

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW083018\
 Data File : VW005023.D
 Acq On : 30 Aug 2018 17:15
 Operator : SY/AP
 Sample : J4648-02
 Misc : 6.45G/5ML/MSVOA W/SOIL
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 SB-1-@-8-FT

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_W\METHOD\82W082918S.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	1.794	18	34	56	rBV	4118770	7417369	100.00%	15.970%
2	1.959	57	61	72	rVB2	100866	170402	2.30%	0.367%
3	2.355	117	126	135	rVB	170981	292744	3.95%	0.630%
4	2.818	195	202	212	rVB	207124	384094	5.18%	0.827%
5	3.026	228	236	248	rVB	579405	1221448	16.47%	2.630%
6	3.391	287	296	310	rVB	179983	408420	5.51%	0.879%
7	4.123	401	416	417	rBV3	83554	217679	2.93%	0.469%
8	4.385	447	459	468	rBV	80229	229271	3.09%	0.494%
9	4.909	530	545	553	rBV4	296819	1280860	17.27%	2.758%
10	5.446	618	633	648	rBV	253987	857685	11.56%	1.847%
11	5.946	703	715	728	rBV2	49771	151552	2.04%	0.326%
12	6.720	831	842	850	rBV3	25293	83782	1.13%	0.180%
13	6.879	861	868	875	rBV4	31857	89955	1.21%	0.194%
14	6.988	875	886	896	rBV	1680647	4951170	66.75%	10.660%
15	7.153	905	913	933	rVB	238027	790085	10.65%	1.701%
16	7.641	985	993	998	rBV4	43619	126852	1.71%	0.273%
17	7.702	998	1003	1012	rVB2	58191	155106	2.09%	0.334%
18	7.891	1020	1034	1037	rBV	156385	384546	5.18%	0.828%
19	7.958	1037	1045	1054	rVV	2552156	6304138	84.99%	13.574%
20	8.037	1054	1058	1070	rVV2	203247	490037	6.61%	1.055%
21	8.165	1070	1079	1082	rVV2	59577	129565	1.75%	0.279%
22	8.226	1082	1089	1097	rVV	320134	780868	10.53%	1.681%
23	8.317	1097	1104	1114	rVV2	191812	479929	6.47%	1.033%
24	8.439	1114	1124	1130	rVV3	282712	684870	9.23%	1.475%
25	8.512	1130	1136	1141	rVV3	306445	681129	9.18%	1.467%
26	8.580	1141	1147	1157	rVV	434567	961156	12.96%	2.069%
27	8.848	1182	1191	1203	rBV	640551	1513460	20.40%	3.259%
28	9.079	1221	1229	1232	rBV	119654	245742	3.31%	0.529%
29	9.122	1232	1236	1254	rVB	187426	386887	5.22%	0.833%
30	9.342	1257	1272	1279	rBV	1041611	2754136	37.13%	5.930%
31	9.524	1296	1302	1308	rVB	78614	146322	1.97%	0.315%
32	9.610	1308	1316	1323	rVB	106354	220529	2.97%	0.475%
33	9.756	1335	1340	1348	rVB2	151290	310540	4.19%	0.669%
34	9.854	1348	1356	1363	rBV2	703321	1312643	17.70%	2.826%

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35	10.018	1377	1383	1388	rVB	43137	85528	1.15%	0.184%
36	10.085	1388	1394	1401	rBV2	84886	171637	2.31%	0.370%
37	10.207	1407	1414	1420	rBV	377986	703105	9.48%	1.514%
38	10.329	1427	1434	1441	rBV	751964	1455185	19.62%	3.133%
39	10.500	1451	1462	1473	rBV2	152457	448063	6.04%	0.965%
40	10.671	1484	1490	1495	rBV	195078	362827	4.89%	0.781%
41	10.750	1495	1503	1512	rVB4	355899	1013165	13.66%	2.181%
42	11.024	1541	1548	1551	rBV	95328	168282	2.27%	0.362%
43	11.341	1595	1600	1606	rVB2	59320	95127	1.28%	0.205%
44	11.634	1641	1648	1654	rBV	461920	811303	10.94%	1.747%
45	11.695	1654	1658	1661	rVV2	53312	99829	1.35%	0.215%
46	11.731	1661	1664	1669	rVB2	59468	91846	1.24%	0.198%
47	11.847	1674	1683	1699	rVB	341032	783150	10.56%	1.686%
48	12.170	1725	1736	1746	rBV	163844	313963	4.23%	0.676%
49	12.469	1780	1785	1790	rBV	64486	102058	1.38%	0.220%
50	12.621	1803	1810	1819	rVB	298685	508848	6.86%	1.096%
51	12.945	1859	1863	1870	rVB2	147642	251324	3.39%	0.541%
52	13.109	1883	1890	1898	rBV	258433	427863	5.77%	0.921%
53	13.256	1909	1914	1926	rVB	280425	456746	6.16%	0.983%
54	13.566	1958	1965	1968	rBV	258237	464994	6.27%	1.001%
55	13.768	1987	1998	2008	rBV	435326	722273	9.74%	1.555%
56	14.219	2067	2072	2079	rVB	118968	187840	2.53%	0.404%
57	14.816	2162	2170	2175	rBV2	51777	104509	1.41%	0.225%

