

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW011019\
 Data File : VW008182.D
 Acq On : 11 Jan 2019 07:46
 Operator : SY/AP
 Sample : K1102-01
 Misc : 5.09G/5ML/MSVOA W/SOIL
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 CPSB-18(15-20)

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\82W010819S.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	6.879	805	816	825	rVV	700875	2465278	1.00%	0.481%
2	7.939	973	990	1000	rBV4	6775979	27348606	11.10%	5.336%
3	8.037	1000	1006	1018	rVV	3693026	10296440	4.18%	2.009%
4	8.165	1018	1027	1032	rVV	5890723	13886278	5.64%	2.709%
5	8.226	1032	1037	1048	rVB	2368219	6347141	2.58%	1.238%
6	8.439	1061	1072	1078	rBV2	7164809	17855662	7.25%	3.484%
7	8.512	1078	1084	1089	rVV	6423391	14873428	6.04%	2.902%
8	8.579	1089	1095	1107	rVV	10509652	25171149	10.22%	4.911%
9	8.701	1107	1115	1126	rVB	10894891	19866637	8.06%	3.876%
10	9.159	1179	1190	1204	rVB	1066294	2798682	1.14%	0.546%
11	9.354	1204	1222	1237	rBV3	79967882	246411506	100.00%	48.074%
12	9.524	1244	1250	1255	rVB	2225798	3891298	1.58%	0.759%
13	9.616	1255	1265	1278	rVB2	5086189	12603819	5.11%	2.459%
14	9.756	1278	1288	1300	rBV2	5100360	10709243	4.35%	2.089%
15	9.902	1300	1312	1316	rBV	1759583	3536736	1.44%	0.690%
16	9.963	1316	1322	1334	rVB2	4145093	9588327	3.89%	1.871%
17	10.098	1334	1344	1358	rBV2	3651068	9566054	3.88%	1.866%
18	10.323	1374	1381	1385	rBV	6403057	12337761	5.01%	2.407%
19	10.390	1385	1392	1398	rVB2	5931114	13876615	5.63%	2.707%
20	10.506	1404	1411	1419	rVB3	2063449	4888553	1.98%	0.954%
21	10.671	1432	1438	1446	rVB	1815317	3443334	1.40%	0.672%
22	11.658	1590	1600	1606	rBV2	1834672	3364811	1.37%	0.656%
23	12.140	1671	1679	1686	rBV4	1208402	3075429	1.25%	0.600%
24	12.786	1780	1785	1793	rVB3	1661769	4076513	1.65%	0.795%
25	12.945	1802	1811	1816	rVV6	1601644	4628750	1.88%	0.903%
26	13.457	1888	1895	1899	rBV3	1426953	3073432	1.25%	0.600%
27	13.883	1960	1965	1970	rBV2	2554892	4436079	1.80%	0.865%
28	14.103	1997	2001	2008	rVB	2120479	3289327	1.33%	0.642%
29	14.213	2013	2019	2024	rVB5	1325596	2479241	1.01%	0.484%
30	14.396	2043	2049	2052	rBV2	1943022	3390118	1.38%	0.661%
31	14.804	2110	2116	2122	rBV3	3086892	5878650	2.39%	1.147%
32	15.340	2198	2204	2210	rVB2	1847044	3116552	1.26%	0.608%

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Instrument :
MSVOA_W
ClientSampleId :
CPSB-18(15-20)

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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Title : SW846 8260

