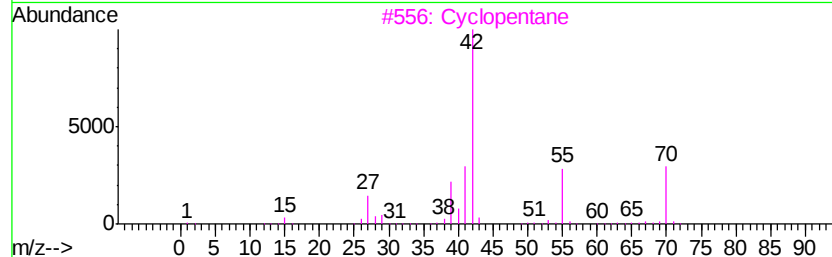
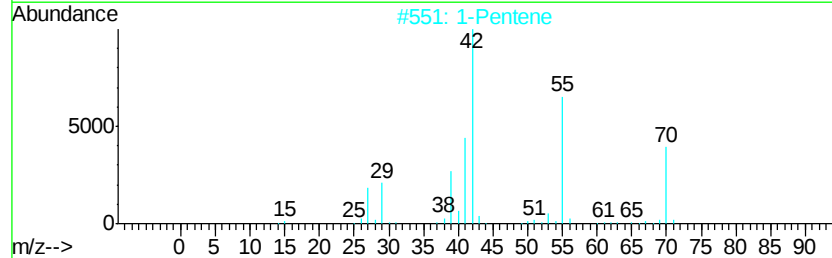
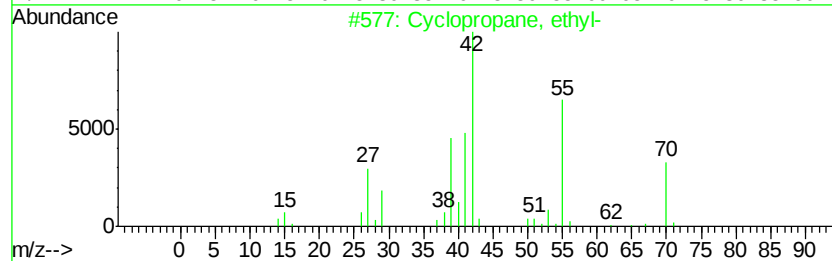
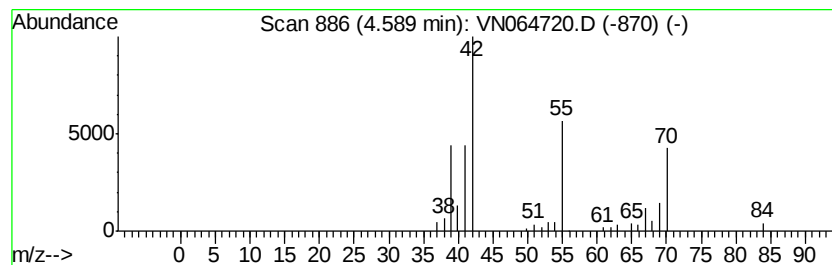
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Quant Title : SW846 8260

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

Peak Number 6 Cyclopropane, ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.59	13.23 ug/l	213676	Pentafluorobenzene	7.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopropane, ethyl-	70	C5H10	001191-96-4	64
2		1-Pentene	70	C5H10	000109-67-1	52
3		Cyclopentane	70	C5H10	000287-92-3	50
4		Cyclobutane, methyl-	70	C5H10	000598-61-8	47
5		2H-Tetrazole, 2,5-dimethyl-	98	C3H6N4	004135-93-7	33



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN111820\
Data File : VN064720.D
Acq On : 18 Nov 2020 20:36
Operator : JC/MD
Sample : L4783-20 40X
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampled :
MW-1

