

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_D\DATA\VD072820\
 Data File : VD066077.D
 Acq On : 28 Jul 2020 13:49
 Operator : VA/SY
 Sample : L3181-06
 Misc : 8.39G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 CL-01-072720

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_D\METHOD\82D072120S.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	4.961	507	517	532	rBV2	25123	81550	1.23%	0.363%
2	7.914	1009	1019	1024	rBV	181078	422242	6.38%	1.878%
3	7.979	1024	1030	1039	rVB	335407	746740	11.28%	3.321%
4	8.332	1077	1090	1104	rVB3	140226	340153	5.14%	1.513%
5	8.867	1173	1181	1193	rBV	458429	934891	14.12%	4.158%
6	10.344	1423	1432	1439	rBV	722764	1324621	20.01%	5.892%
7	11.649	1645	1654	1664	rBV	751021	1288581	19.47%	5.731%
8	11.855	1682	1689	1700	rBV	44857	101605	1.53%	0.452%
9	12.185	1735	1745	1754	rBV4	41741	97808	1.48%	0.435%
10	12.638	1815	1822	1831	rVB	660805	1098584	16.60%	4.886%
11	13.279	1925	1931	1935	rBV	85649	130720	1.97%	0.581%
12	13.585	1976	1983	1992	rVV	836494	1428868	21.58%	6.355%
13	13.785	2012	2017	2028	rVB2	80270	155448	2.35%	0.691%
14	13.908	2032	2038	2042	rBV6	51206	93424	1.41%	0.416%
15	13.967	2042	2048	2054	rBV	227727	386385	5.84%	1.719%
16	14.067	2059	2065	2070	rBV5	51463	90938	1.37%	0.404%
17	14.126	2070	2075	2078	rVB4	50428	73635	1.11%	0.328%
18	14.167	2078	2082	2088	rVB4	60979	101668	1.54%	0.452%
19	14.237	2088	2094	2098	rBV3	56398	116085	1.75%	0.516%
20	14.296	2100	2104	2110	rVB6	47148	81813	1.24%	0.364%
21	14.385	2110	2119	2121	rBV4	78803	173961	2.63%	0.774%
22	14.420	2121	2125	2134	rVB7	107726	222651	3.36%	0.990%
23	14.532	2140	2144	2147	rBV4	56287	83648	1.26%	0.372%
24	14.579	2150	2152	2157	rVB4	72078	99476	1.50%	0.442%
25	14.726	2170	2177	2181	rBV4	43262	107925	1.63%	0.480%
26	14.837	2191	2196	2199	rBV4	194702	363846	5.50%	1.618%
27	14.879	2199	2203	2212	rVB3	381768	771433	11.65%	3.431%
28	14.996	2219	2223	2229	rVB6	103529	184040	2.78%	0.819%
29	15.414	2282	2294	2305	rBV	2878159	6619812	100.00%	29.444%
30	15.514	2306	2311	2315	rBV3	119807	226185	3.42%	1.006%
31	15.714	2337	2345	2351	rBV8	145408	439901	6.65%	1.957%
32	15.790	2354	2358	2363	rVB4	198155	361027	5.45%	1.606%
33	16.355	2449	2454	2462	rVB	377797	688569	10.40%	3.063%
34	16.508	2474	2480	2492	rVB	655621	1495827	22.60%	6.653%

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Instrument :
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ClientSampleId :
CL-01-072720

