

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110222\
 Data File : VY011264.D
 Acq On : 02 Nov 2022 15:49
 Operator : KP/MD
 Sample : N5395-08
 Misc : 5.12g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 S-2-103122

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y110122S.M

Title : SW846 8260

Signal : TIC: VY011264.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.686	136	142	148	rVB	11405	21618	2.35%	0.506%
2	3.942	339	348	357	rBV	15787	42799	4.65%	1.003%
3	4.704	464	473	489	rBV4	8330	28354	3.08%	0.664%
4	6.795	806	816	828	rBV2	3718	11955	1.30%	0.280%
5	6.984	837	847	861	rBV3	4969	16002	1.74%	0.375%
6	7.716	956	967	973	rBV	162121	394971	42.89%	9.253%
7	7.789	973	979	994	rVB	246841	554274	60.18%	12.985%
8	8.136	1026	1036	1047	rBV	126034	268168	29.12%	6.282%
9	8.685	1116	1126	1138	rBV	348263	671808	72.95%	15.738%
10	10.179	1361	1371	1378	rBV	544175	920975	100.00%	21.575%
11	11.489	1578	1586	1597	rBV	290089	472969	51.36%	11.080%
12	11.989	1662	1668	1675	rBV9	3237	9826	1.07%	0.230%
13	12.111	1677	1688	1695	rBV5	7428	25169	2.73%	0.590%
14	12.233	1705	1708	1714	rVB3	6618	10020	1.09%	0.235%
15	12.477	1741	1748	1755	rBV2	188046	324326	35.22%	7.598%
16	12.642	1768	1775	1782	rVB5	11068	23840	2.59%	0.558%
17	12.727	1782	1789	1795	rBV8	4043	12644	1.37%	0.296%
18	12.812	1795	1803	1810	rBV10	8424	25939	2.82%	0.608%
19	12.940	1816	1824	1828	rVB7	4772	11313	1.23%	0.265%
20	12.995	1828	1833	1837	rBV3	7939	13649	1.48%	0.320%
21	13.068	1841	1845	1850	rVB5	7077	11825	1.28%	0.277%
22	13.166	1858	1861	1867	rBV7	5557	12222	1.33%	0.286%
23	13.282	1876	1880	1884	rBV4	8712	11085	1.20%	0.260%
24	13.422	1894	1903	1908	rBV	112862	181472	19.70%	4.251%
25	13.477	1908	1912	1917	rVB3	7661	14764	1.60%	0.346%
26	13.745	1951	1956	1960	rBV6	10935	21857	2.37%	0.512%
27	13.959	1986	1991	2004	rVB2	20440	45390	4.93%	1.063%
28	14.135	2015	2020	2027	rVB9	5661	13378	1.45%	0.313%
29	14.196	2027	2030	2035	rBV7	5593	10151	1.10%	0.238%
30	14.251	2035	2039	2046	rBV8	9004	20104	2.18%	0.471%
31	14.410	2056	2065	2070	rBV9	12525	33686	3.66%	0.789%
32	14.733	2114	2118	2124	rVB7	6374	11628	1.26%	0.272%
33	15.214	2194	2197	2202	rVB6	6547	9669	1.05%	0.227%
34	15.592	2257	2259	2266	rVB8	5683	10809	1.17%	0.253%