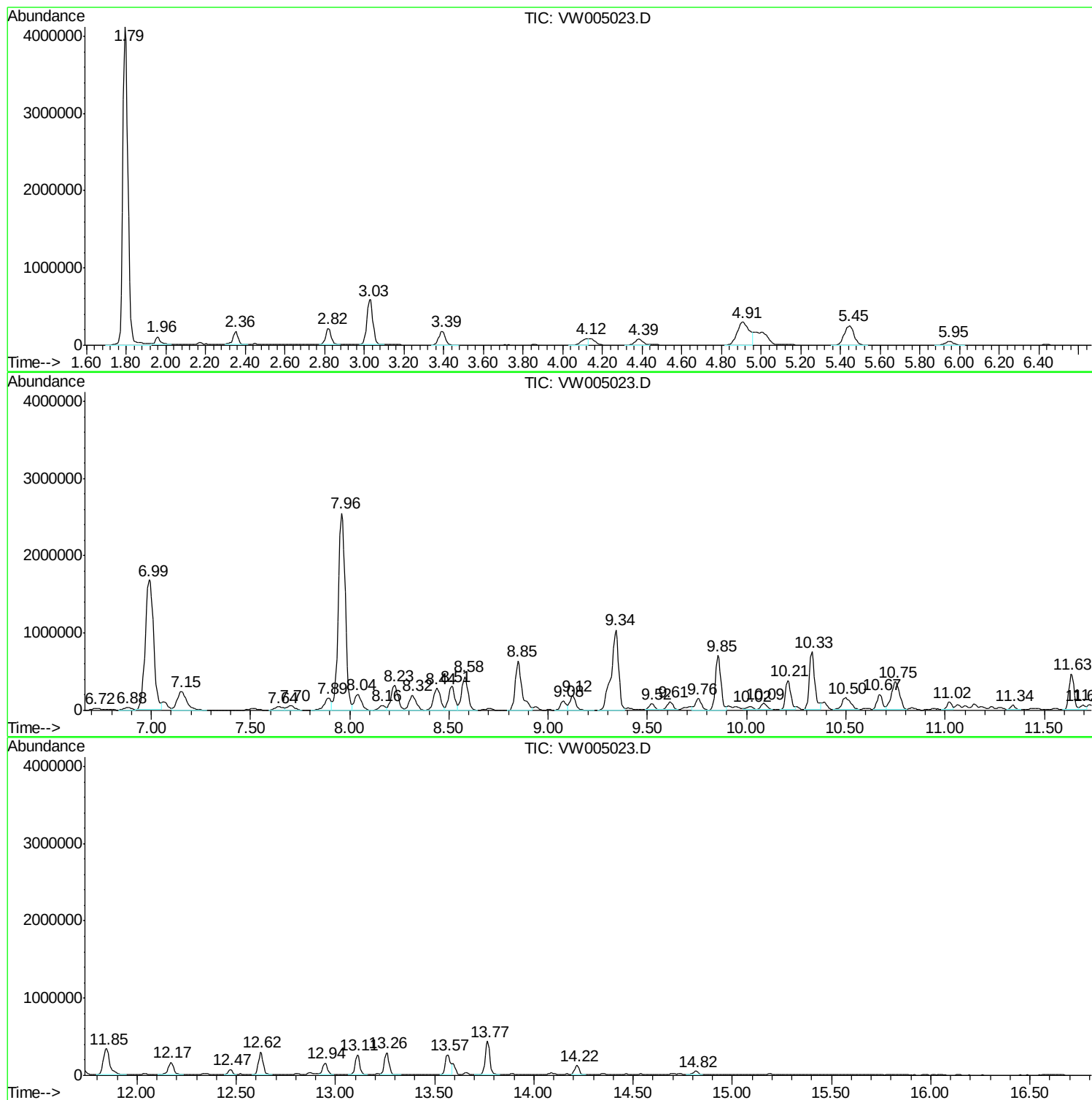
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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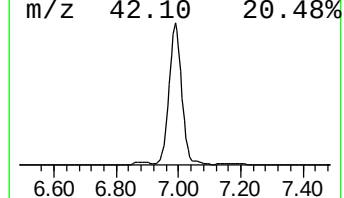
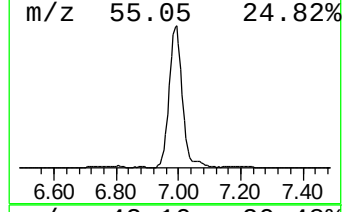
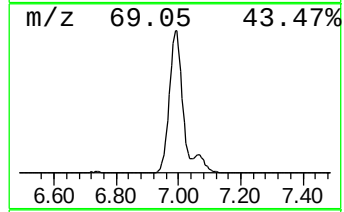
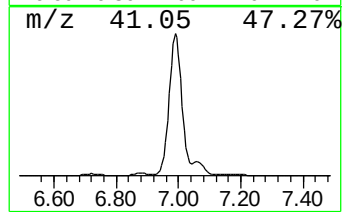
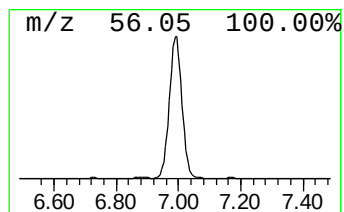
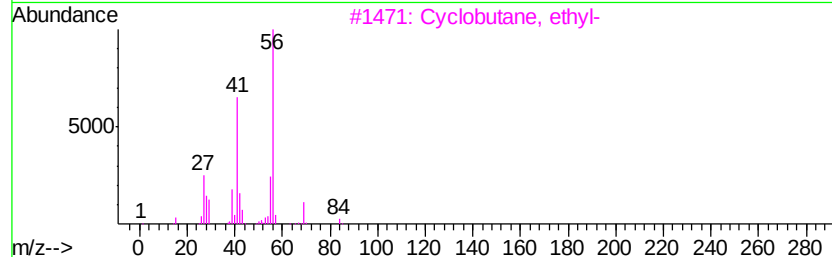
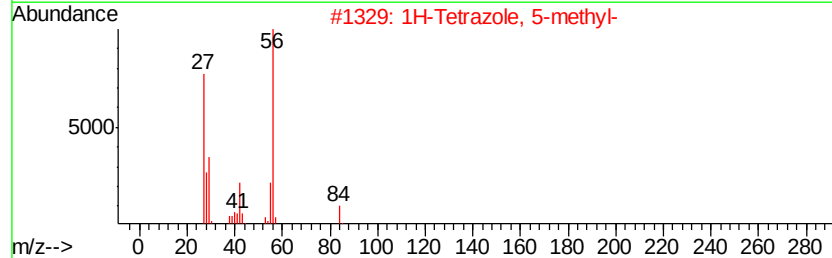
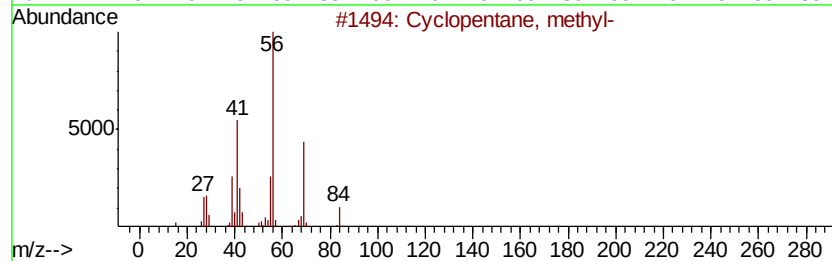
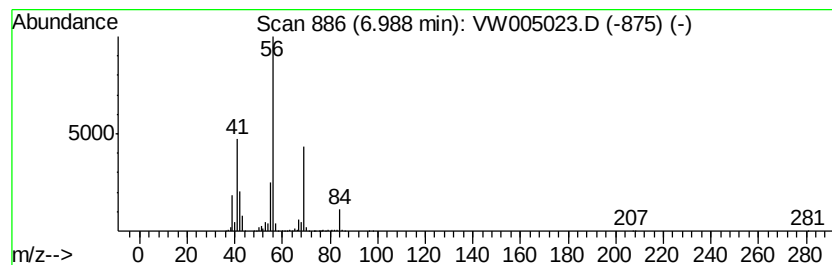
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W082918S.M
Quant Title : SW846 8260

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

Peak Number 2 Cyclopentane, methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.99	39.27 ug/l	4951170	Pentafluorobenzene	7.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	80
3		Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
4		Cyclobutane, butyl-	112	C8H16	013152-44-8	50
5		Cyclohexane	84	C6H12	000110-82-7	47