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

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35	16.714	2507	2515	2527	rVB	517829	1548510	23.39%	6.888%
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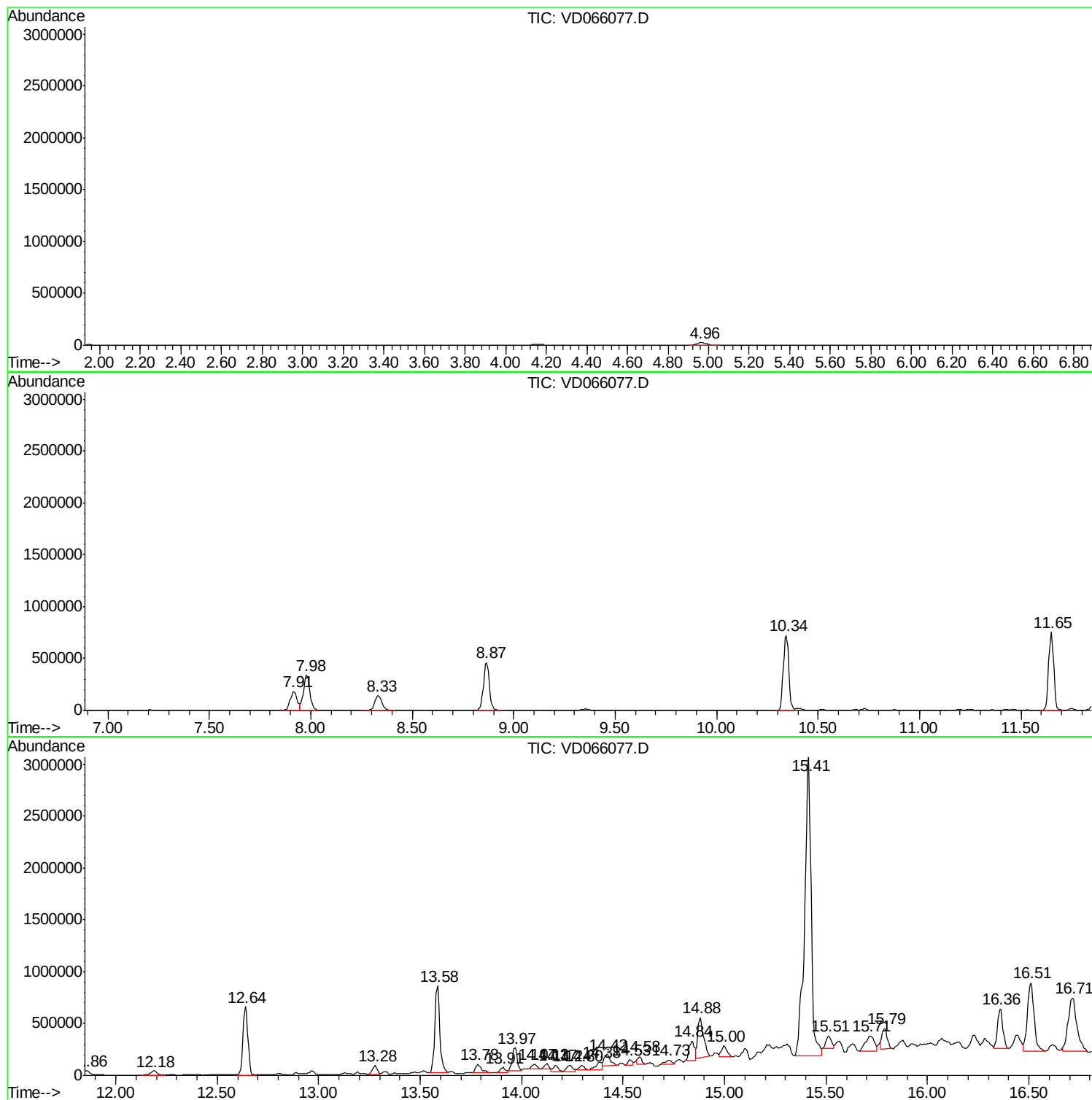
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Data Path : Z:\V0ASRV\HPCHEM1\MSV0A_D\DATA\VD072820\
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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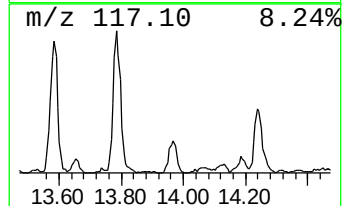
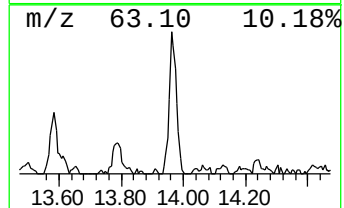
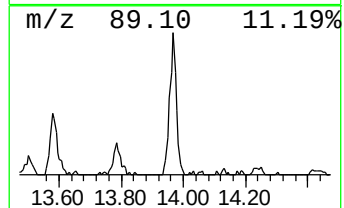
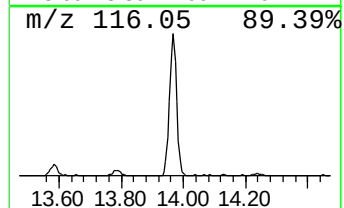
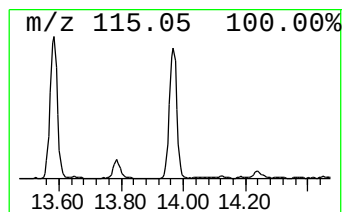
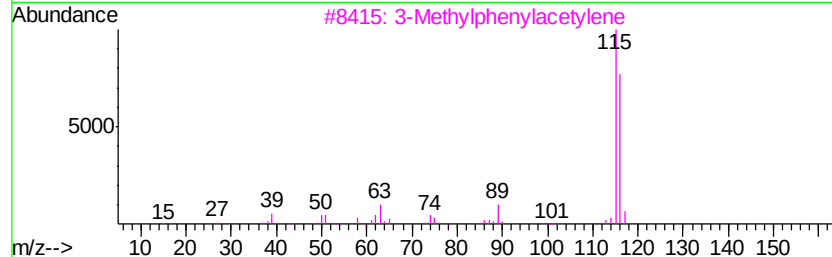
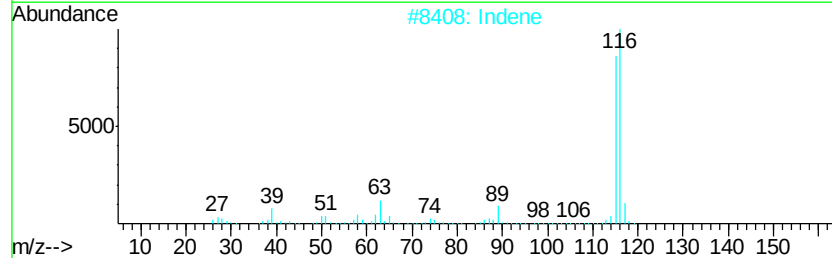
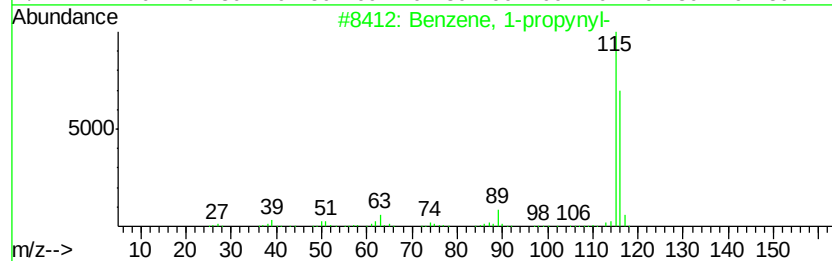
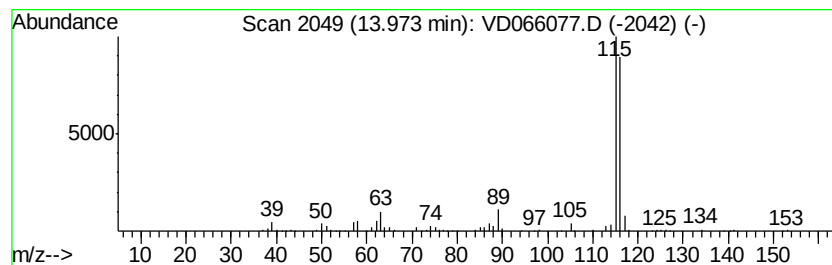
Instrument :
MSVOA_D
ClientSampleId :
CL-01-072720

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Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-propynyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	13.52 ug/l	386385	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-propynyl-	116 C9H8	000673-32-5 95
2		Indene	116 C9H8	000095-13-6 95
3		3-Methylphenylacetylene	116 C9H8	000766-82-5 94
4		Benzene, 1,2-propadienyl-	116 C9H8	002327-99-3 91
5		Benzene, 1-ethynyl-4-methyl-	116 C9H8	000766-97-2 87



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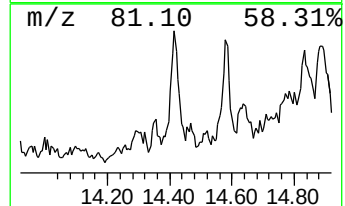
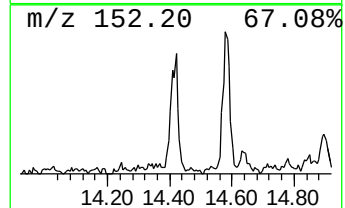
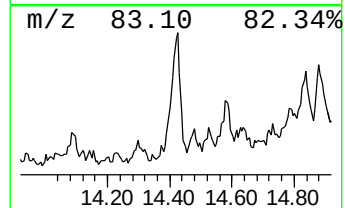
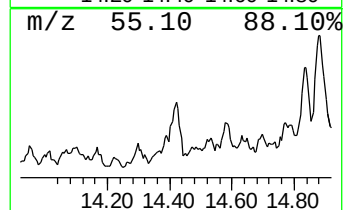
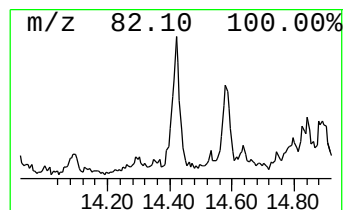
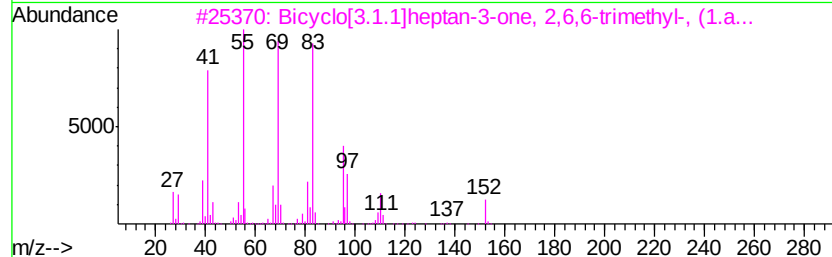
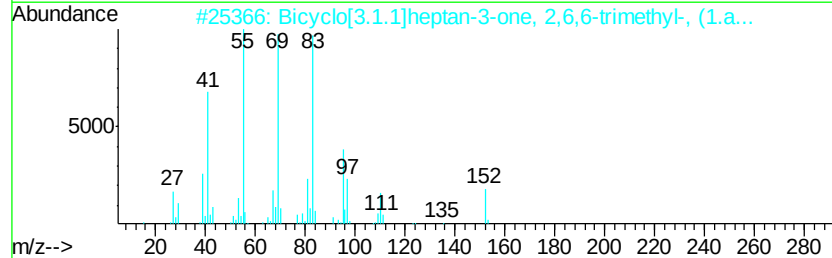
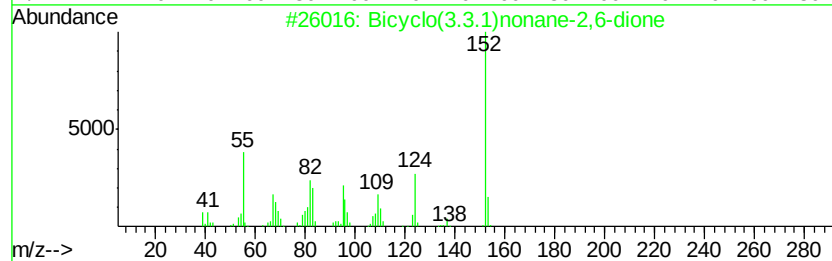
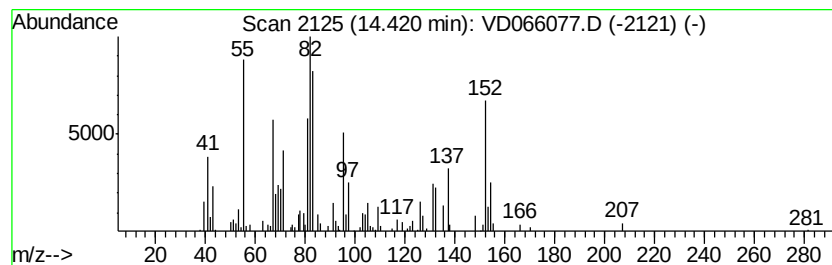
Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D072120S.M
Quant Title : SW846 8260