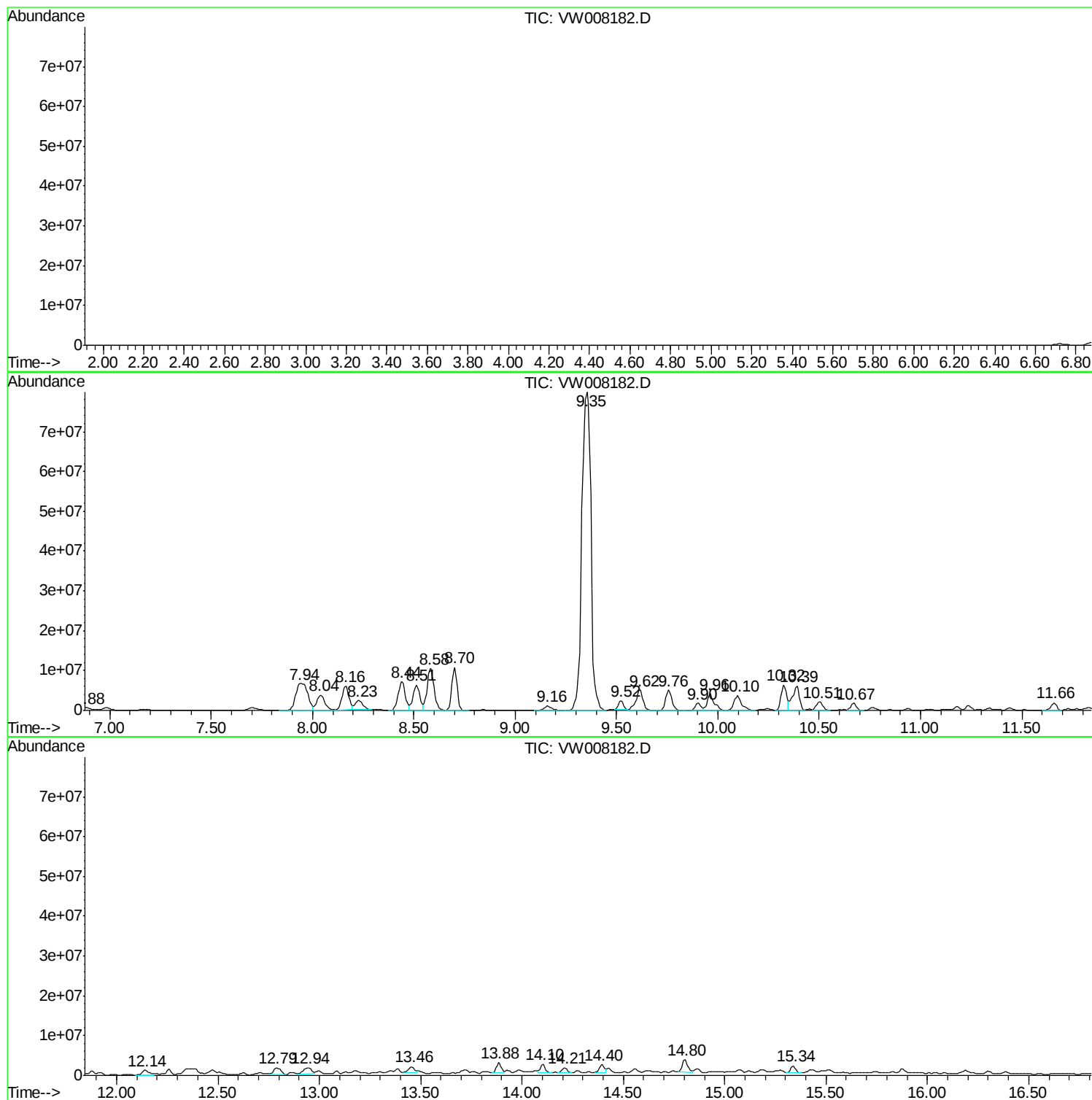
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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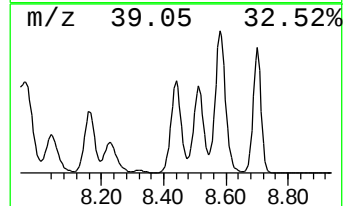
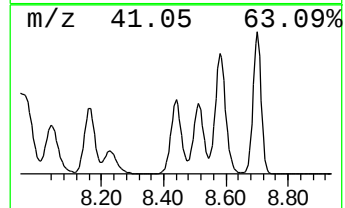
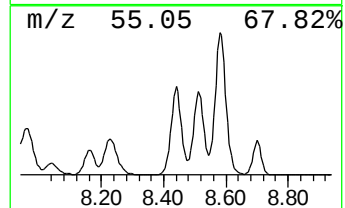
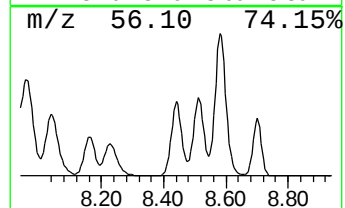
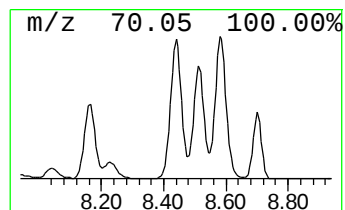
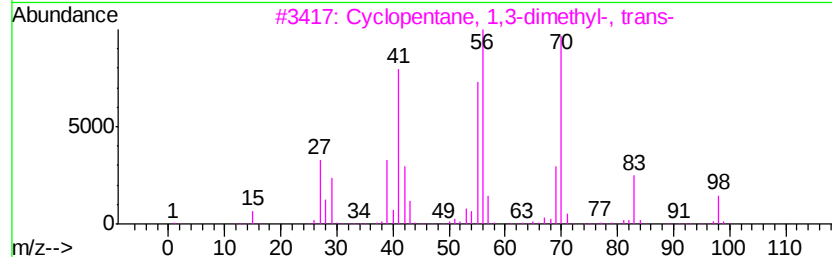
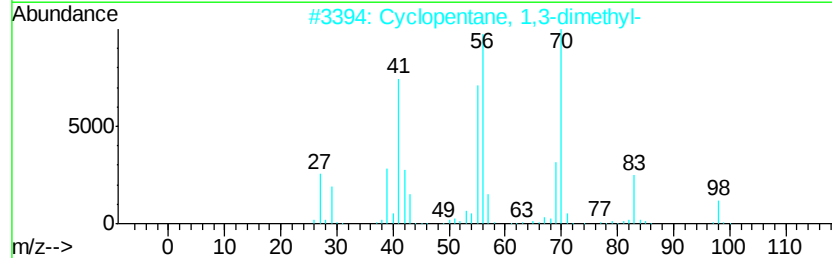
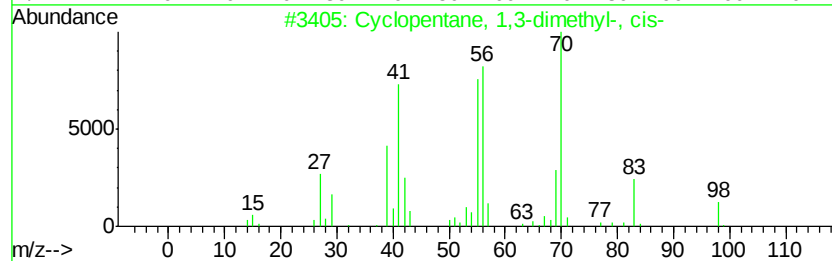
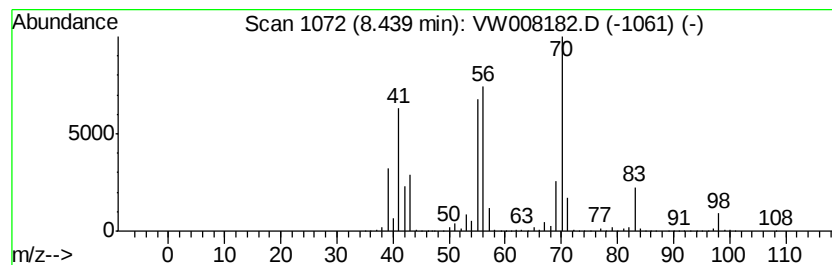
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.44	44.94 ug/l	17855700	1,4-Difluorobenzene	8.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	97
2		Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	95
3		Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	93
4		Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	76
5		Isopropylcyclobutane	98	C7H14	000872-56-0	53



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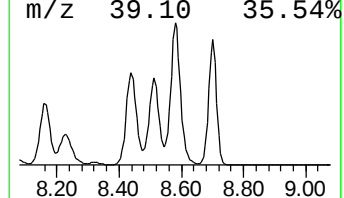
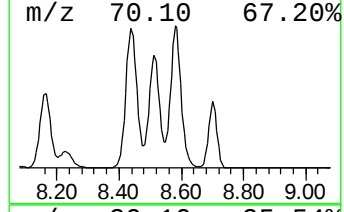
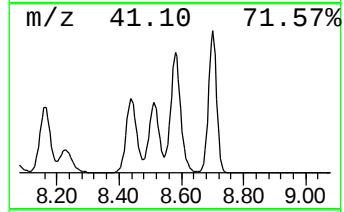
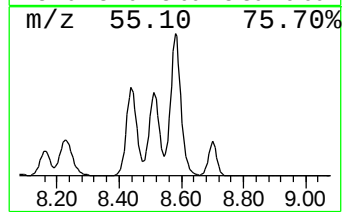
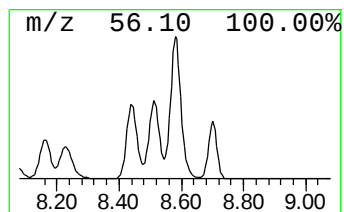
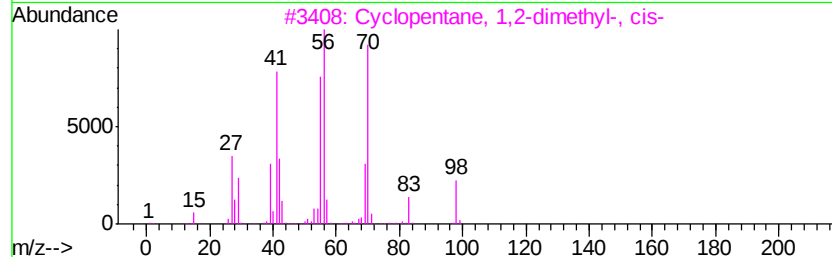
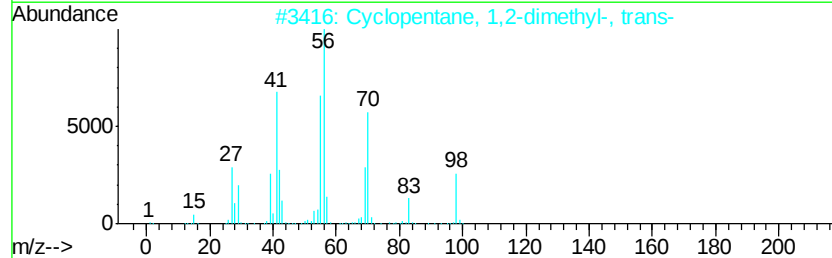
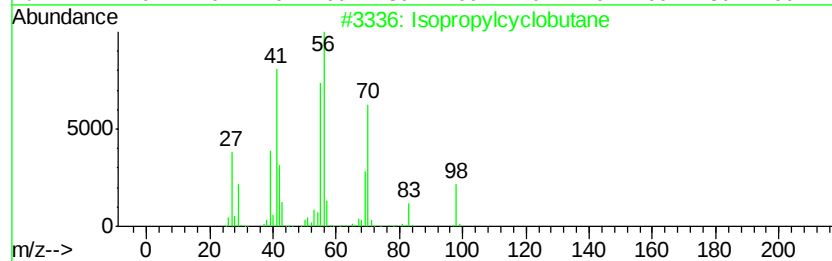
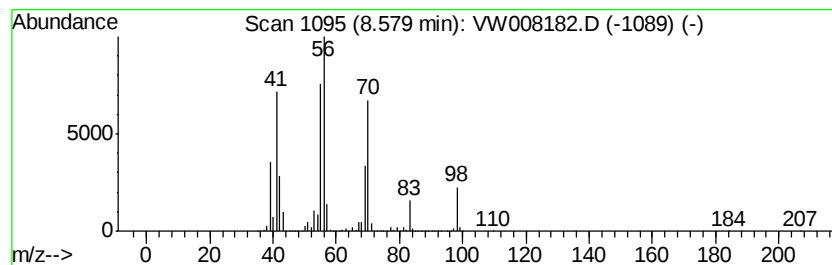
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.M
Quant Title : SW846 8260