Sum of corrected areas: 4268659

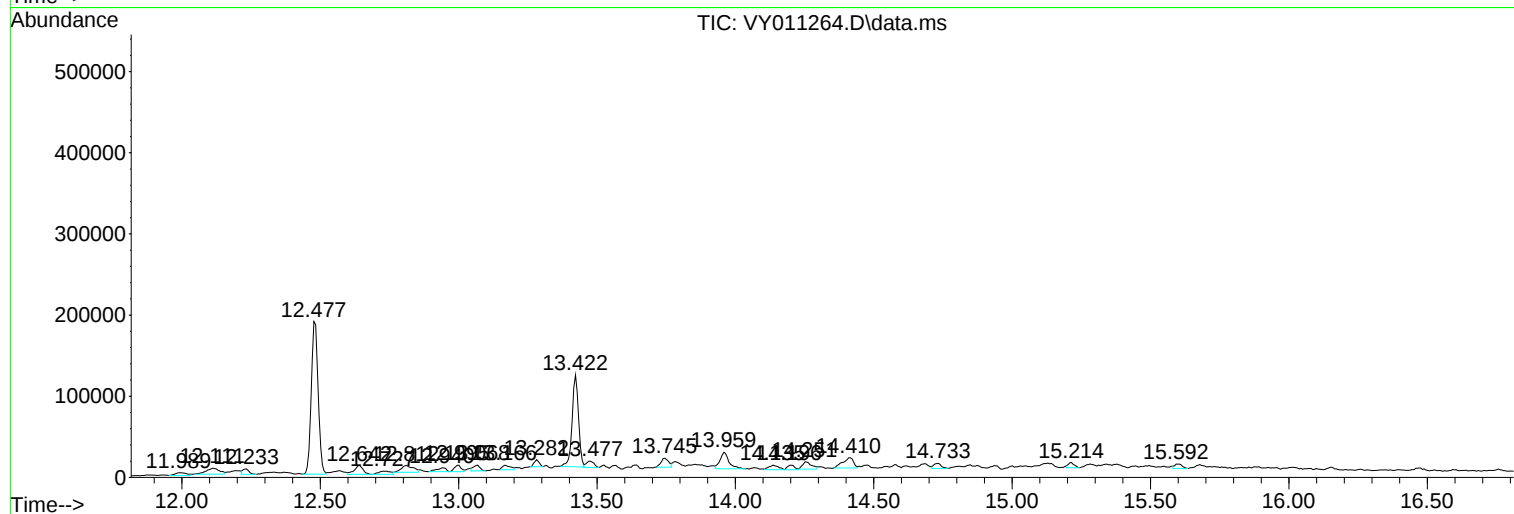
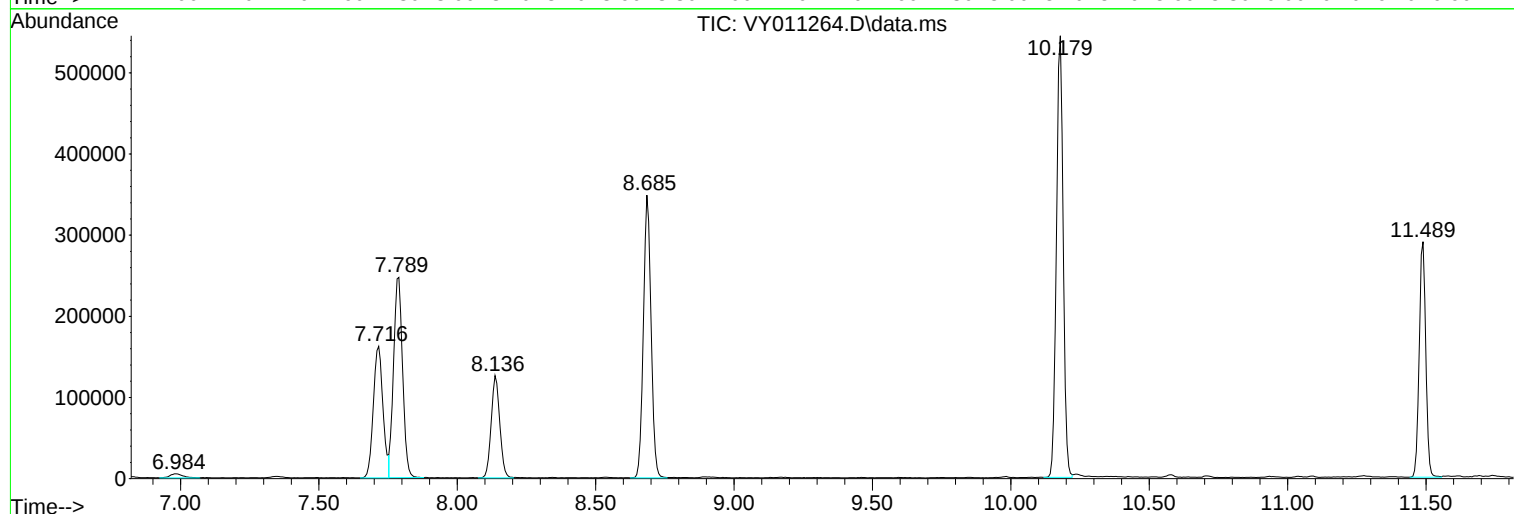