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

Peak Number 2 Bicyclo(3.3.1)nonane-2,6-dione Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.42	7.79 ug/l	222651	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Bicyclo(3.3.1)nonane-2,6-dione	152 C9H12O2	016473-11-3 52
2		Bicyclo[3.1.1]heptan-3-one, 2,6,...	152 C10H16O	015358-88-0 38
3		Bicyclo[3.1.1]heptan-3-one, 2,6,...	152 C10H16O	000547-60-4 38
4		1H-Pyrrole, 2,3-dihydro-1-methyl-	83 C5H9N	033838-11-8 27
5		Cyclohexane, (1-methylethyl)-	126 C9H18	000696-29-7 27



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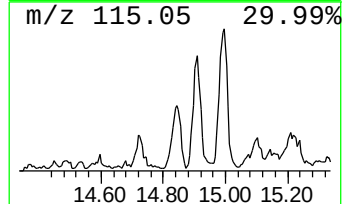
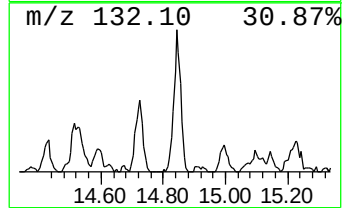
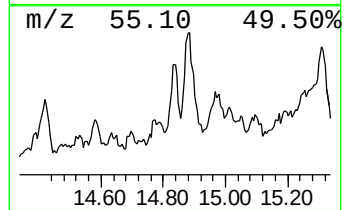
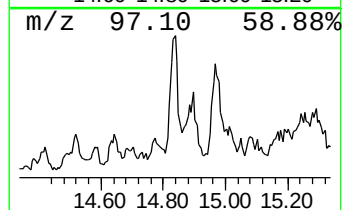
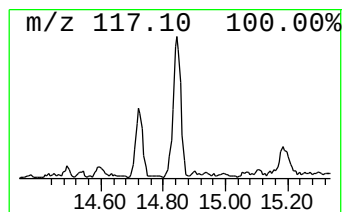
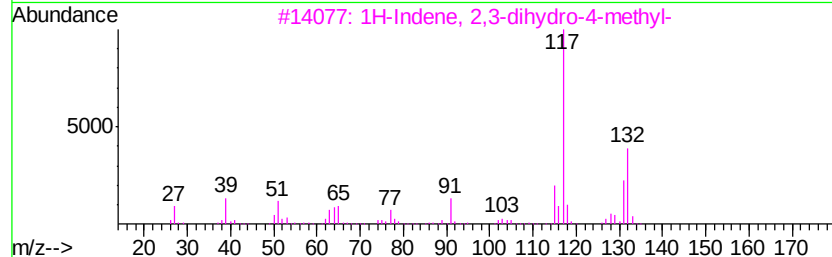
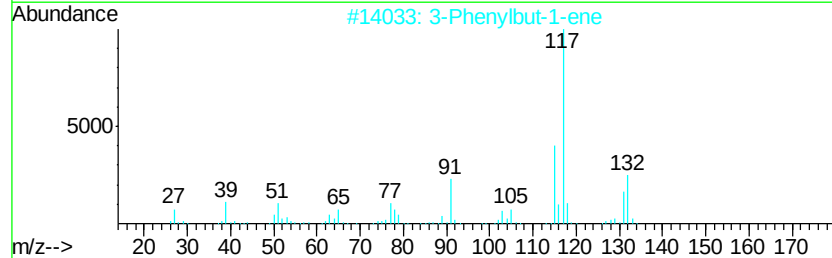
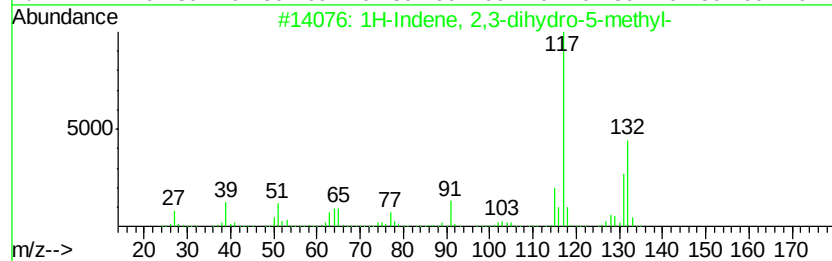
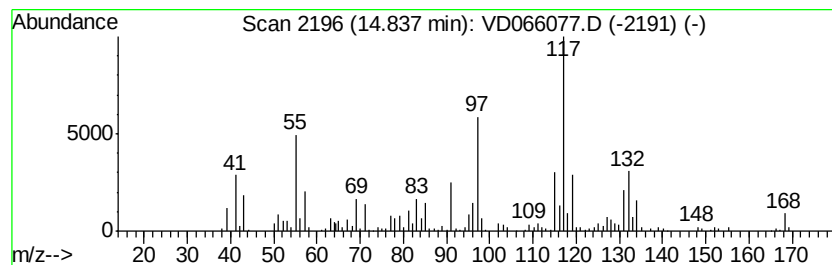
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.84	12.73 ug/l	363846	1,4-Dichlorobenzene-d4	13.58	
Hit# of	5	Tentative ID	MW MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	64
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	64
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	52
4	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	50
5	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	47