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

Peak Number 2 Isopropylcyclobutane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.58	63.35 ug/l	25171100	1,4-Difluorobenzene	8.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Isopropylcvclobutane	98	C7H14	000872-56-0	96
2		Cyclopentane, 1,2-dimethyl-, trans-	98	C7H14	000822-50-4	95
3		Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	91
4		1-Heptene	98	C7H14	000592-76-7	86
5		Cyclopentane, 1,2-dimethyl-	98	C7H14	002452-99-5	86



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CPSB-18(15-20)

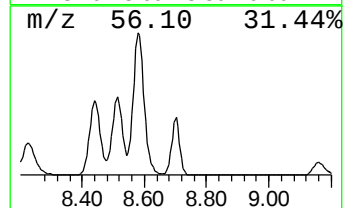
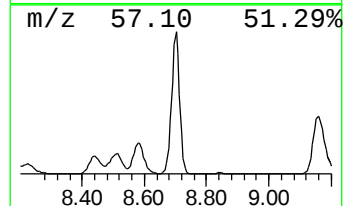
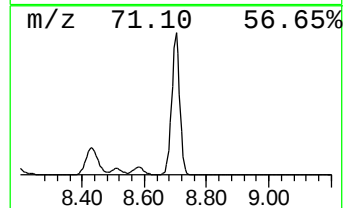
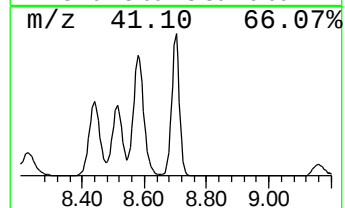
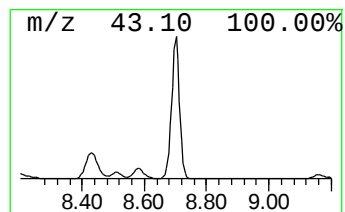
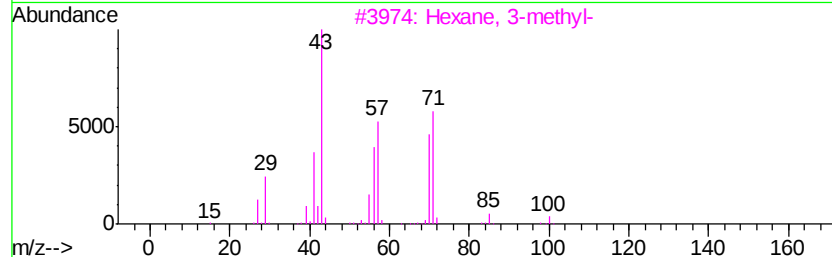
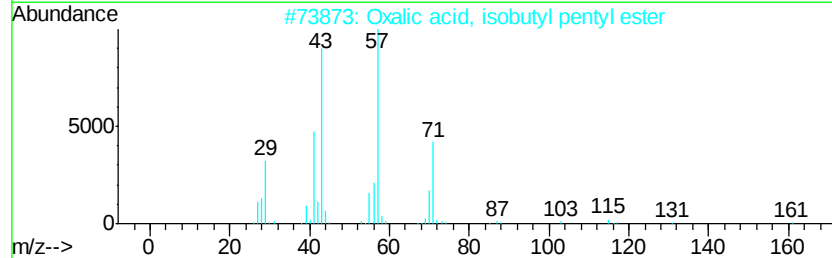
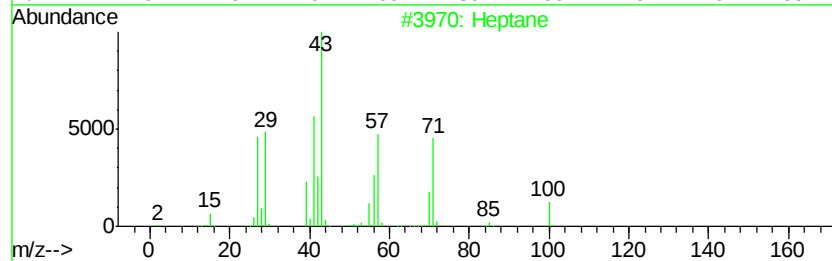
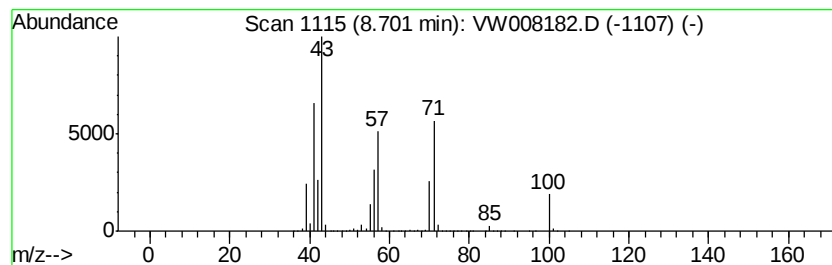
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Heptane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.70	50.00 ug/l	19866600	1,4-Difluorobenzene	8.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptane	100	C7H16	000142-82-5	91
2		Oxalic acid, isobutyl pentyl ester	216	C11H20O4	1000309-37-0	45
3		Hexane, 3-methyl-	100	C7H16	000589-34-4	40
4		Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	40
5		Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	38