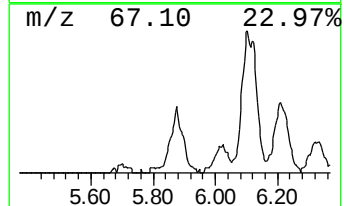
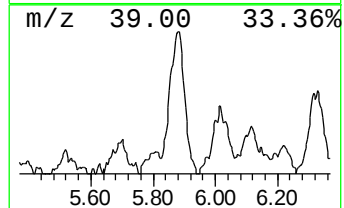
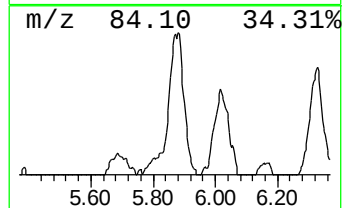
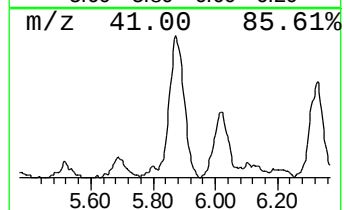
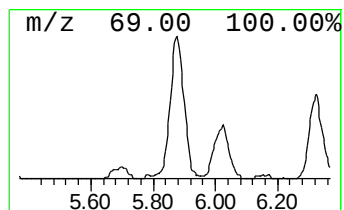
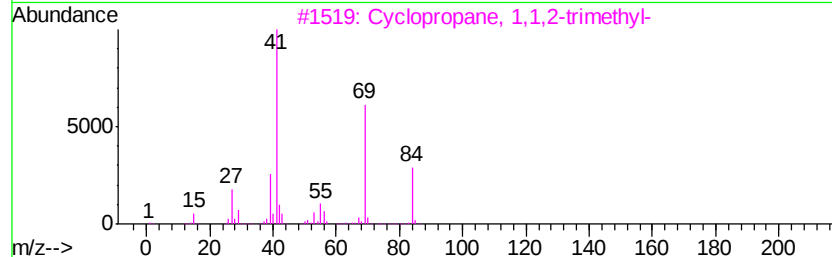
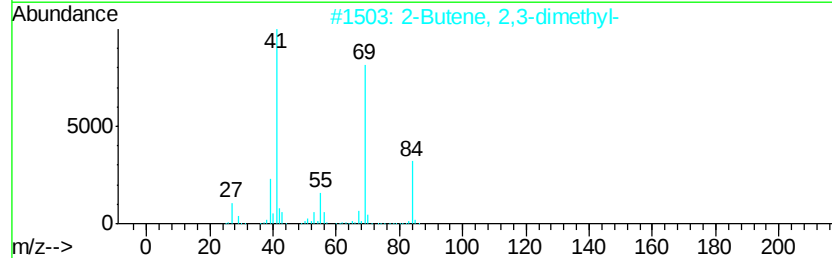
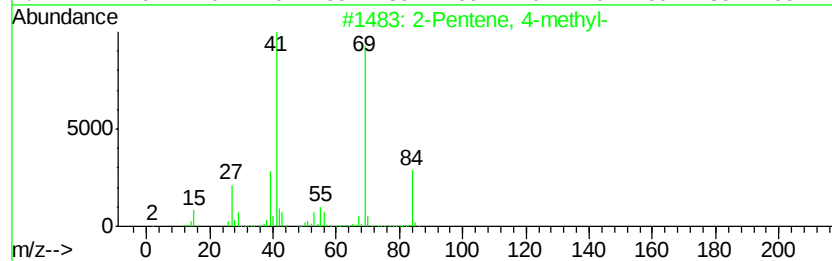
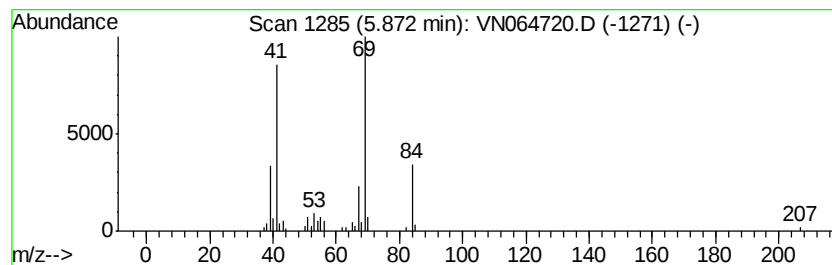
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Quant Title : SW846 8260