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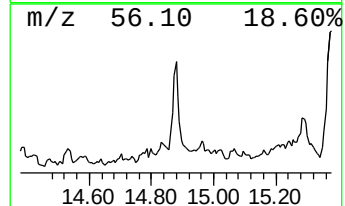
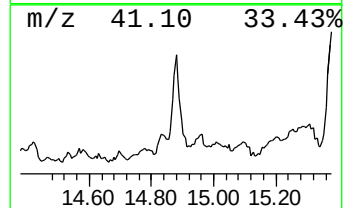
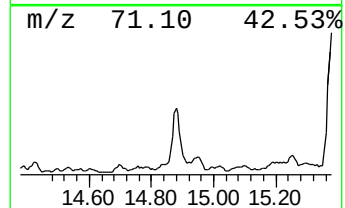
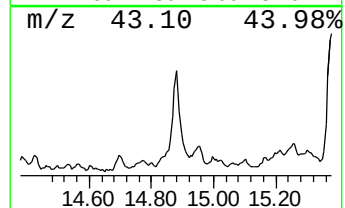
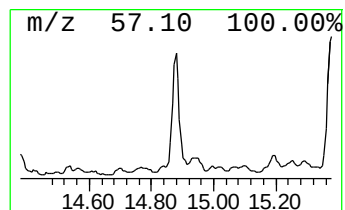
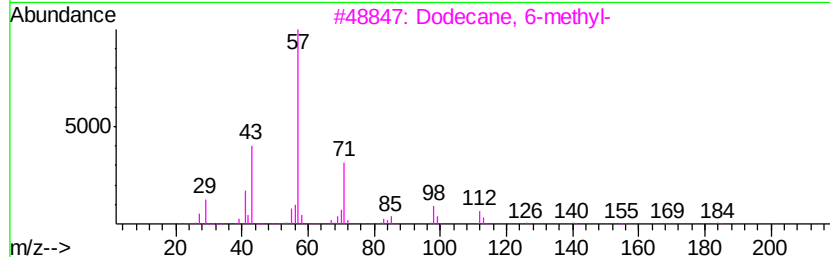
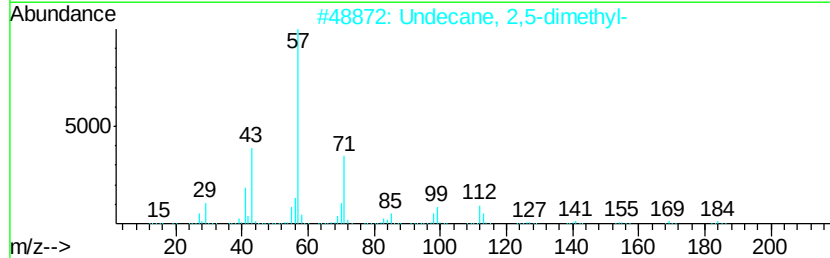
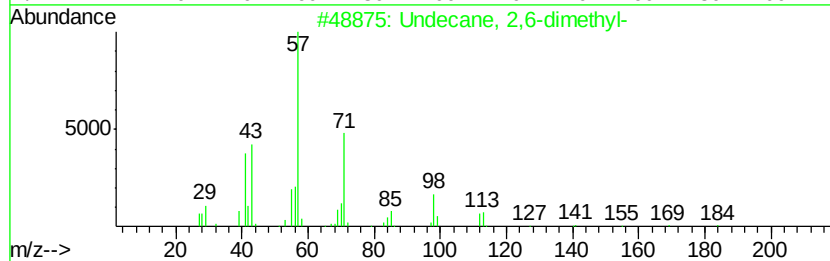
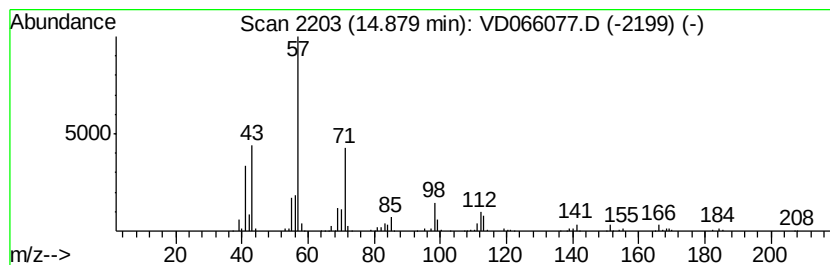
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 Undecane, 2,6-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.88	26.99 ug/l	771433	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Undecane, 2,6-dimethyl-	184 C13H28	017301-23-4	93
2	Undecane, 2,5-dimethyl-	184 C13H28	017301-22-3	81
3	Dodecane, 6-methyl-	184 C13H28	006044-71-9	70
4	Undecane, 3,6-dimethyl-	184 C13H28	017301-28-9	59
5	Decane, 2,6,8-trimethyl-	184 C13H28	062108-26-3	58



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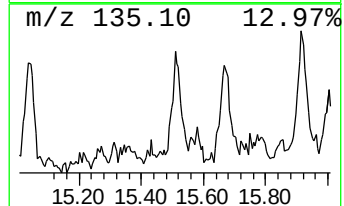
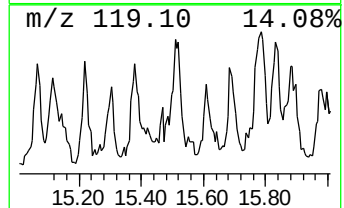
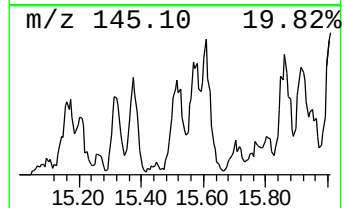
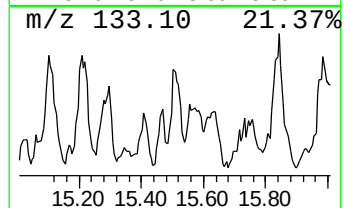
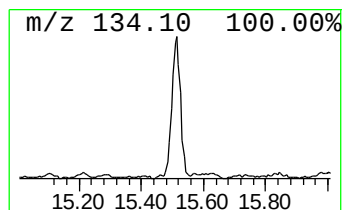
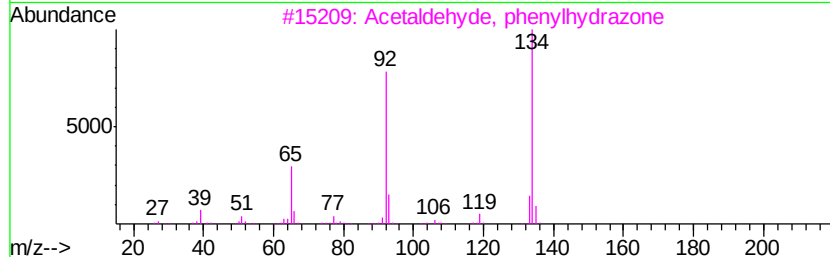
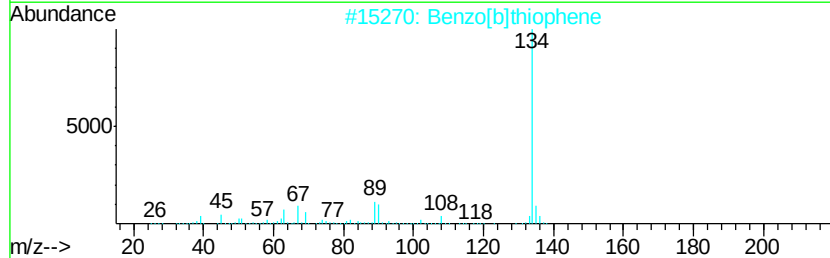
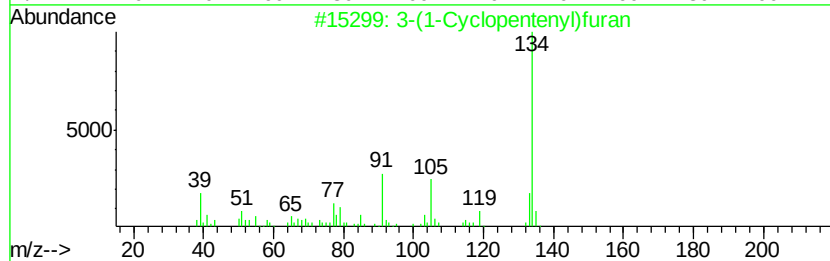
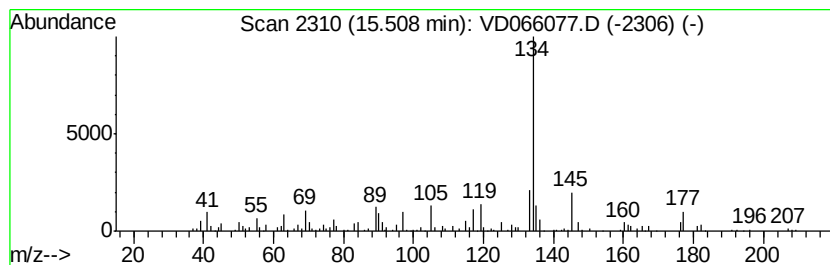
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TIC Integration Parameters: LSCINT.P