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 Misc : 5.09G/5ML/MSVOA W/SOIL
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
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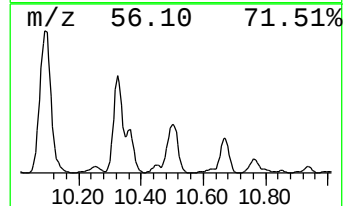
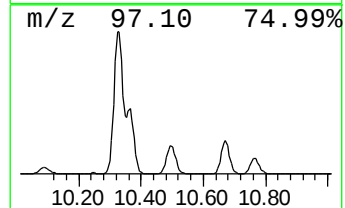
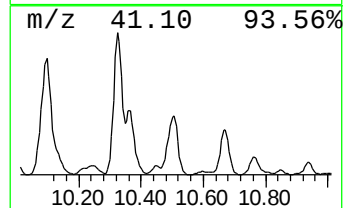
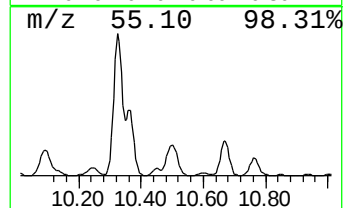
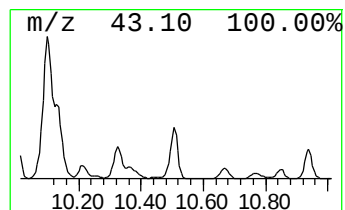
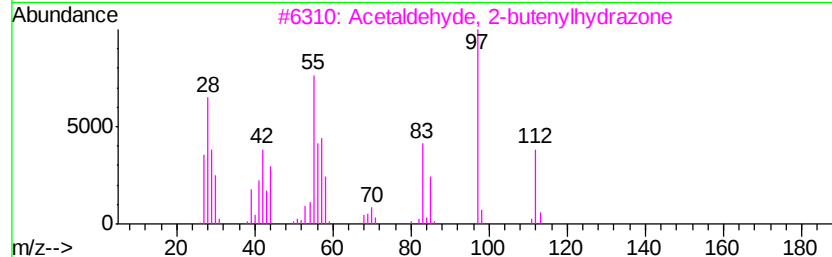
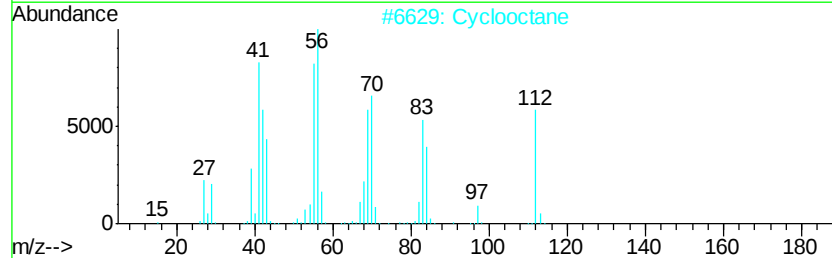
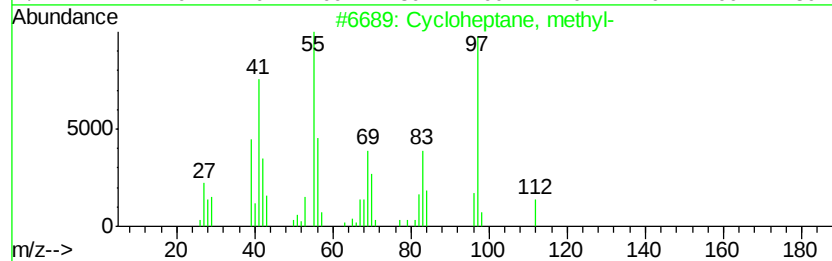
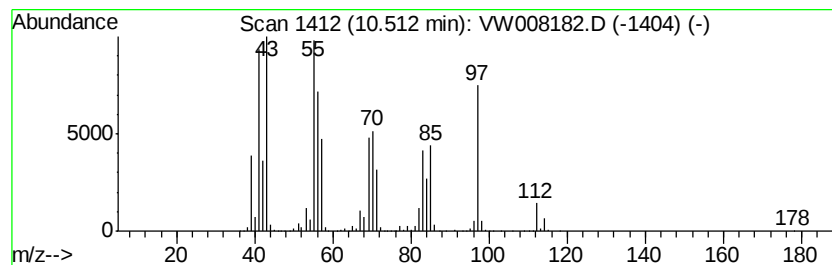
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 Quant Title : SW846 8260

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

 Peak Number 4 Cycloheptane, methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.51	72.64 ug/l	4888550	Chlorobenzene-d5	11.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cycloheptane, methyl-	112	C8H16	004126-78-7	76
2		Cyclooctane	112	C8H16	000292-64-8	50
3		Acetaldehyde, 2-butenylhydrazone	112	C6H12N2	075268-07-4	49
4		Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	46
5		3-Octene, (Z)-	112	C8H16	014850-22-7	45



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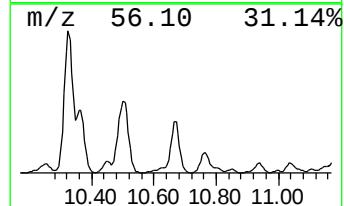
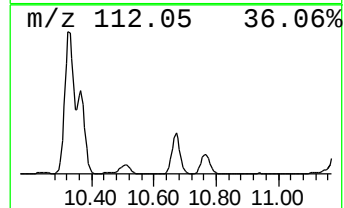
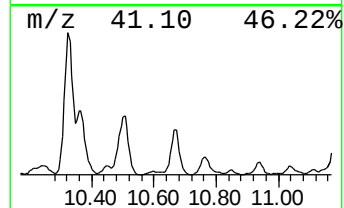
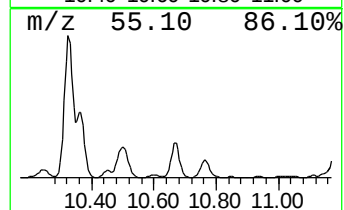
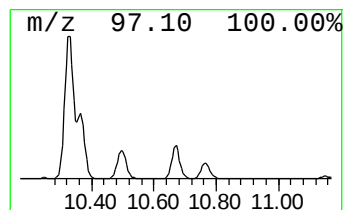
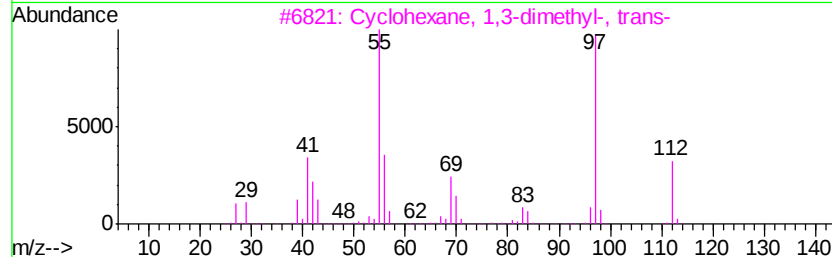
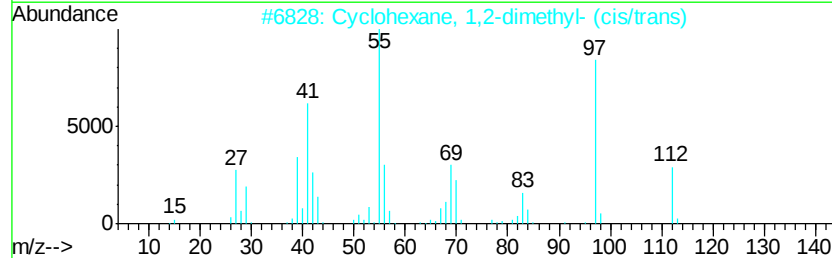
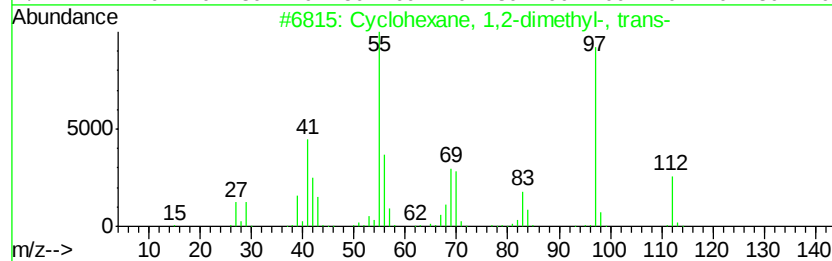
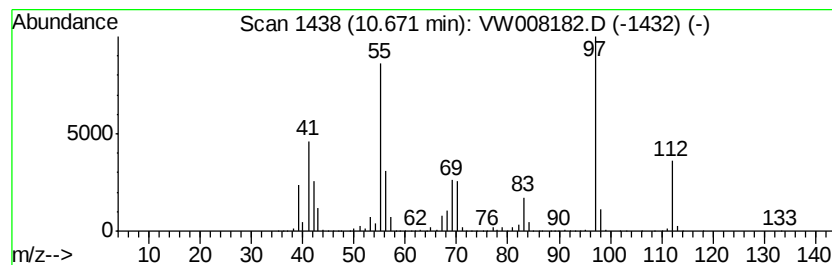
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Quant Title : SW846 8260