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

Peak Number 7 2-Pentene, 4-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.87	10.06 ug/l	162449	Pentafluorobenzene	7.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentene, 4-methyl-	84	C6H12	004461-48-7	80
2		2-Butene, 2,3-dimethyl-	84	C6H12	000563-79-1	80
3		Cyclopropane, 1,1,2-trimethyl-	84	C6H12	004127-45-1	72
4		2-Pentene, 2-methyl-	84	C6H12	000625-27-4	72
5		2-Pentene, 4-methyl-, (Z)-	84	C6H12	000691-38-3	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN111820\
Data File : VN064720.D
Acq On : 18 Nov 2020 20:36
Operator : JC/MD
Sample : L4783-20 40X
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1

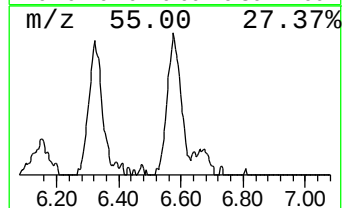
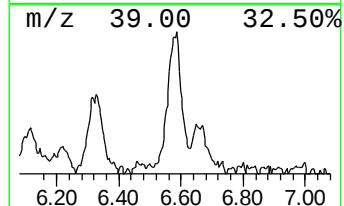
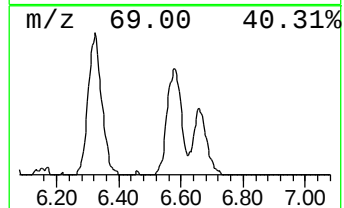
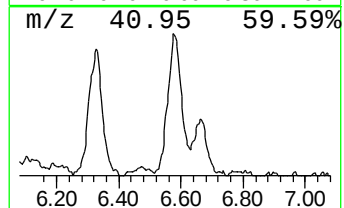
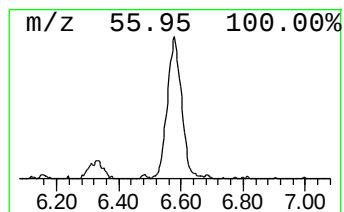
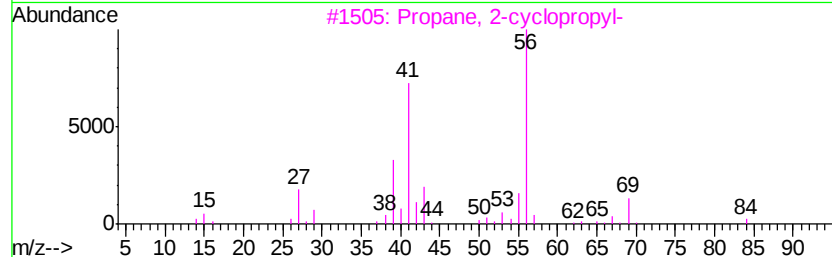
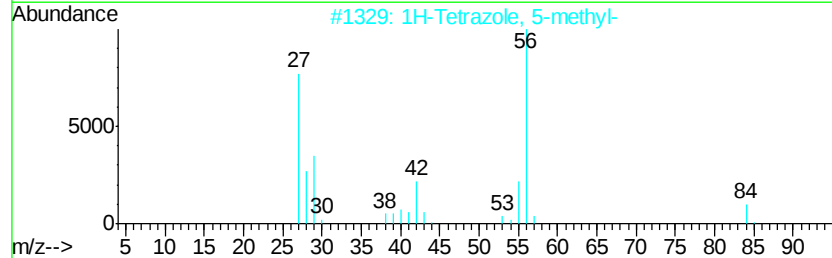
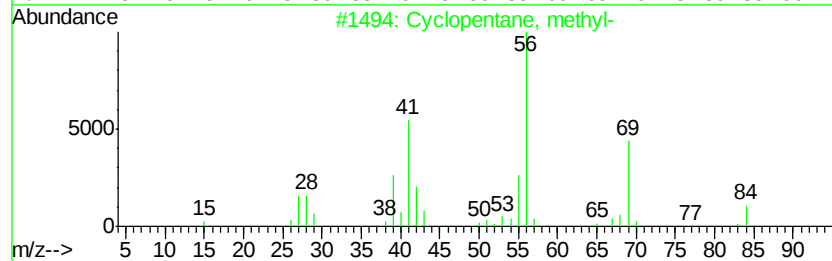
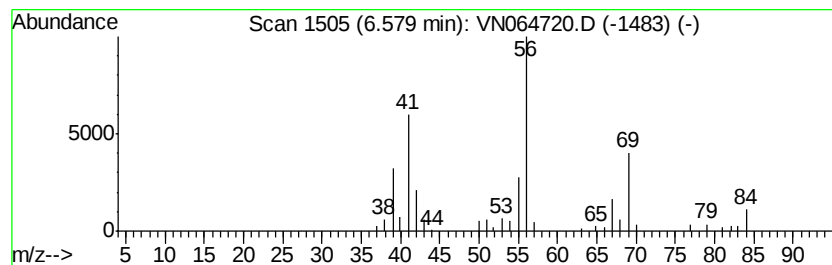
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Quant Title : SW846 8260