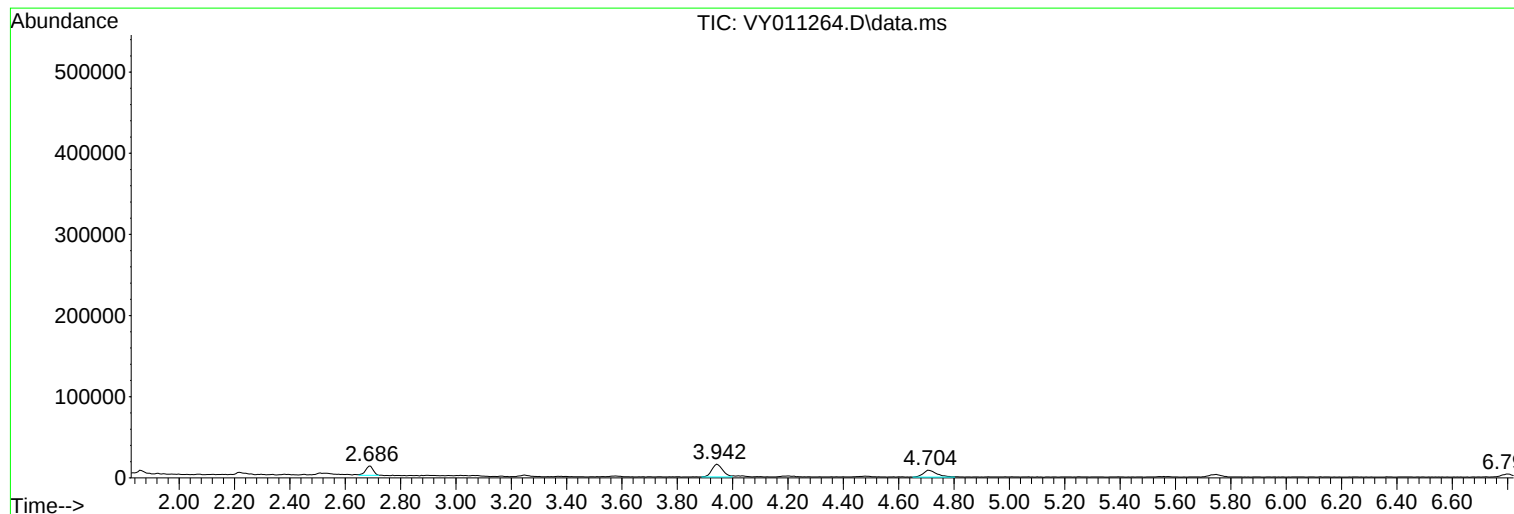
Instrument :
MSVOA_Y
ClientSampleId :
S-2-103122