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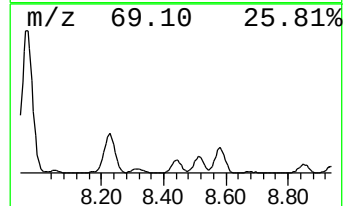
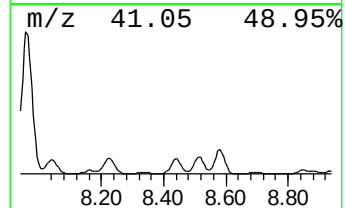
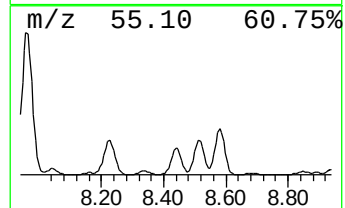
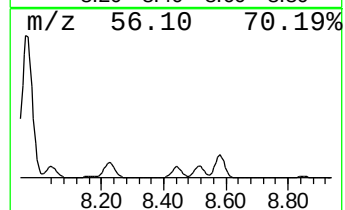
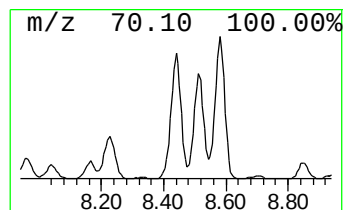
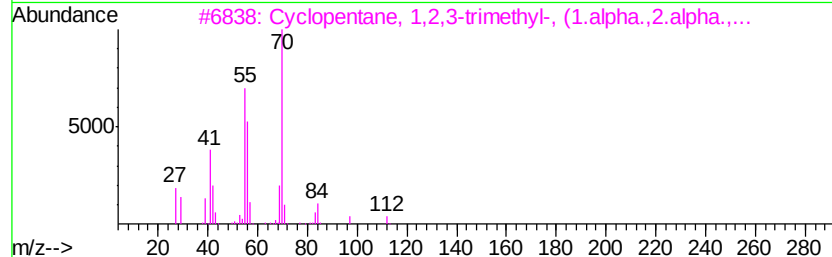
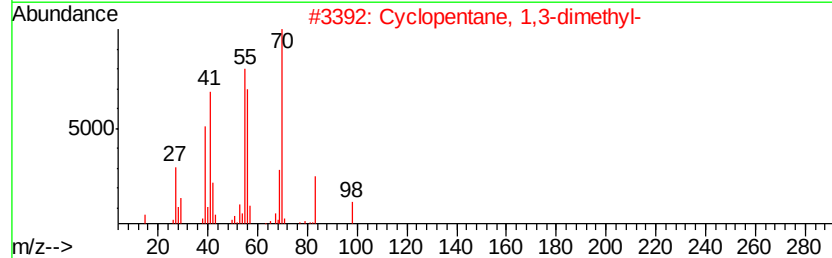
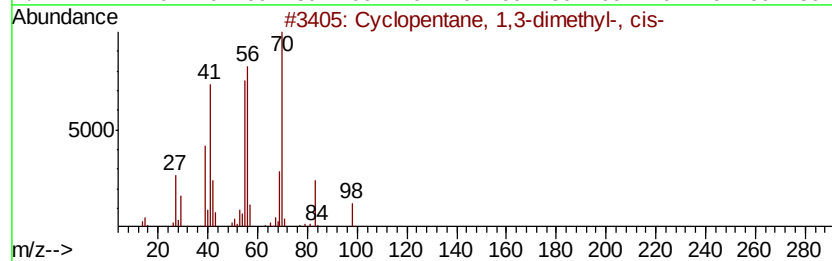
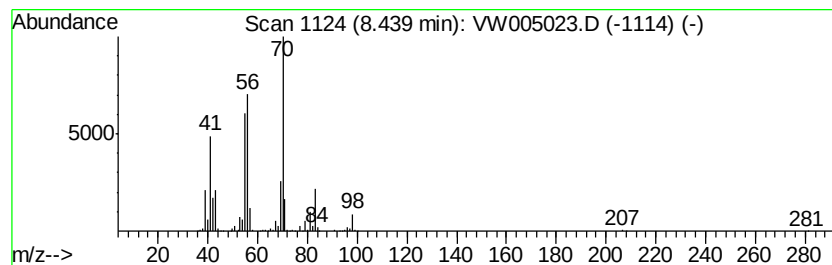
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W082918S.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Cyclopentane, 1,3-dimethyl-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.44	22.63 ug/l	684870	1,4-Difluorobenzene	8.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	90
2		Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	87
3		Cyclopentane, 1,2,3-trimethyl-, ...	112	C8H16	002613-69-6	72
4		Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	70
5		Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	53