Peak Number 5 3-(1-Cyclopentenyl)furan Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.51	7.91 ug/l	226185	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-(1-Cyclopentenyl)furan	134	C9H10O	115754-79-5	50
2		Benzo[b]thiophene	134	C8H6S	000095-15-8	49
3		Acetaldehyde, phenylhydrazone	134	C8H10N2	000935-07-9	49
4		4(1H)-Pyrimidinone, tetrahydro-1...	306	C16H22N2O2S	088655-36-1	47
5		Benzo[c]thiophene	134	C8H6S	000270-82-6	47



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_D\DATA\VD072820\
Data File : VD066077.D
Acq On : 28 Jul 2020 13:49
Operator : VA/SY
Sample : L3181-06
Misc : 8.39G/5.00ml/MSVOA_D/SOIL
ALS Vial : 7 Sample Multiplier: 1

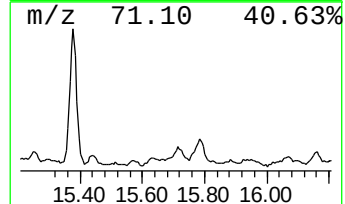
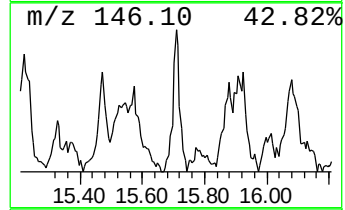
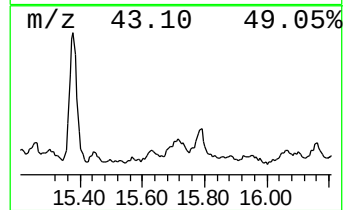
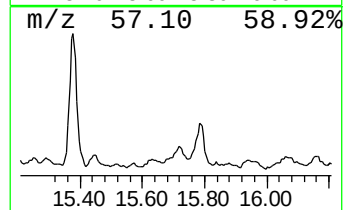
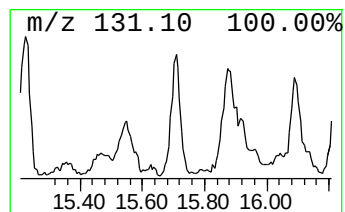
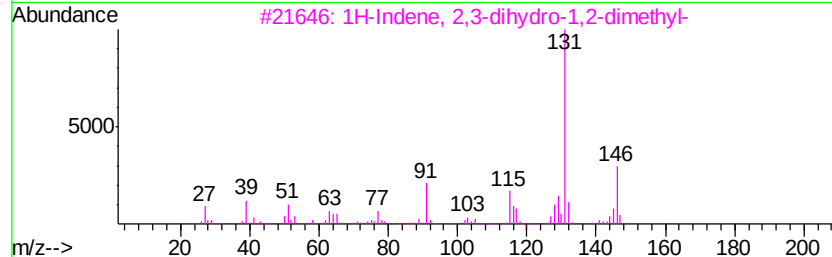
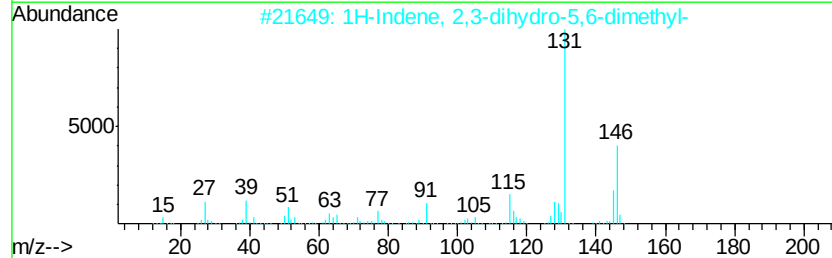
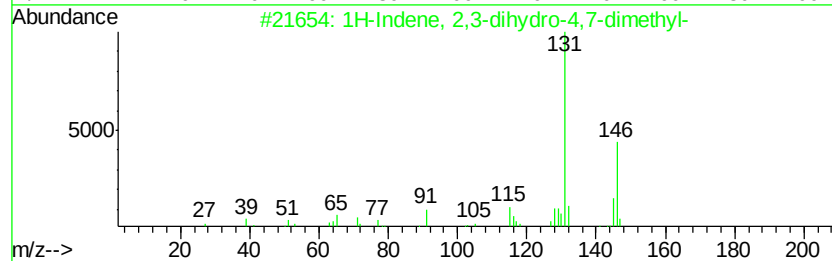
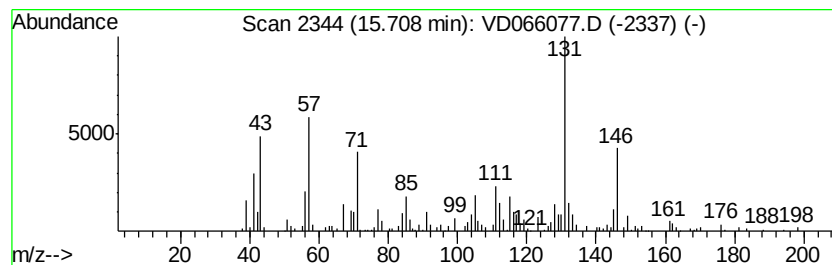
Instrument :
MSVOA_D
ClientSampleId :
CL-01-072720

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D072120S.M
Quant Title : SW846 8260

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

Peak Number 6 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.71	15.39 ug/l	439901	1,4-Dichlorobenzene-d4	13.58	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	86
2	1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	50
3	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	50
4	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	46
5	1H-Indene, 2,3-dihydro-4,6-dimet...	146	C11H14	001685-82-1	45



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_D\DATA\VD072820\
Data File : VD066077.D
Acq On : 28 Jul 2020 13:49
Operator : VA/SY
Sample : L3181-06
Misc : 8.39G/5.00ml/MSVOA_D/SOIL
ALS Vial : 7 Sample Multiplier: 1