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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y110122S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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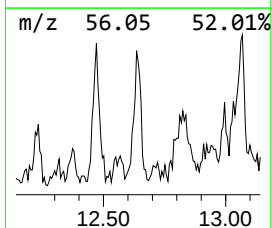
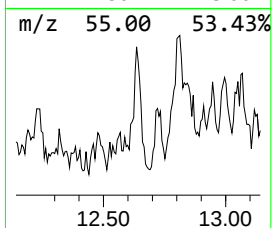
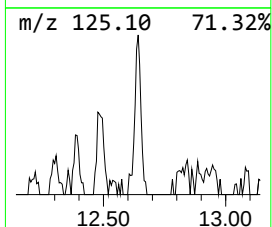
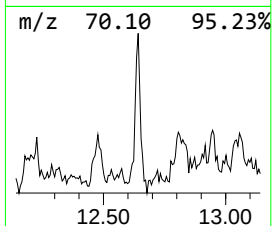
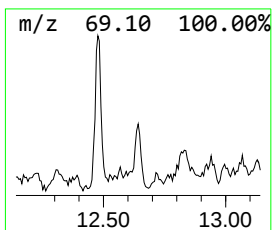
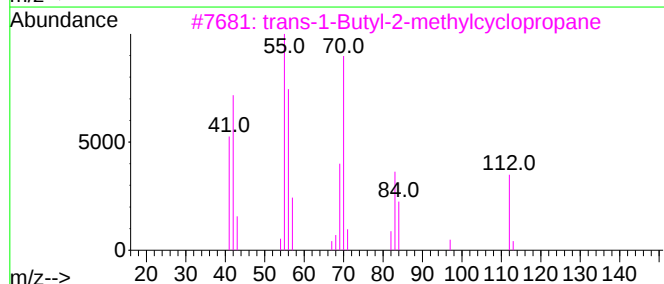
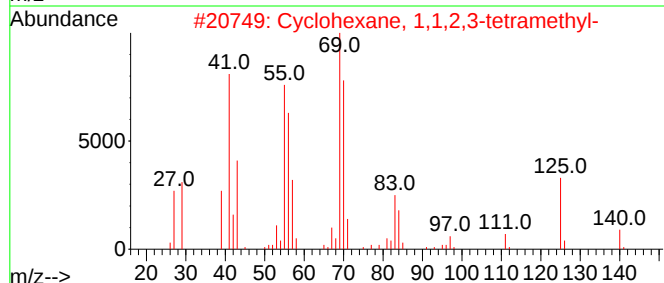
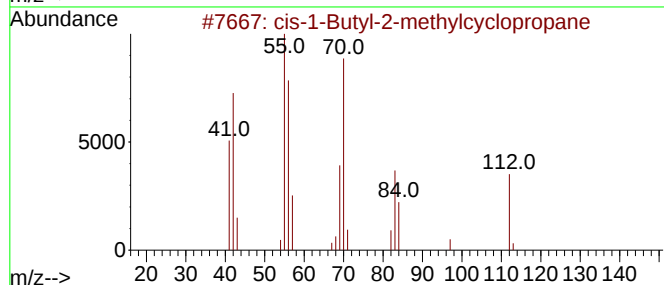
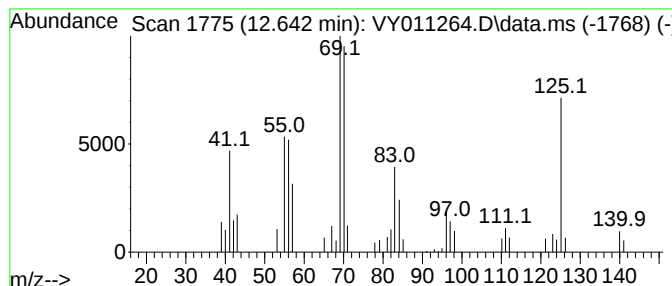
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y110122S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 cis-1-Butyl-2-methylcyclopr... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.642	6.57 ug/l	23840	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			cis-1-Butyl-2-methylcyclopropane	112	C8H16	038851-69-3	55
2			Cyclohexane, 1,1,2,3-tetramethyl-	140	C10H20	006783-92-2	52
3			trans-1-Butyl-2-methylcyclopropane	112	C8H16	038851-70-6	49
4			2-Octene, 2,6-dimethyl-	140	C10H20	004057-42-5	46
5			Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	38