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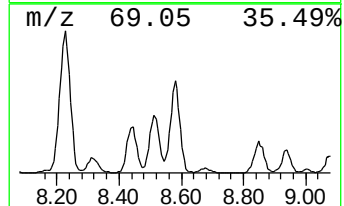
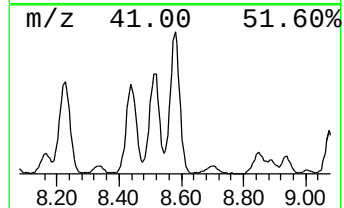
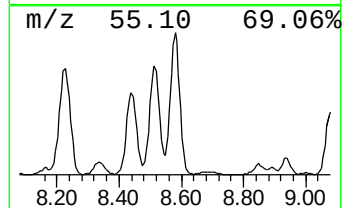
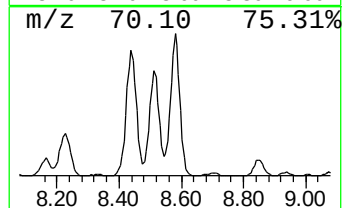
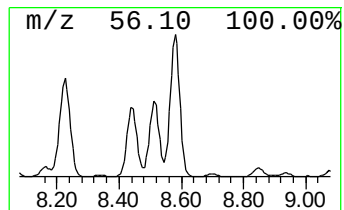
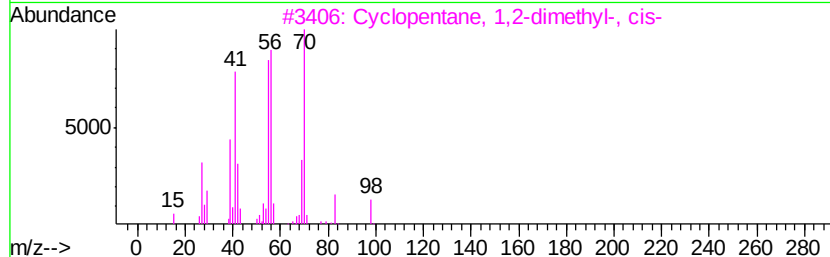
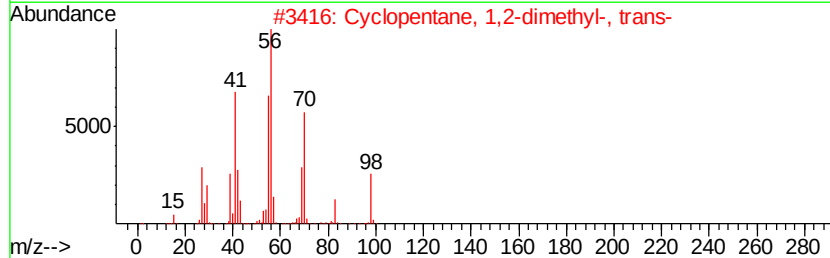
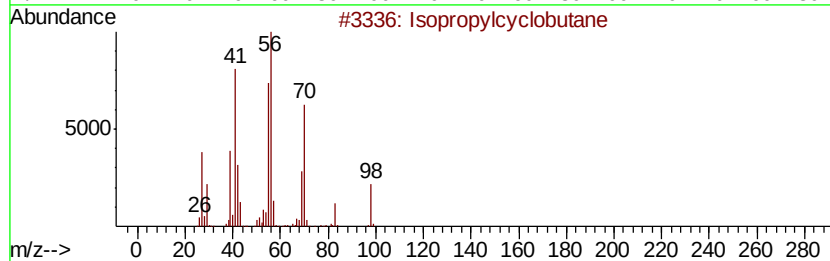
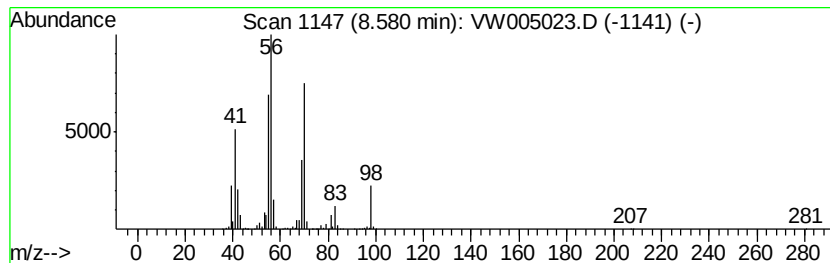
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 Isopropylcyclobutane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.58	31.75 ug/l	961156	1,4-Difluorobenzene	8.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Isopropylcyclobutane	98	C7H14	000872-56-0	94
2		Cyclopentane, 1,2-dimethyl-, trans-	98	C7H14	000822-50-4	91
3		Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	78
4		Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	64
5		1-Heptene	98	C7H14	000592-76-7	53



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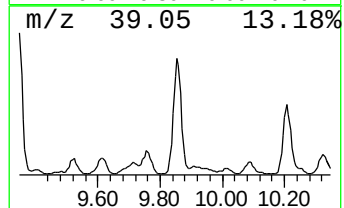
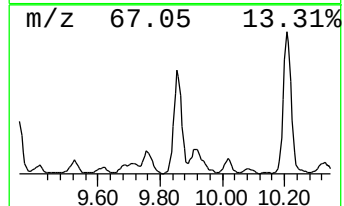
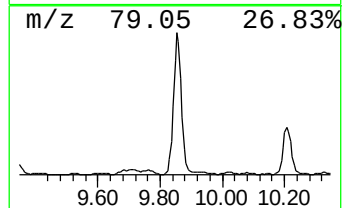
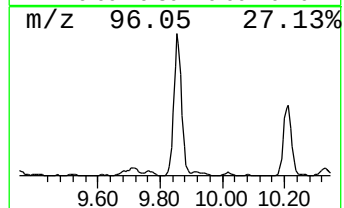
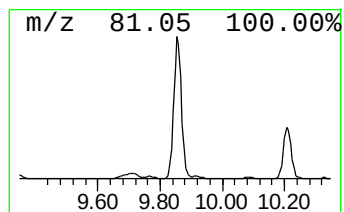
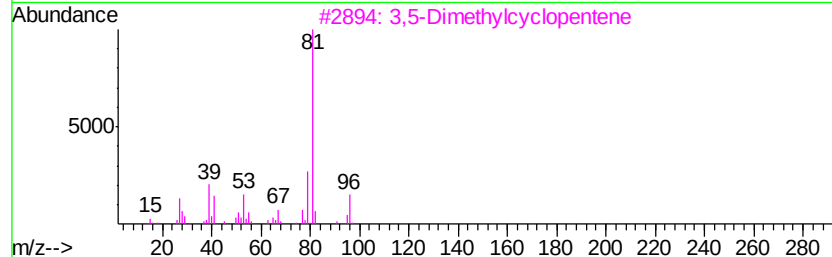
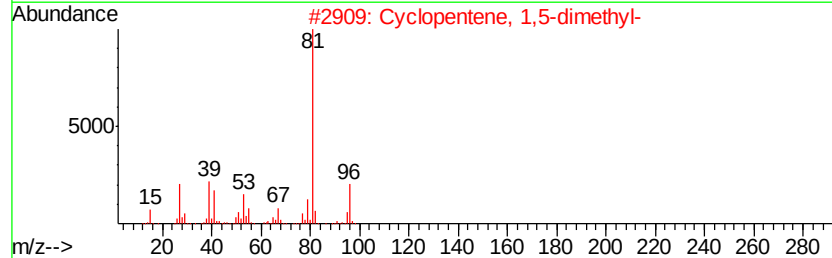
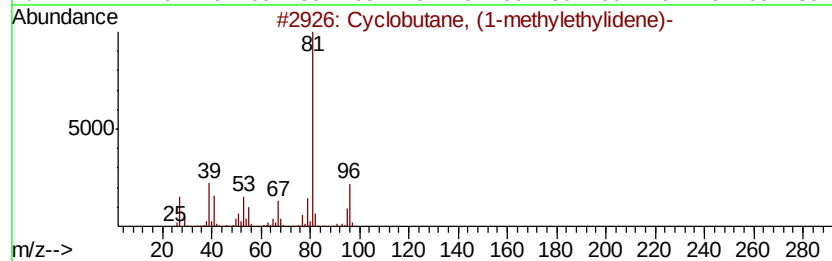
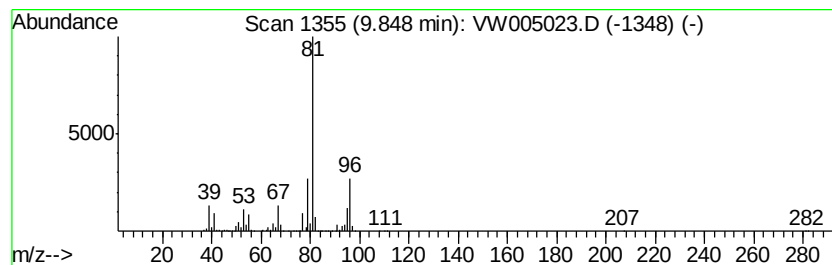
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Cyclobutane, (1-methylethyl... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.85	43.37 ug/l	1312640	1,4-Difluorobenzene	8.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclobutane, (1-methylethylidene)-	96	C7H12	001528-22-9	87
2		Cyclopentene, 1,5-dimethyl-	96	C7H12	016491-15-9	87
3		3,5-Dimethylcyclopentene	96	C7H12	007459-71-4	78
4		Cyclopropane, 1-methyl-1-isopropyl-	96	C7H12	003422-07-9	78
5		Cyclopentene, 4,4-dimethyl-	96	C7H12	019037-72-0	64