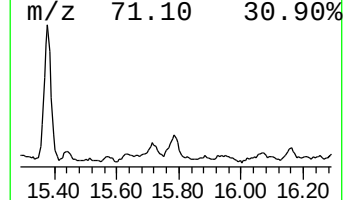
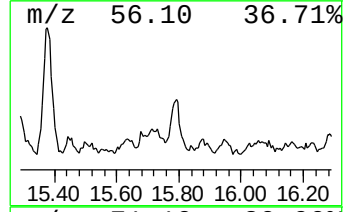
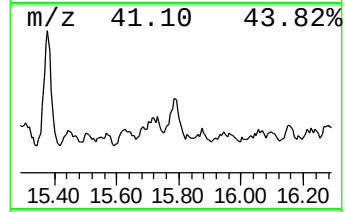
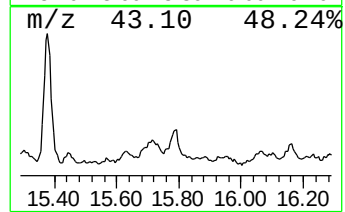
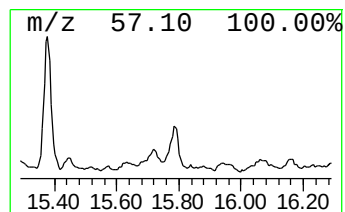
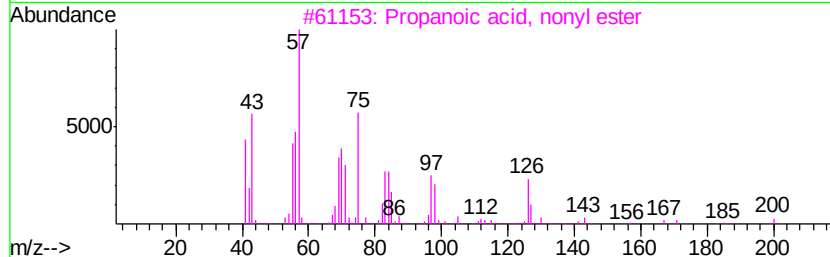
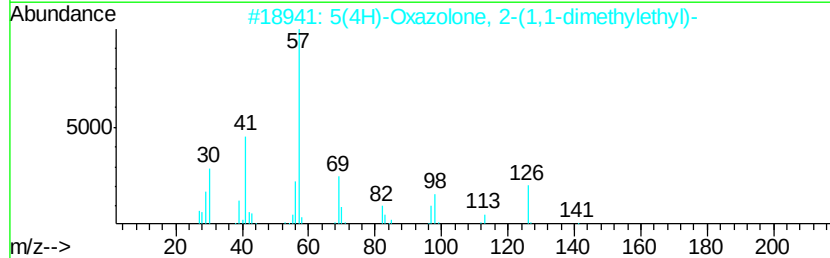
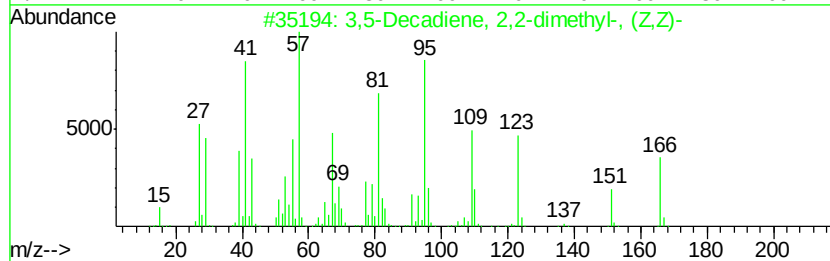
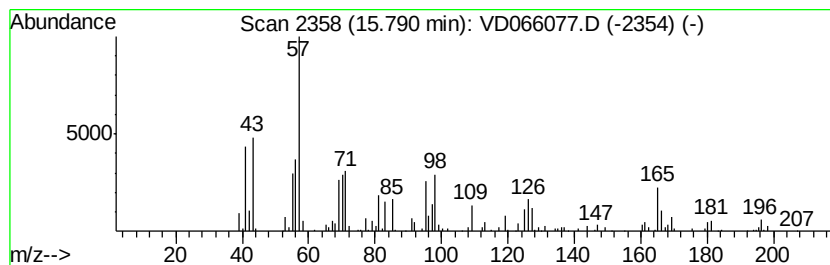
Instrument :
MSVOA_D
ClientSampleId :
CL-01-072720

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D072120S.M
Quant Title : SW846 8260

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

Peak Number 7 unknown15.79 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.79	12.63 ug/l	361027	1,4-Dichlorobenzene-d4	13.58	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,5-Decadiene, 2,2-dimethyl-, (Z...	166	C12H22	055638-50-1	18
2	5(4H)-Oxazolone, 2-(1,1-dimethyl...	141	C7H11NO2	081375-96-4	10
3	Propanoic acid, nonyl ester	200	C12H24O2	053184-67-1	10
4	Isobutyl nonyl carbonate	244	C14H28O3	959311-27-4	10
5	tert-Nonyl mercaptan	160	C9H20S	025360-10-5	10



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_D\DATA\VD072820\
Data File : VD066077.D
Acq On : 28 Jul 2020 13:49
Operator : VA/SY
Sample : L3181-06
Misc : 8.39G/5.00ml/MSVOA_D/SOIL
ALS Vial : 7 Sample Multiplier: 1