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

Peak Number 8 Cyclopentane, methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	10.96 ug/l	176934	Pentafluorobenzene	7.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	87
2		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64
3		Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	53
4		Cyclobutane, ethyl-	84	C6H12	004806-61-5	50
5		Cyclobutane	56	C4H8	000287-23-0	47



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN111820\
Data File : VN064720.D
Acq On : 18 Nov 2020 20:36
Operator : JC/MD
Sample : L4783-20 40X
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1

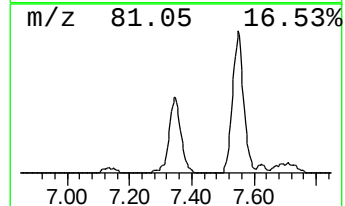
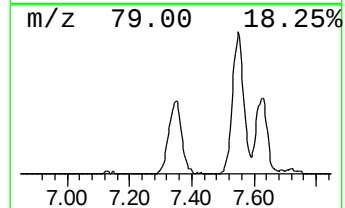
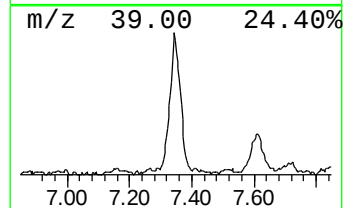
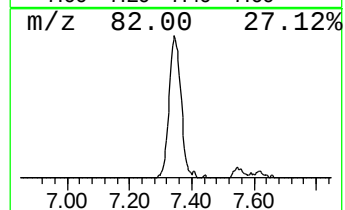
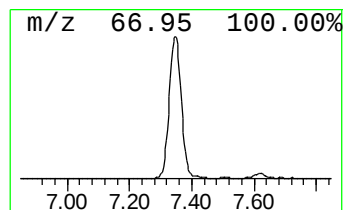
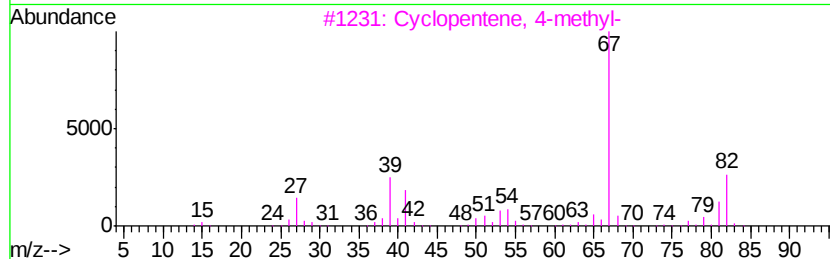
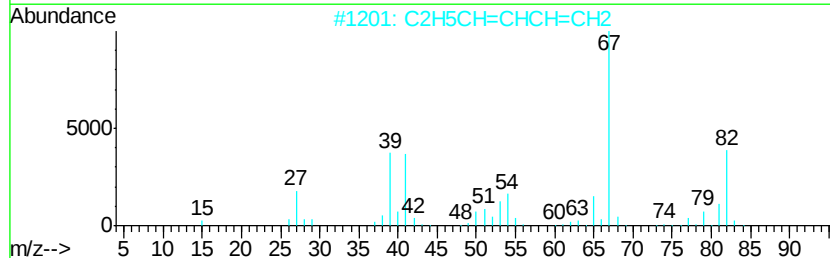
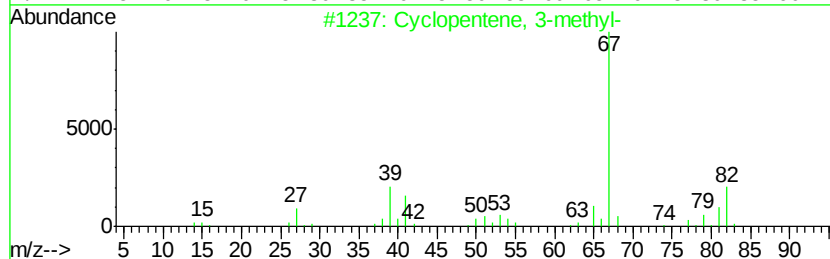
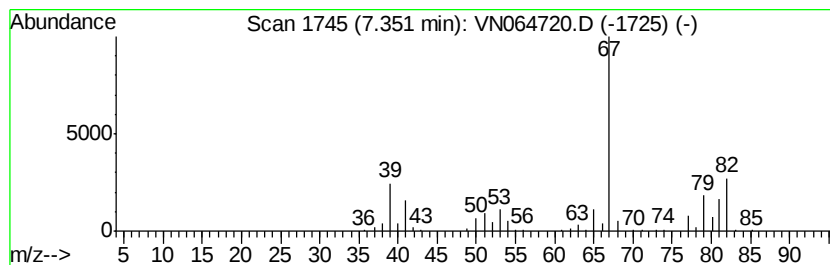
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Quant Title : SW846 8260