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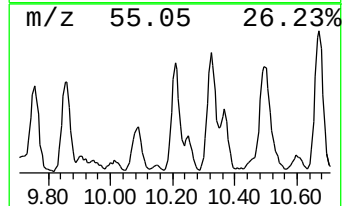
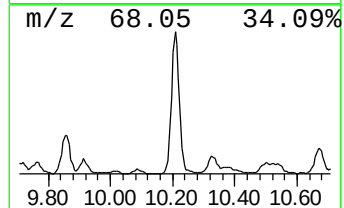
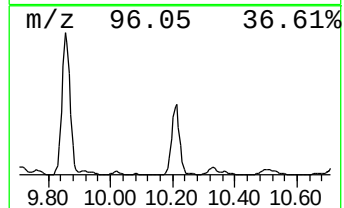
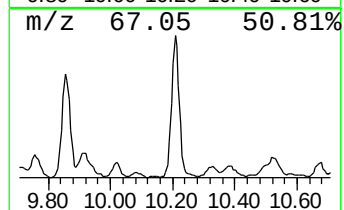
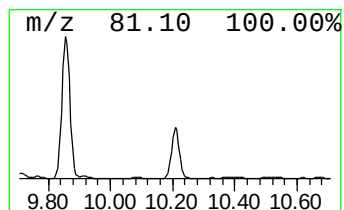
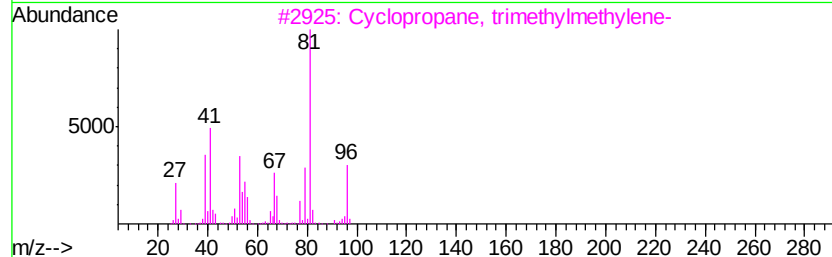
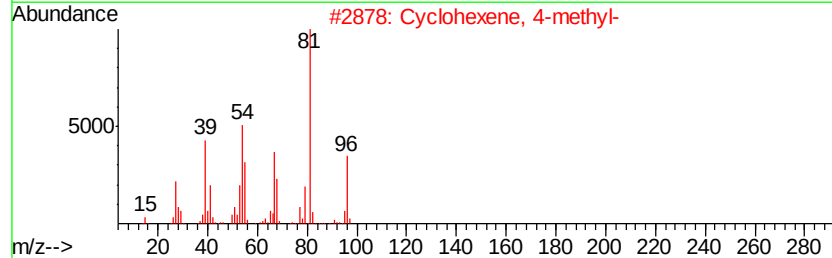
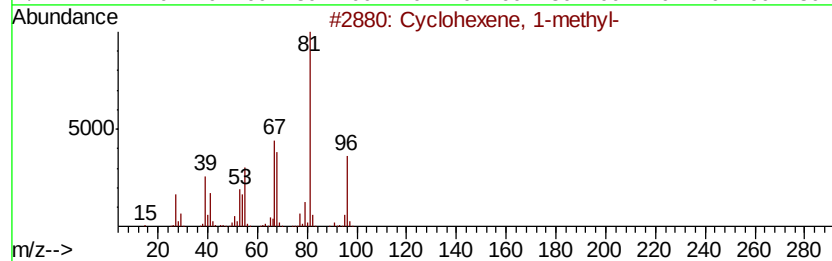
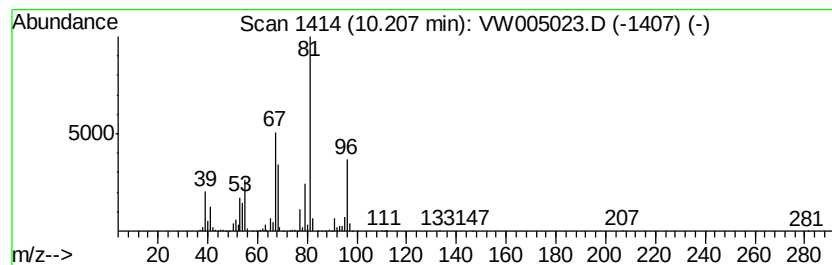
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Quant Title : SW846 8260

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

Peak Number 6 Cyclohexene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.21	23.23 ug/l	703105	1,4-Difluorobenzene	8.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexene, 1-methyl-	96	C7H12	000591-49-1	95
2		Cyclohexene, 4-methyl-	96	C7H12	000591-47-9	87
3		Cyclopropane, trimethylmethylen-	96	C7H12	034462-28-7	87
4		Cyclohexene, 3-methyl-	96	C7H12	000591-48-0	81
5		2-Hexyne, 4-methyl-	96	C7H12	020198-49-6	80



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW083018\
Data File : VW005023.D
Acq On : 30 Aug 2018 17:15
Operator : SY/AP
Sample : J4648-02
Misc : 6.45G/5ML/MSVOA W/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
SB-1-@-8-FT