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ClientSampleId :
S-2-103122

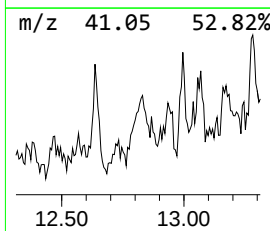
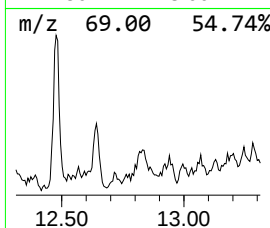
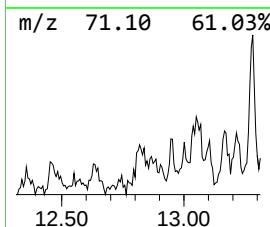
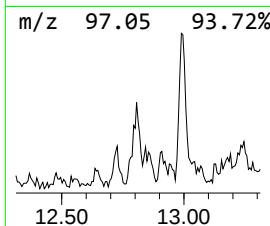
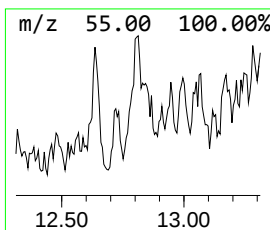
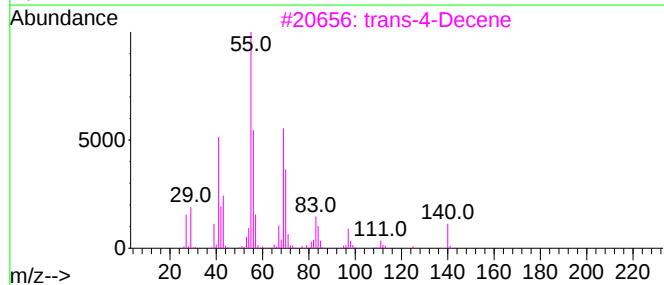
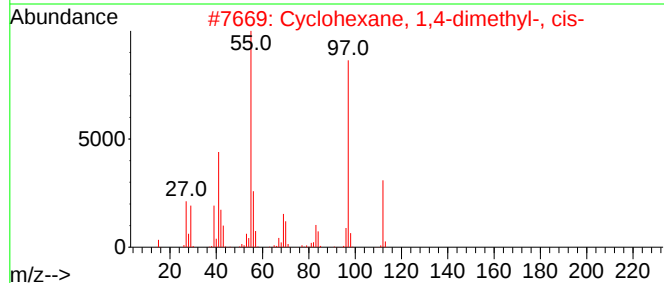
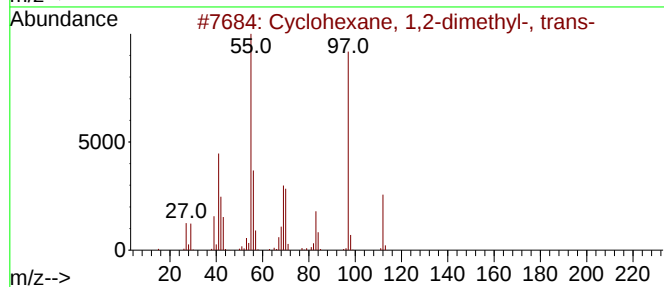
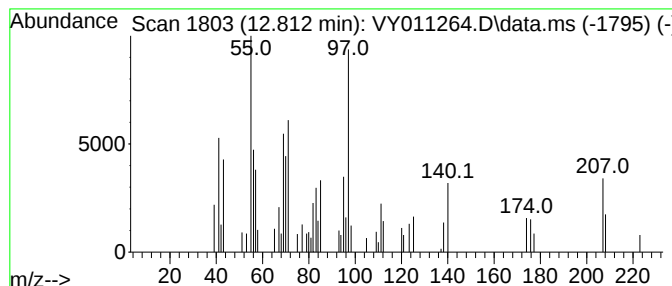
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y110122S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Cyclohexane, 1,2-dimethyl-,... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.812	7.15 ug/l	25939	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,2-dimethyl-, trans-	112	C8H16	006876-23-9	50
2			Cyclohexane, 1,4-dimethyl-, cis-	112	C8H16	000624-29-3	25
3			trans-4-Decene	140	C10H20	019398-89-1	25
4			1-(2-Furyl)-3-buten-1-ol	138	C8H10O2	1000311-94-9	25
5			2-Cyclohexen-1-ol, 1-methyl-	112	C7H12O	023758-27-2	25