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

Peak Number 9 Cyclopentene, 3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.35	18.19 ug/l	293754	Pentafluorobenzene	7.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentene, 3-methyl-	82	C6H10	001120-62-3	90
2		C2H5CH=CHCH=CH2	82	C6H10	000592-48-3	90
3		Cyclopentene, 4-methyl-	82	C6H10	001759-81-5	90
4		Cyclopentene, 1-methyl-	82	C6H10	000693-89-0	87
5		4-Methyl-1,3-pentadiene	82	C6H10	000926-56-7	80



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN111820\
Data File : VN064720.D
Acq On : 18 Nov 2020 20:36
Operator : JC/MD
Sample : L4783-20 40X
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1

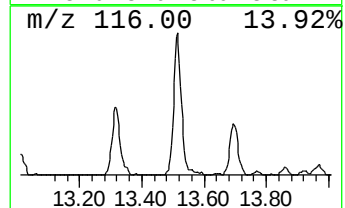
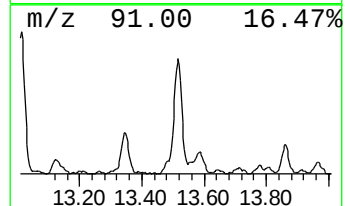
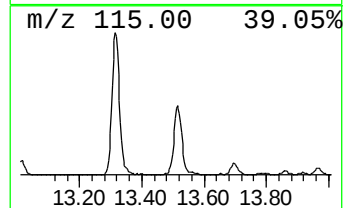
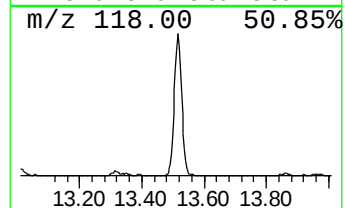
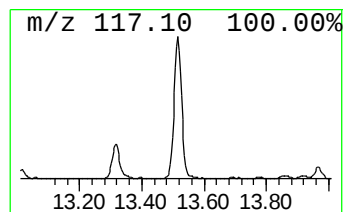
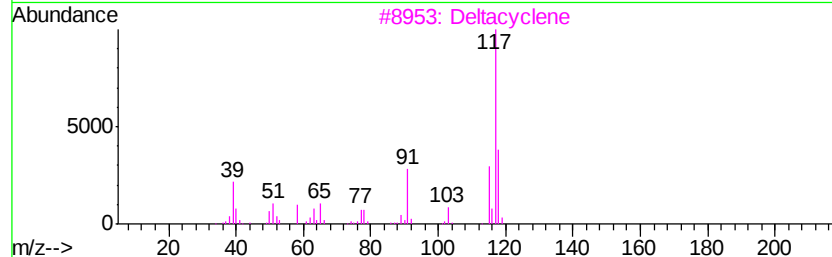
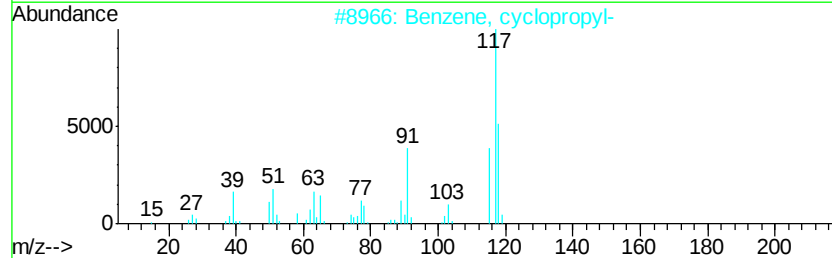
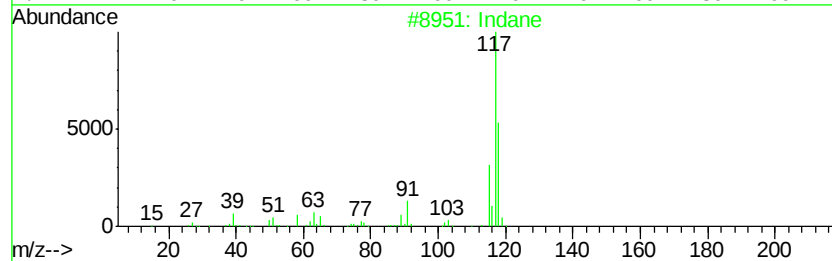
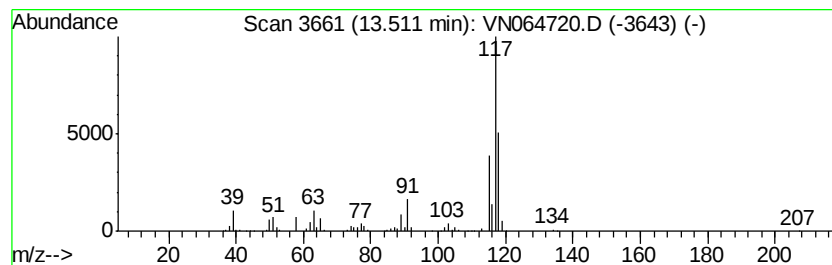
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N111820W.M
Quant Title : SW846 8260

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