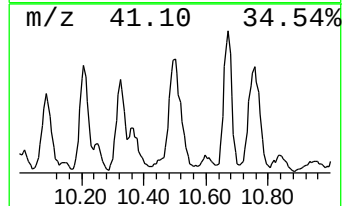
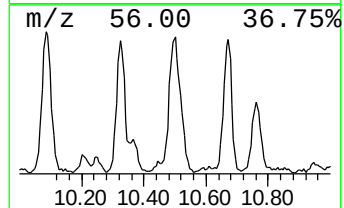
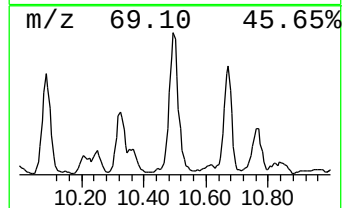
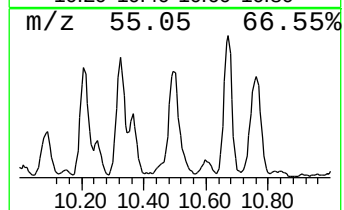
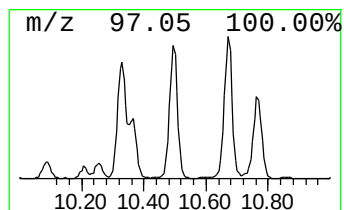
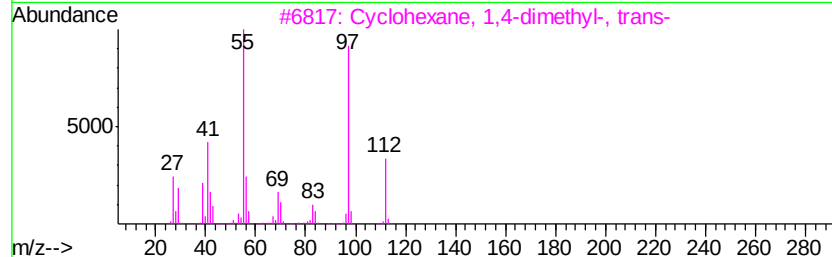
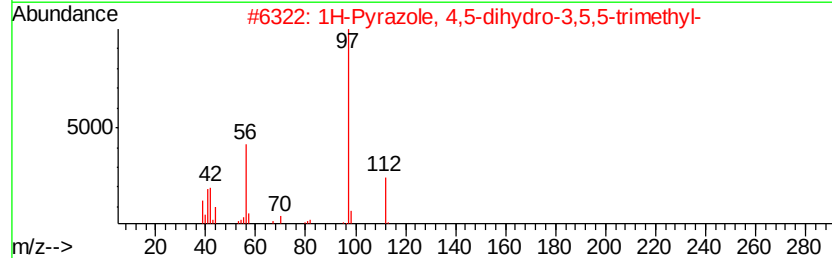
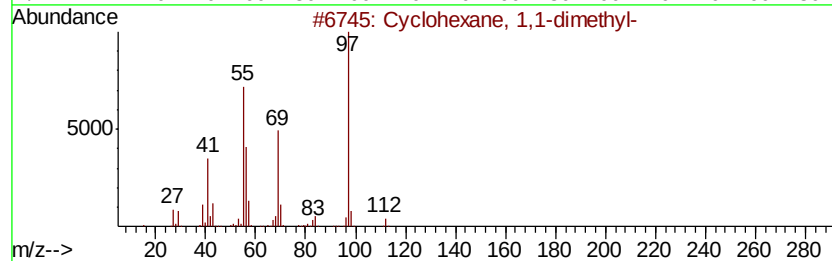
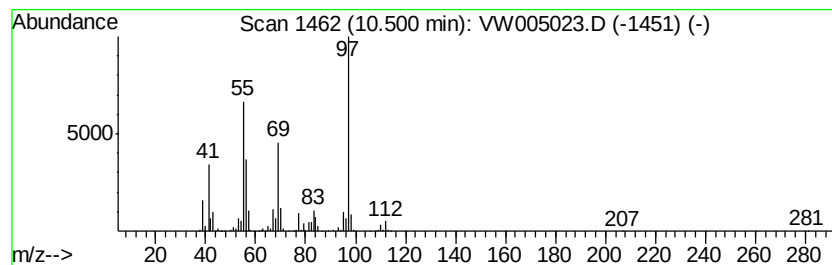
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Quant Title : SW846 8260

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

Peak Number 7 Cyclohexane, 1,1-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.50	27.61 ug/l	448063	Chlorobenzene-d5	11.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1,1-dimethyl-	112	C8H16	000590-66-9	94
2		1H-Pyrazole, 4,5-dihydro-3,5,5-t...	112	C6H12N2	003975-85-7	64
3		Cyclohexane, 1,4-dimethyl-, trans-	112	C8H16	002207-04-7	59
4		Oxalic acid, cyclohexylmethyl oc...	298	C17H30O4	1000309-68-4	59
5		Cyclohexane, 1-bromo-2-methyl-	176	C7H13Br	006294-39-9	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW083018\
Data File : VW005023.D
Acq On : 30 Aug 2018 17:15
Operator : SY/AP
Sample : J4648-02
Misc : 6.45G/5ML/MSVOA W/SOIL
ALS Vial : 15 Sample Multiplier: 1

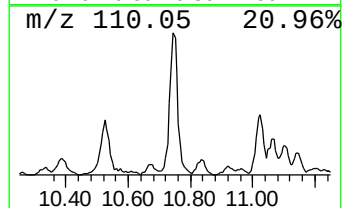
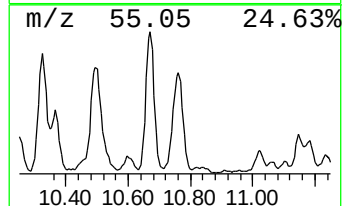
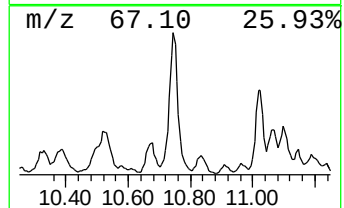
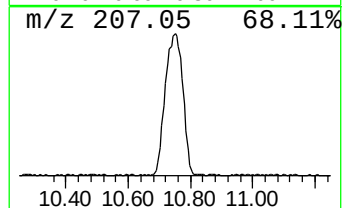
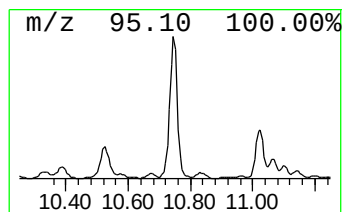
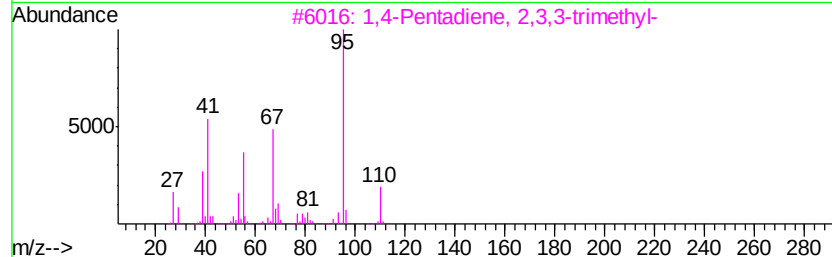
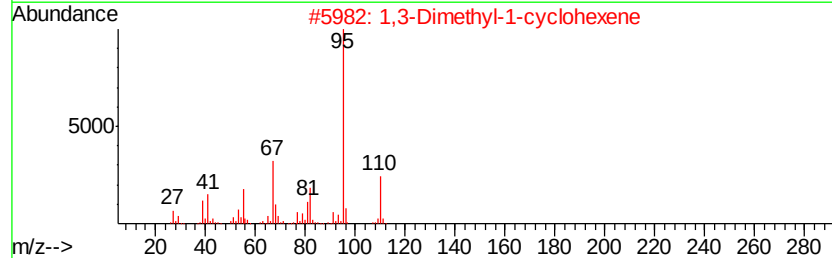
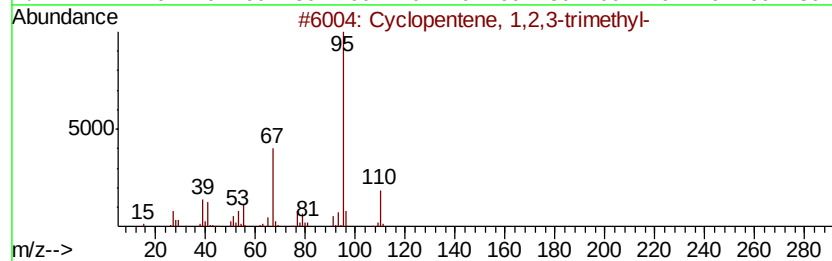
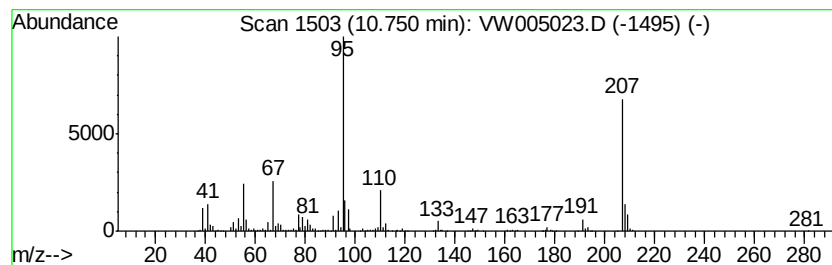
Instrument :
MSVOA_W
ClientSampleId :
SB-1-@-8-FT