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

Peak Number 5 Cyclohexane, 1,2-dimethyl-,... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.67	51.17 ug/l	3443330	Chlorobenzene-d5	11.63

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, 1,2-dimethyl-, trans-	112	C8H16	006876-23-9	94
2		Cyclohexane, 1,2-dimethyl- (cis/...	112	C8H16	000583-57-3	90
3		Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	87
4		Cyclohexane, 1,4-dimethyl-, cis-	112	C8H16	000624-29-3	74
5		Cyclohexane, 1,4-dimethyl-	112	C8H16	000589-90-2	74



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW011019\
Data File : VW008182.D
Acq On : 11 Jan 2019 07:46
Operator : SY/AP
Sample : K1102-01
Misc : 5.09G/5ML/MSVOA W/SOIL
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
CPSB-18(15-20)

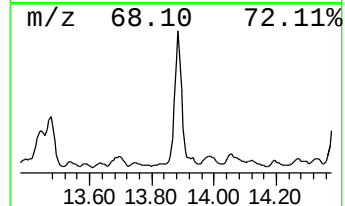
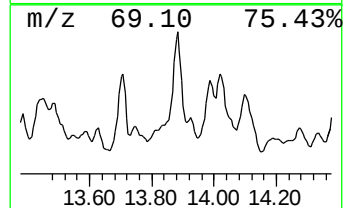
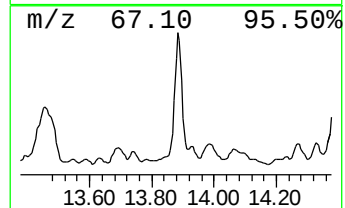
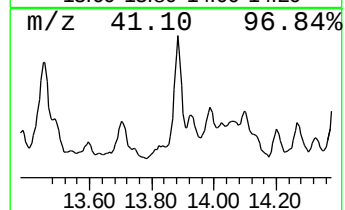
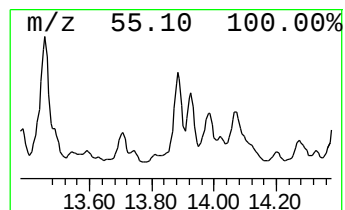
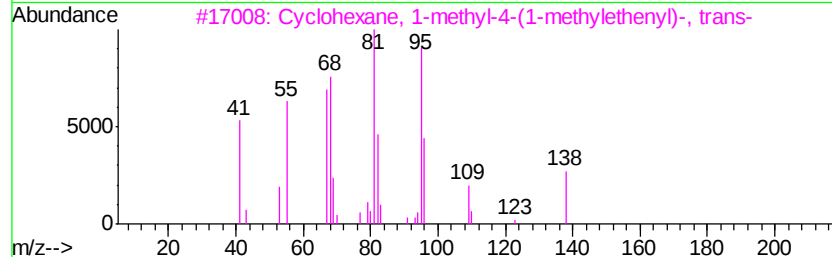
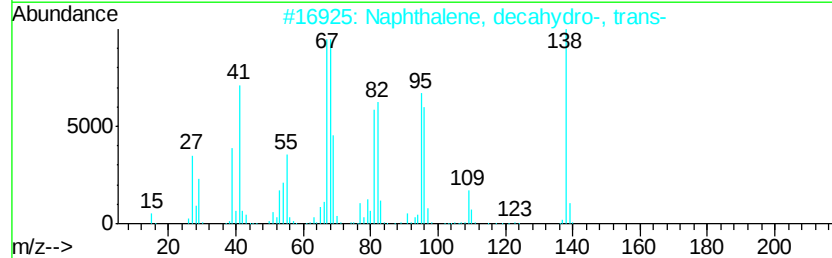
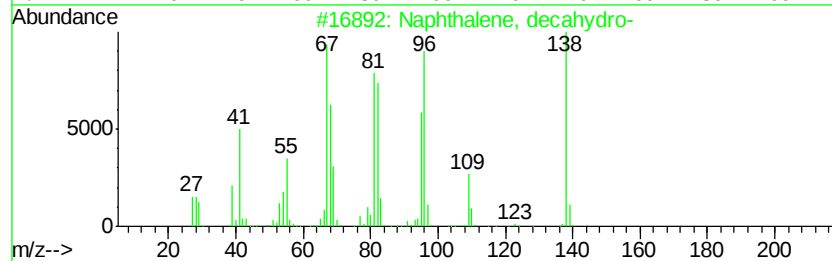
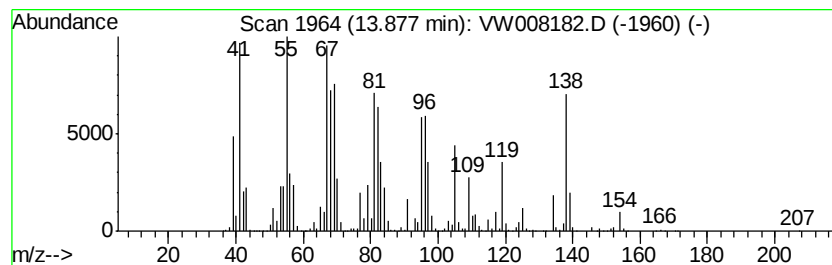
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Quant Title : SW846 8260