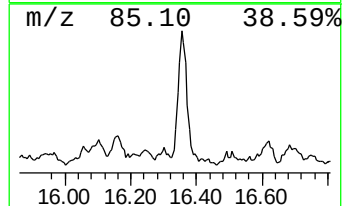
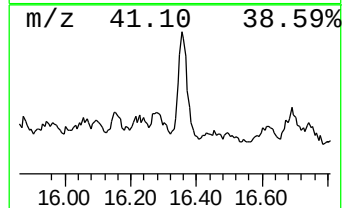
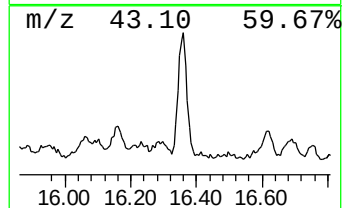
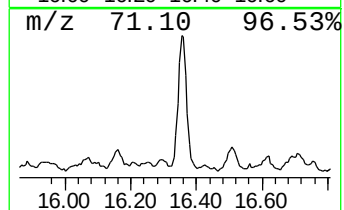
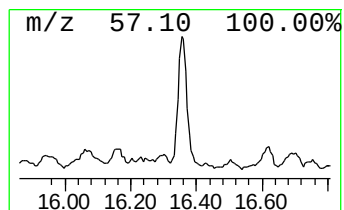
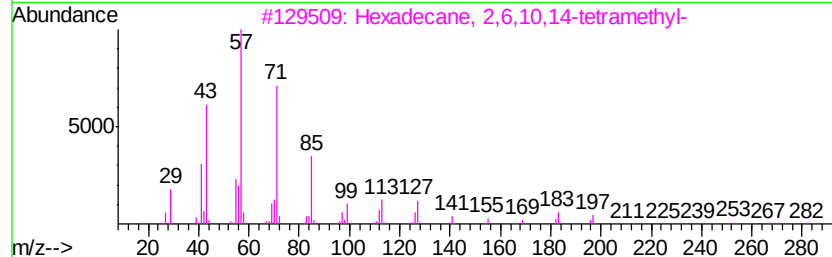
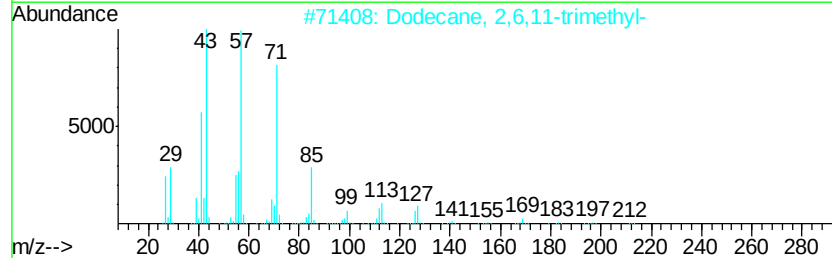
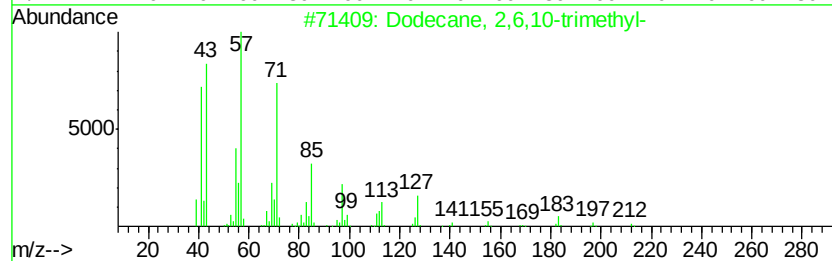
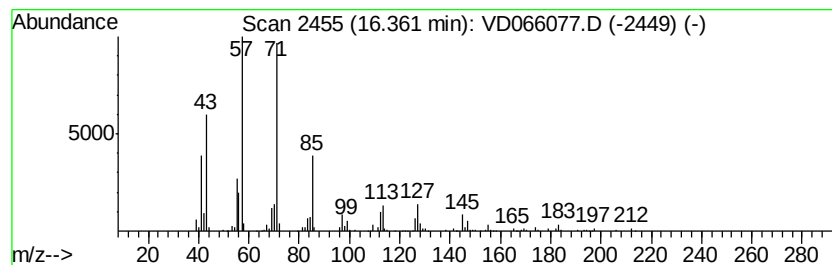
Instrument :
MSVOA_D
ClientSampleId :
CL-01-072720

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D072120S.M
Quant Title : SW846 8260

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

Peak Number 8 Dodecane, 2,6,10-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.36	24.09 ug/l	688569	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Dodecane, 2,6,10-trimethyl-	212 C15H32	003891-98-3	86
2	Dodecane, 2,6,11-trimethyl-	212 C15H32	031295-56-4	86
3	Hexadecane, 2,6,10,14-tetramethyl-	282 C20H42	000638-36-8	86
4	Nonane, 3-methyl-5-propyl-	184 C13H28	031081-18-2	80
5	Decane, 5-propyl-	184 C13H28	017312-62-8	74



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_D\DATA\VD072820\
Data File : VD066077.D
Acq On : 28 Jul 2020 13:49
Operator : VA/SY
Sample : L3181-06
Misc : 8.39G/5.00ml/MSVOA_D/SOIL
ALS Vial : 7 Sample Multiplier: 1

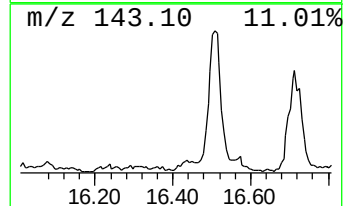
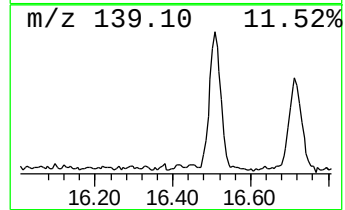
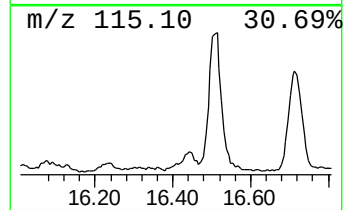
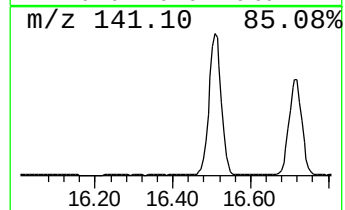
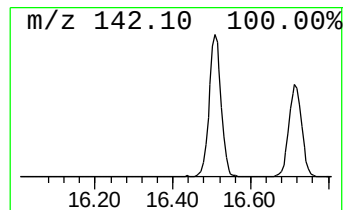
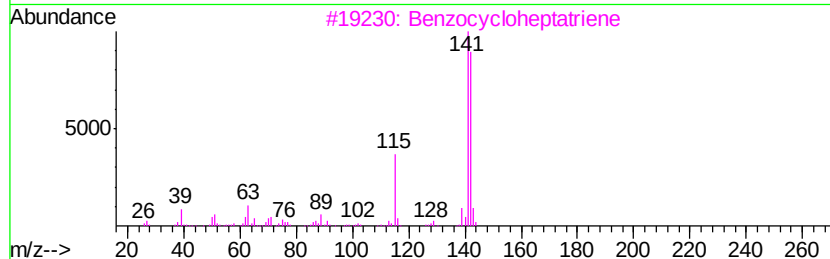
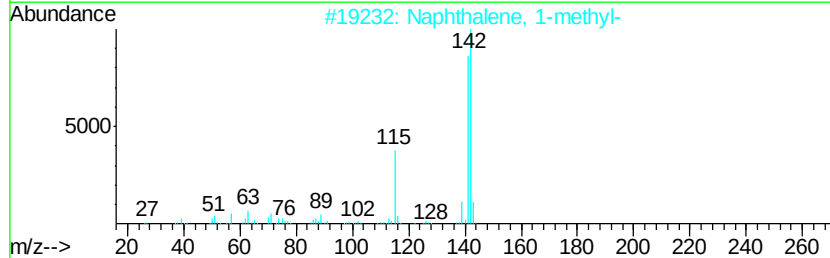
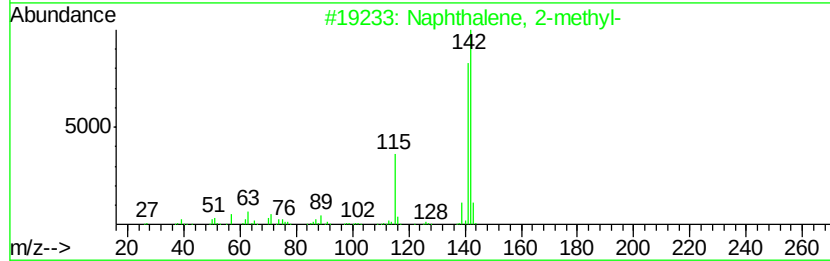
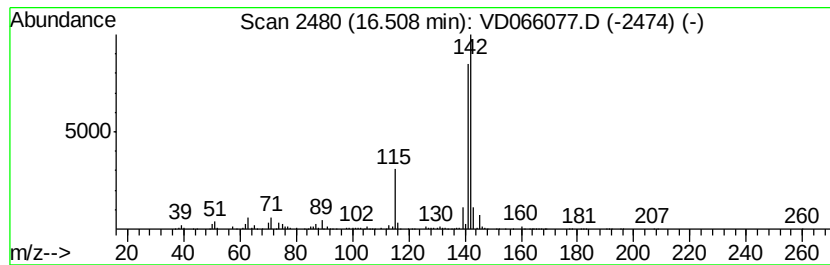
Instrument :
MSVOA_D
ClientSampleId :
CL-01-072720