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W082918S.M
Quant Title : SW846 8260

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

Peak Number 8 Cyclopentene, 1,2,3-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.75	62.44 ug/l	1013170	Chlorobenzene-d5	11.63		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentene, 1,2,3-trimethyl-	110	C8H14	000473-91-6	53
2		1,3-Dimethyl-1-cyclohexene	110	C8H14	002808-76-6	50
3		1,4-Pentadiene, 2,3,3-trimethyl-	110	C8H14	000756-02-5	47
4		exo-2-Bromonorbornane	174	C7H11Br	002534-77-2	43
5		Furan, 2-ethyl-5-methyl-	110	C7H10O	001703-52-2	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW083018\
Data File : VW005023.D
Acq On : 30 Aug 2018 17:15
Operator : SY/AP
Sample : J4648-02
Misc : 6.45G/5ML/MSVOA W/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
SB-1-@-8-FT

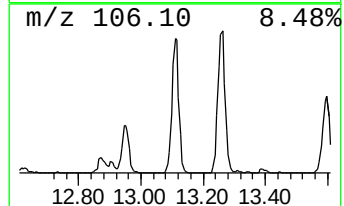
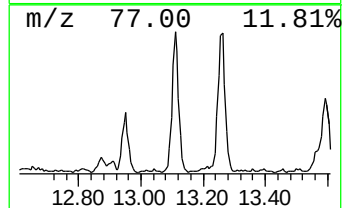
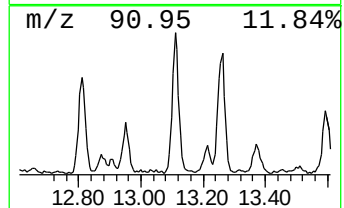
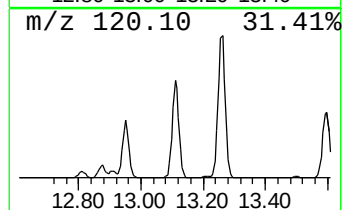
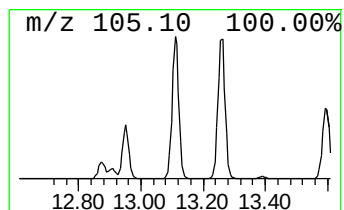
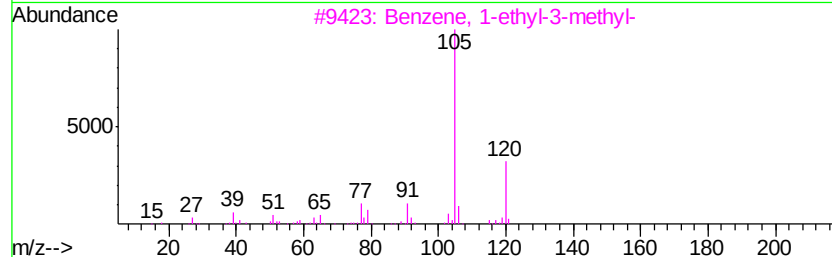
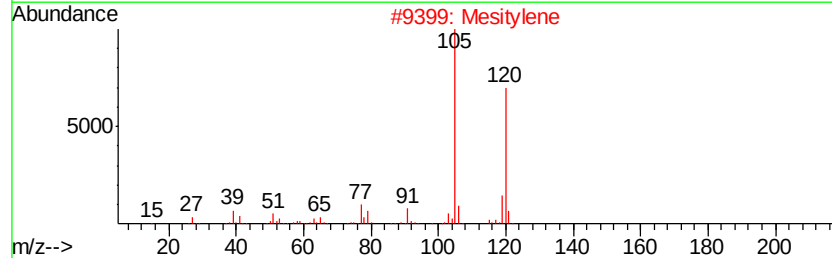
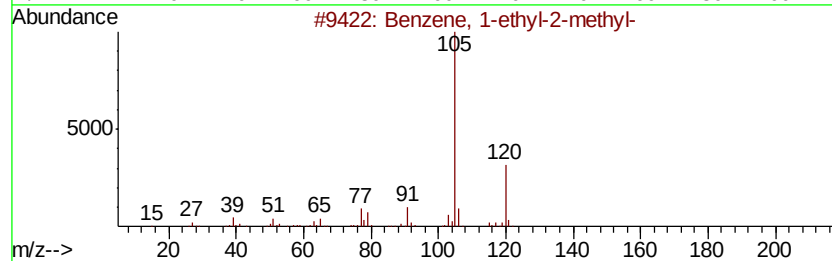
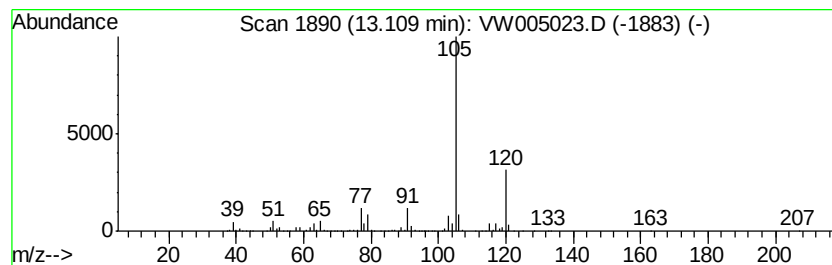
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W082918S.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-ethyl-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.11	46.01 ug/l	427863	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Mesitylene	120	C9H12	000108-67-8	91
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW083018\
Data File : VW005023.D
Acq On : 30 Aug 2018 17:15
Operator : SY/AP
Sample : J4648-02
Misc : 6.45G/5ML/MSVOA W/SOIL
ALS Vial : 15 Sample Multiplier: 1