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

Peak Number 6 Naphthalene, decahydro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.88	72.17 ug/l	4436080	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, decahydro-	138	C10H18	000091-17-8	96
2		Naphthalene, decahydro-, trans-	138	C10H18	000493-02-7	95
3		Cyclohexane, 1-methyl-4-(1-methyl-...	138	C10H18	001124-25-0	91
4		Naphthalene, decahydro-, cis-	138	C10H18	000493-01-6	86
5		Cyclohexane, 1-methyl-3-(1-methyl-...	138	C10H18	024399-15-3	70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW011019\
Data File : VW008182.D
Acq On : 11 Jan 2019 07:46
Operator : SY/AP
Sample : K1102-01
Misc : 5.09G/5ML/MSVOA W/SOIL
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
CPSB-18(15-20)

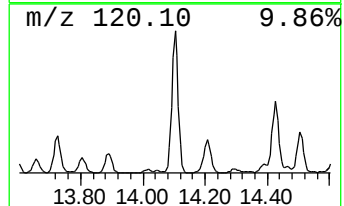
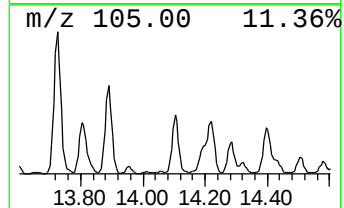
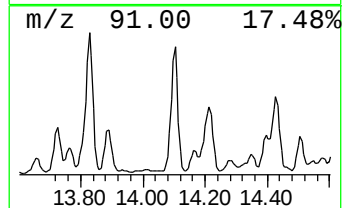
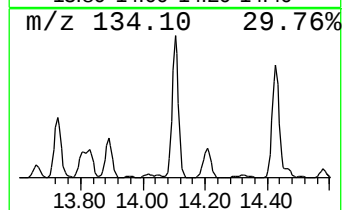
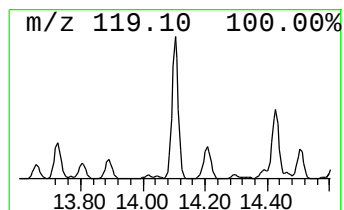
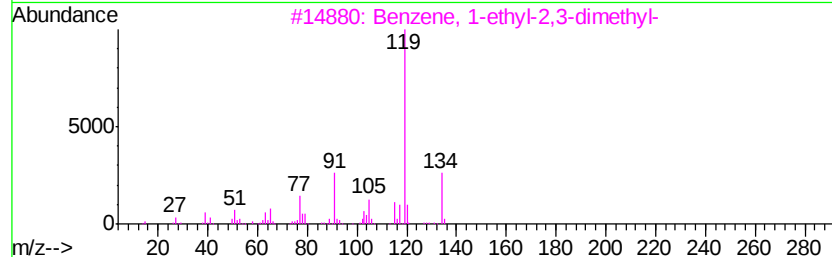
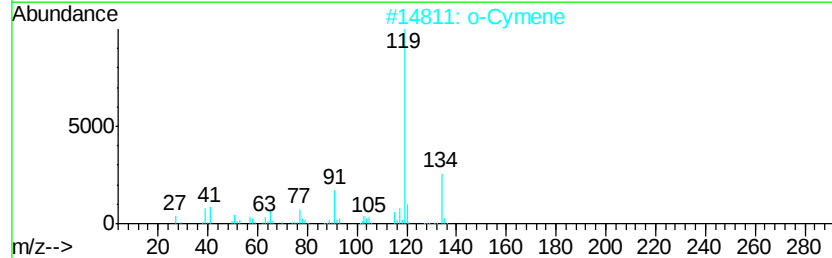
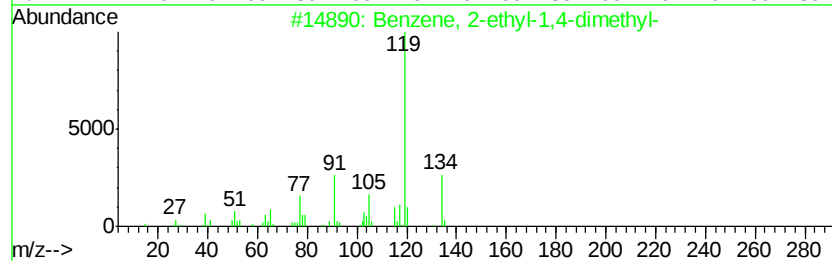
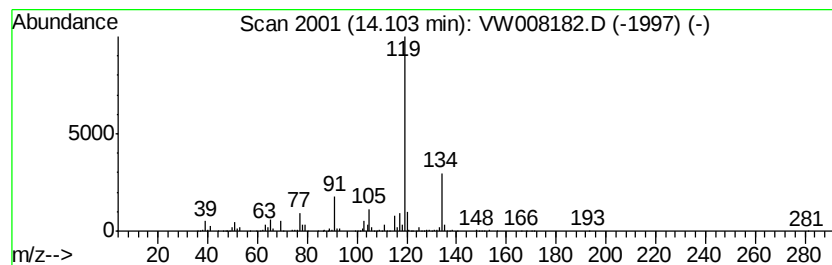
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.10	53.51 ug/l	3289330	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2		o-Cymene	134	C10H14	000527-84-4	95
3		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW011019\
Data File : VW008182.D
Acq On : 11 Jan 2019 07:46
Operator : SY/AP
Sample : K1102-01
Misc : 5.09G/5ML/MSVOA W/SOIL
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
CPSB-18(15-20)