Peak Number 10 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.51	18.81 ug/l	665253	1,4-Dichlorobenzene-d4	13.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	95
2	Benzene, cyclopropyl-		118	C9H10	000873-49-4	81
3	Deltacyclene		118	C9H10	007785-10-6	76
4	Benzene, 2-propenyl-		118	C9H10	000300-57-2	74
5	Benzene, 1-ethenyl-2-methyl-		118	C9H10	000611-15-4	68



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN111820\
Data File : VN064720.D
Acq On : 18 Nov 2020 20:36
Operator : JC/MD
Sample : L4783-20 40X
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N111820W.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Cyclopropane, 1,2...	3.13	19.4	ug/l	314070	1	7.62	807290	50.0
2-Pentene, (E)-	3.29	9.6	ug/l	155180	1	7.62	807290	50.0
2-Butene, 2-methyl-	3.52	39.2	ug/l	632279	1	7.62	807290	50.0
Cyclopentene	4.38	12.2	ug/l	197332	1	7.62	807290	50.0
Cyclopropane, ethyl-	4.59	13.2	ug/l	213676	1	7.62	807290	50.0
2-Pentene, 4-methyl-	5.87	10.1	ug/l	162449	1	7.62	807290	50.0
Cyclopentane, met...	6.58	11.0	ug/l	176934	1	7.62	807290	50.0
Cyclopentene, 3-m...	7.35	18.2	ug/l	293754	1	7.62	807290	50.0
Indane	13.51	18.8	ug/l	665253	4	13.32	1768480	50.0