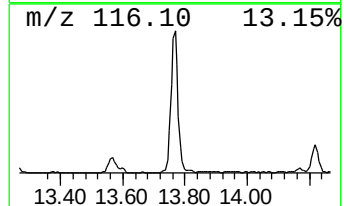
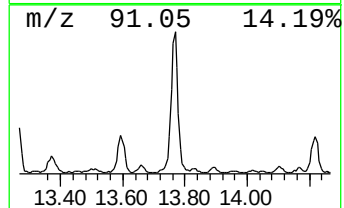
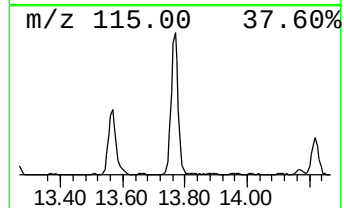
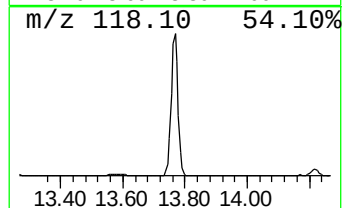
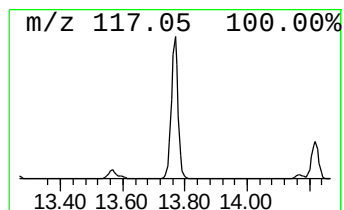
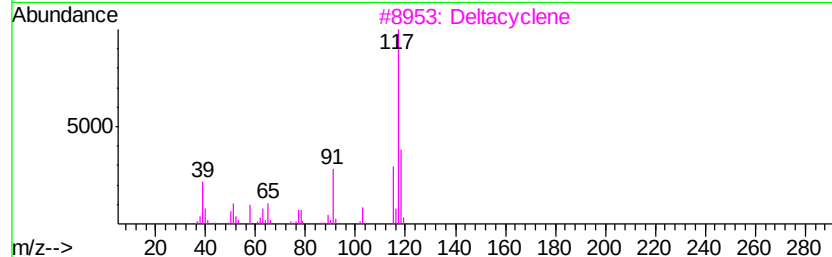
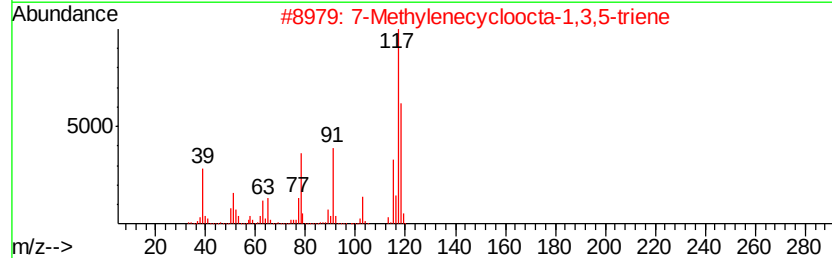
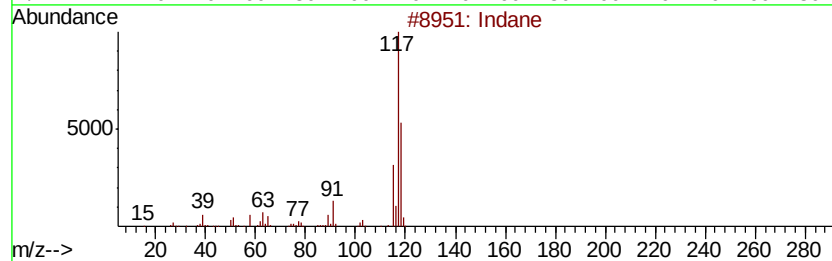
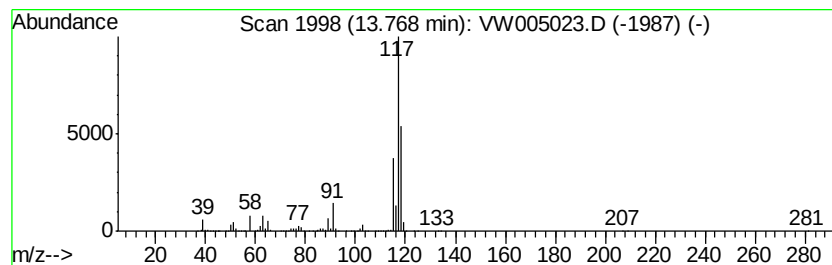
Instrument :
MSVOA_W
ClientSampleId :
SB-1-@-8-FT

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W082918S.M
Quant Title : SW846 8260

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

Peak Number 10 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.77	77.66 ug/l	722273	1,4-Dichlorobenzene-d4	13.57
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Indane		118 C9H10	000496-11-7 95
2	7-Methylenecycloocta-1,3,5-triene		118 C9H10	002570-13-0 74
3	Deltacyclene		118 C9H10	007785-10-6 72
4	Benzene, 1-ethenyl-4-methyl-		118 C9H10	000622-97-9 64
5	1,3-Methanopentalene, 1,2,3,5-te...		118 C9H10	128600-88-4 64



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_W\DATA\VW083018\
Data File : VW005023.D
Acq On : 30 Aug 2018 17:15
Operator : SY/AP
Sample : J4648-02
Misc : 6.45G/5ML/MSVOA_W/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
SB-1-@-8-FT

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W082918S.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
Cyclopentane, met...	6.99	39.3	ug/l	4951170	1	7.96	6304140	50.0
Cyclopentane, 1,3...	8.44	22.6	ug/l	684870	2	8.85	1513460	50.0
Isopropylcyclobutane	8.58	31.8	ug/l	961156	2	8.85	1513460	50.0
Cyclobutane, (1-m...	9.85	43.4	ug/l	1312640	2	8.85	1513460	50.0
Cyclohexene, 1-me...	10.21	23.2	ug/l	703105	2	8.85	1513460	50.0
Cyclohexane, 1,1-...	10.50	27.6	ug/l	448063	3	11.63	811303	50.0
Cyclopentene, 1,2...	10.75	62.4	ug/l	1013170	3	11.63	811303	50.0
Benzene, 1-ethyl-...	13.11	46.0	ug/l	427863	4	13.57	464994	50.0
Indane	13.77	77.7	ug/l	722273	4	13.57	464994	50.0