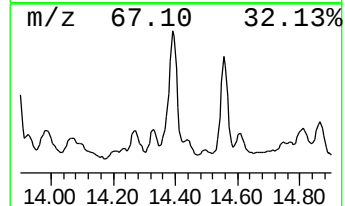
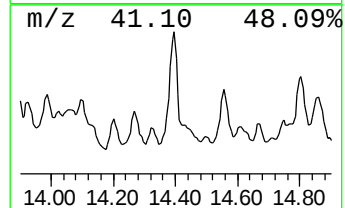
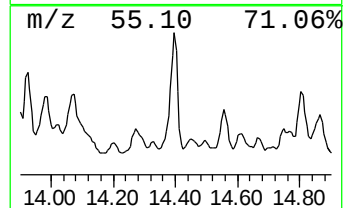
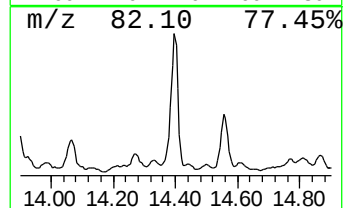
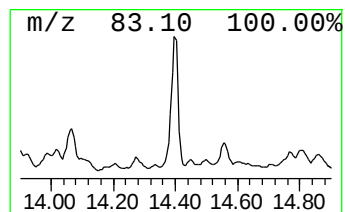
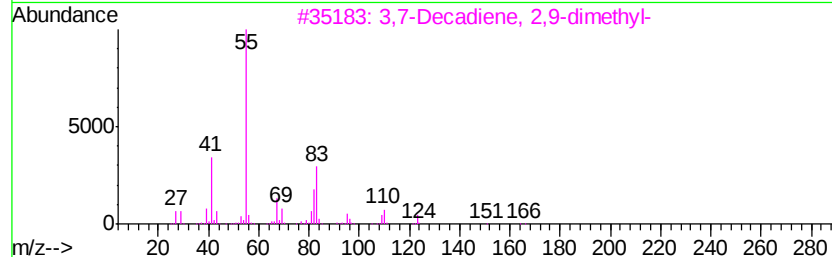
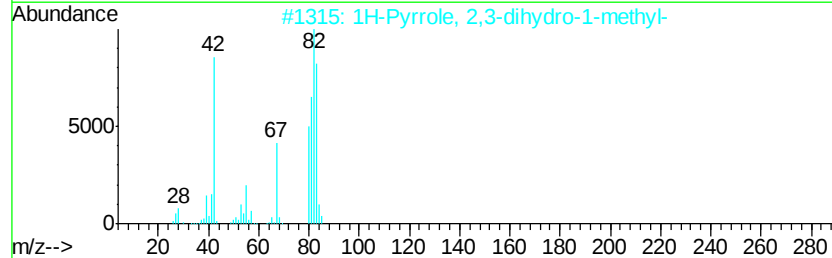
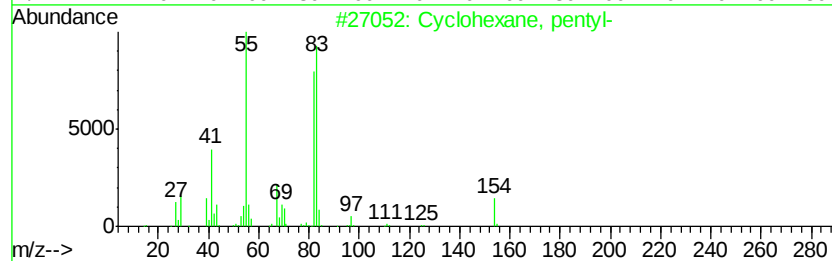
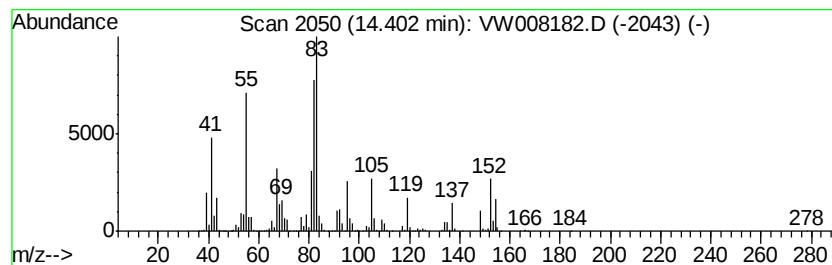
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Quant Title : SW846 8260

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

Peak Number 8 Cyclohexane, pentyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.40	55.15 ug/l	3390120	1,4-Dichlorobenzene-d4	13.56

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, pentyl-	154	C11H22	004292-92-6	55
2		1H-Pyrrole, 2,3-dihydro-1-methyl-	83	C5H9N	033838-11-8	49
3		3,7-Decadiene, 2,9-dimethyl-	166	C12H22	074630-13-0	47
4		Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	43
5		Cyclohexane, 2-propenyl-	124	C9H16	002114-42-3	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW011019\
Data File : VW008182.D
Acq On : 11 Jan 2019 07:46
Operator : SY/AP
Sample : K1102-01
Misc : 5.09G/5ML/MSVOA W/SOIL
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
CPSB-18(15-20)

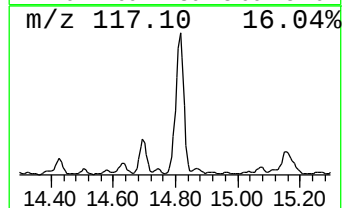
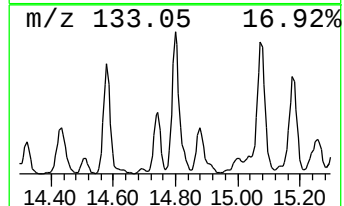
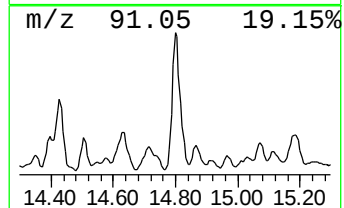
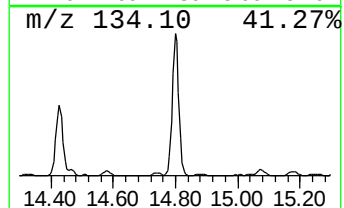
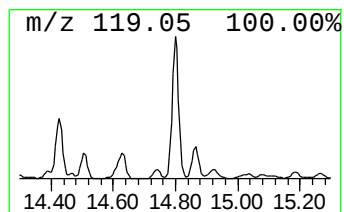
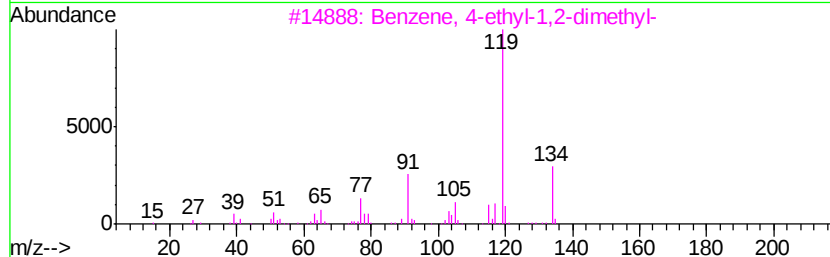
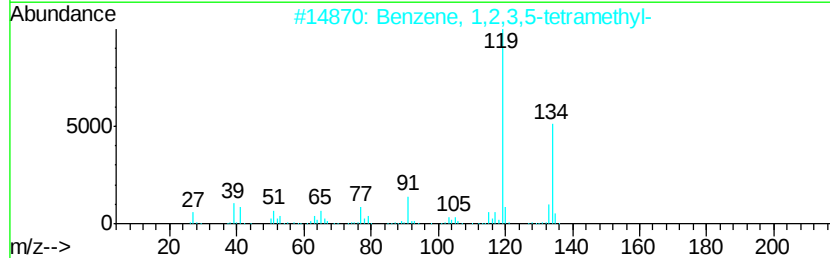
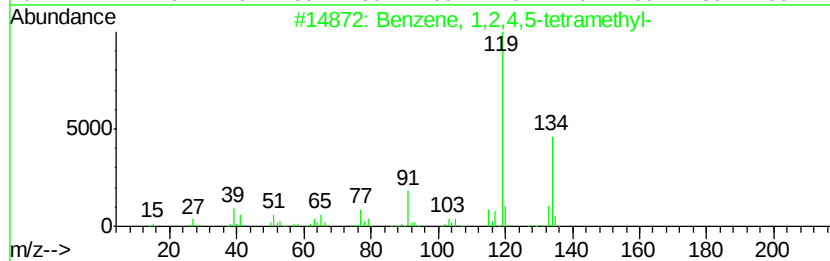
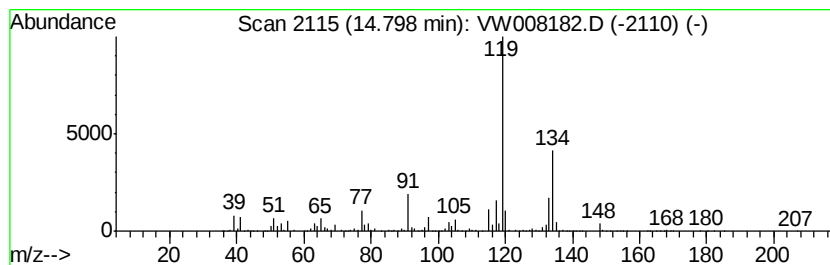
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.80	95.64 ug/l	5878650	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	93
2		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	81
3		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	76
4		1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	76
5		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	76