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D072120S.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.51	52.34 ug/l	1495830	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2-methyl-	142 C11H10	000091-57-6 96
2		Naphthalene, 1-methyl-	142 C11H10	000090-12-0 96
3		Benzocycloheptatriene	142 C11H10	000264-09-5 91
4		1,4-Methanonaphthalene, 1,4-dihy...	142 C11H10	004453-90-1 91
5		1H-Indene, 1-ethylidene-	142 C11H10	002471-83-2 59



Data Path : Z:\V0ASRV\HPCHEM1\MSV0A_D\DATA\VD072820\
Data File : VD066077.D
Acq On : 28 Jul 2020 13:49
Operator : VA/SY
Sample : L3181-06
Misc : 8.39G/5.00ml/MSV0A_D/SOIL
ALS Vial : 7 Sample Multiplier: 1

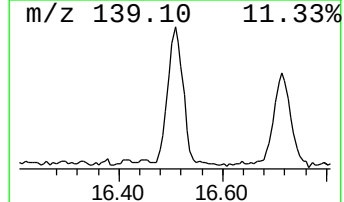
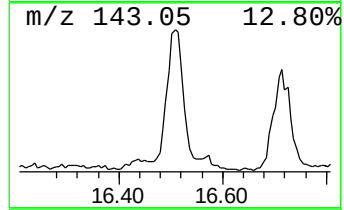
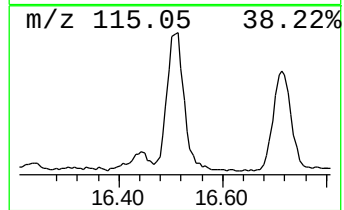
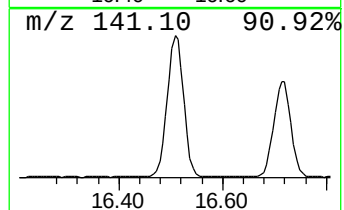
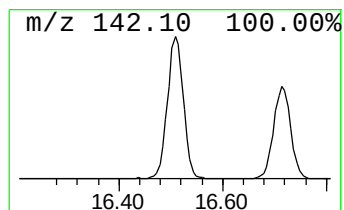
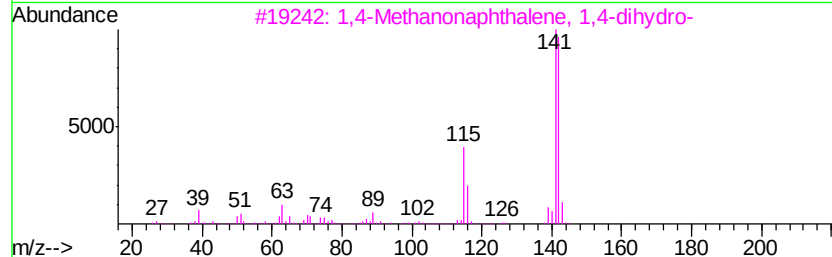
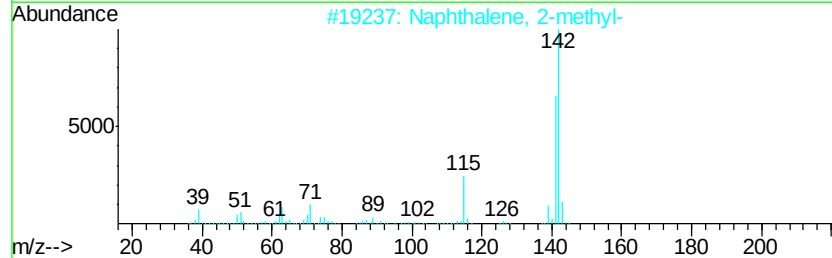
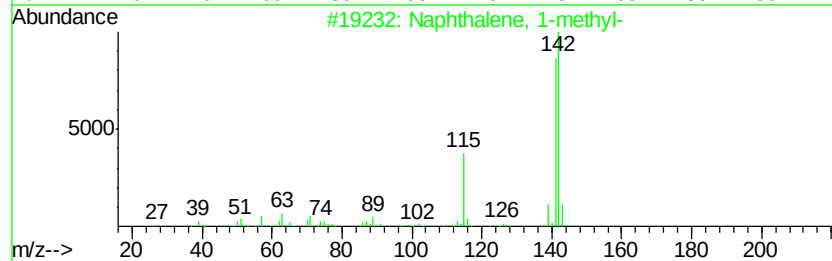
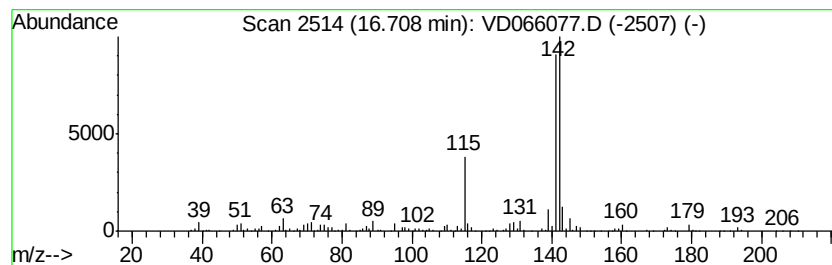
Instrument :
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ClientSampleId :
CL-01-072720

Quant Method : Z:\V0ASRV\HPCHEM1\MSV0A_D\METHOD\82D072120S.M
Quant Title : SW846 8260

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

Peak Number 10 1,4-Methanonaphthalene, 1,4... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.71	54.19 ug/l	1548510	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1-methyl-	142 C11H10	000090-12-0 95
2		Naphthalene, 2-methyl-	142 C11H10	000091-57-6 93
3		1,4-Methanonaphthalene, 1,4-dihy...	142 C11H10	004453-90-1 91
4		Benzocycloheptatriene	142 C11H10	000264-09-5 87
5		Naphthalene, 1-(2-hydroxypropyl)	186 C13H14O	027653-13-0 78



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_D\DATA\VD072820\
Data File : VD066077.D
Acq On : 28 Jul 2020 13:49
Operator : VA/SY
Sample : L3181-06
Misc : 8.39G/5.00ml/MSVOA_D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
CL-01-072720

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D072120S.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
Benzene, 1-propynyl-	13.97	13.5	ug/l	386385	4	13.58	1428870	50.0
Bicyclo(3.3.1)non...	14.42	7.8	ug/l	222651	4	13.58	1428870	50.0
1H-Indene, 2,3-di...	14.84	12.7	ug/l	363846	4	13.58	1428870	50.0
Undecane, 2,6-dim...	14.88	27.0	ug/l	771433	4	13.58	1428870	50.0
3-(1-Cyclopentenyl...	15.51	7.9	ug/l	226185	4	13.58	1428870	50.0
1H-Indene, 2,3-di...	15.71	15.4	ug/l	439901	4	13.58	1428870	50.0
unknown15.79	15.79	12.6	ug/l	361027	4	13.58	1428870	50.0
Dodecane, 2,6,10-...	16.36	24.1	ug/l	688569	4	13.58	1428870	50.0
Naphthalene, 1-me...	16.51	52.3	ug/l	1495830	4	13.58	1428870	50.0
1,4-Methanonaphth...	16.71	54.2	ug/l	1548510	4	13.58	1428870	50.0