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW011019\
Data File : VW008182.D
Acq On : 11 Jan 2019 07:46
Operator : SY/AP
Sample : K1102-01
Misc : 5.09G/5ML/MSVOA W/SOIL
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
CPSB-18(15-20)

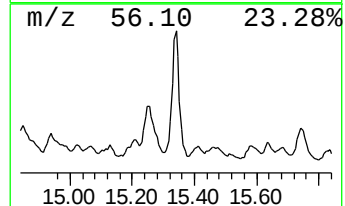
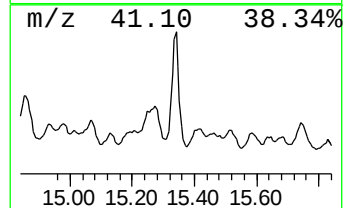
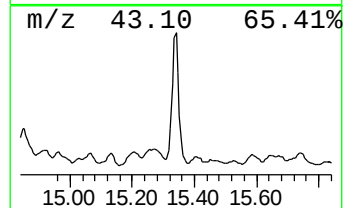
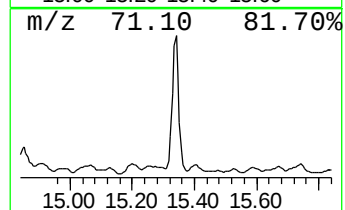
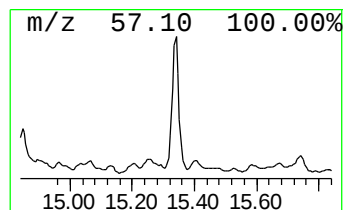
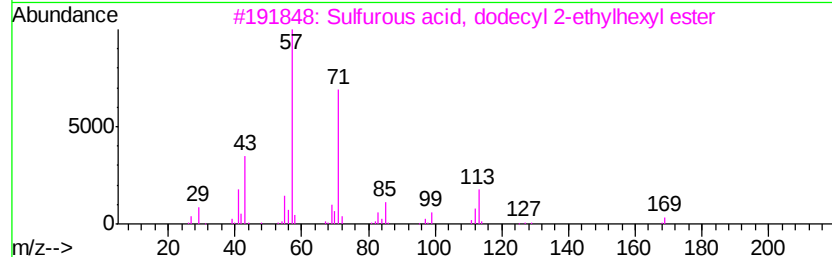
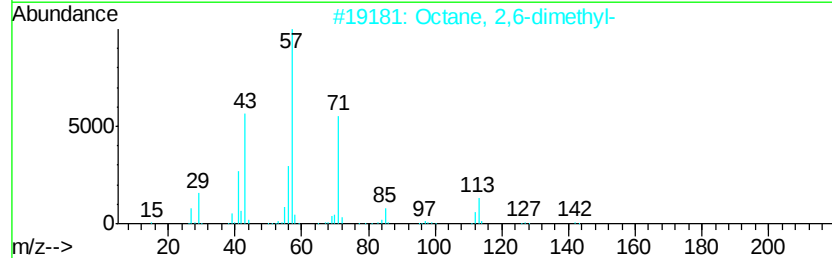
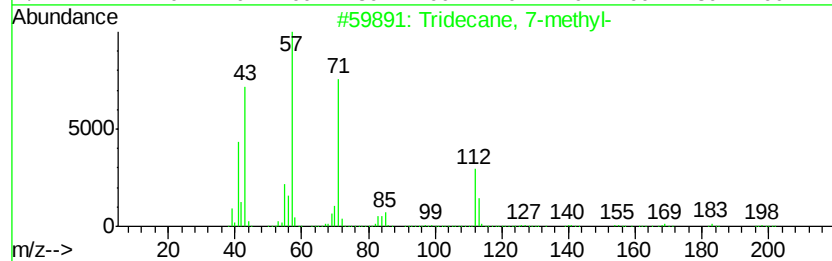
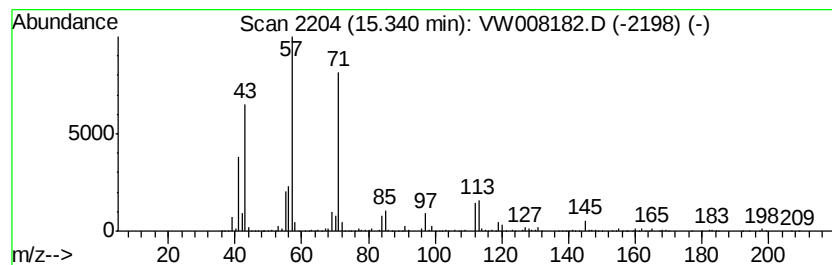
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.M
Quant Title : SW846 8260

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

Peak Number 10 Tridecane, 7-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.34	50.70 ug/l	3116550	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 7-methyl-	198	C14H30	026730-14-3	81
2		Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	72
3		Sulfurous acid, dodecyl 2-ethylh...	362	C20H42O3S	1000309-19-5	64
4		Nonane, 3-methyl-	142	C10H22	005911-04-6	59
5		Nonane, 4-methyl-5-propyl-	184	C13H28	062185-55-1	58



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_W\DATA\VW011019\
Data File : VW008182.D
Acq On : 11 Jan 2019 07:46
Operator : SY/AP
Sample : K1102-01
Misc : 5.09G/5ML/MSVOA_W/SOIL
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
CPSB-18(15-20)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.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, 1,3...	8.44	44.9	ug/l	17855700	2	8.84	19866600	50.0
Isopropylcyclobutane	8.58	63.4	ug/l	25171100	2	8.84	19866600	50.0
Heptane	8.70	50.0	ug/l	19866600	2	8.84	19866600	50.0
Cycloheptane, met...	10.51	72.6	ug/l	4888550	3	11.63	3364810	50.0
Cyclohexane, 1,2-...	10.67	51.2	ug/l	3443330	3	11.63	3364810	50.0
Naphthalene, deca...	13.88	72.2	ug/l	4436080	4	13.56	3073430	50.0
Benzene, 2-ethyl-...	14.10	53.5	ug/l	3289330	4	13.56	3073430	50.0
Cyclohexane, pentyl-	14.40	55.1	ug/l	3390120	4	13.56	3073430	50.0
Benzene, 1,2,4,5-...	14.80	95.6	ug/l	5878650	4	13.56	3073430	50.0
Tridecane, 7-methyl-	15.34	50.7	ug/l	3116550	4	13.56	3073430	50.0