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ClientSampleId :
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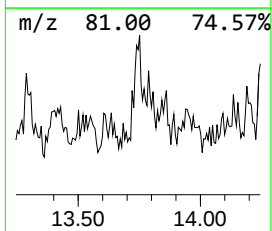
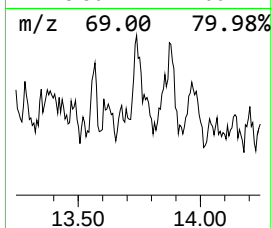
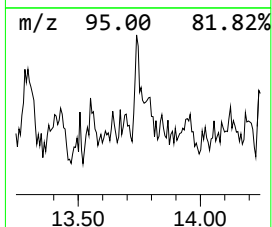
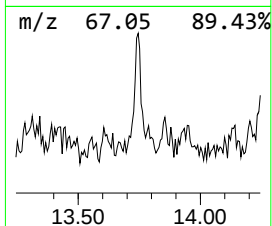
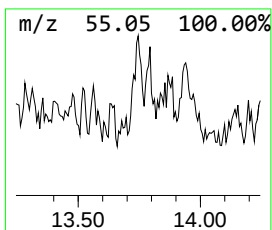
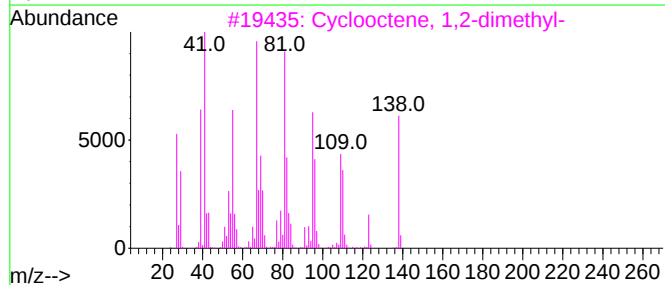
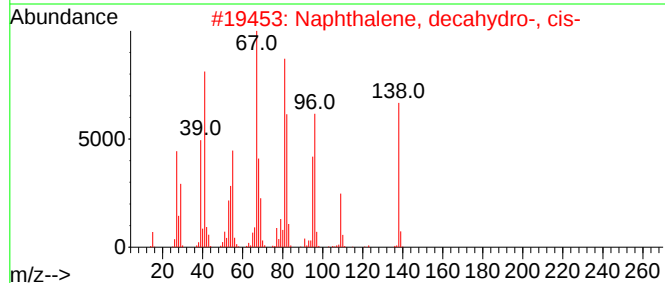
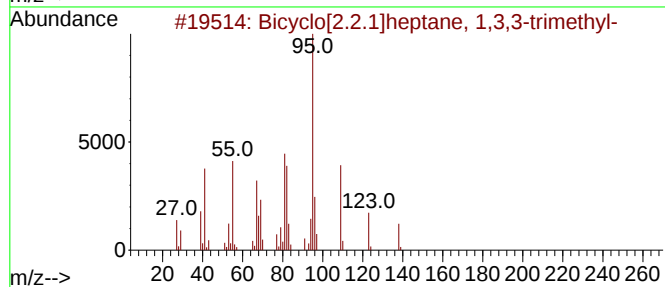
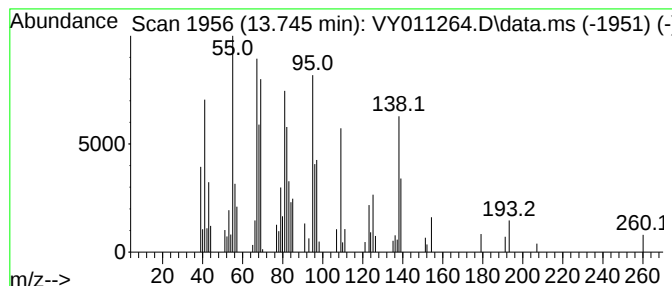
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y110122S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Bicyclo[2.2.1]heptane, 1,3,... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.745	6.02 ug/l	21857	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[2.2.1]heptane, 1,3,3-tri...	138	C10H18	006248-88-0	58
2			Naphthalene, decahydro-, cis-	138	C10H18	000493-01-6	50
3			Cyclooctene, 1,2-dimethyl-	138	C10H18	054299-96-6	46
4			Cyclohexene, 3-methyl-6-(1-methy...	138	C10H18	005256-65-5	46
5			2,6-Dimethylbicyclo[3.2.1]octane	138	C10H18	1000215-28-2	38



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ClientSampleId :
S-2-103122

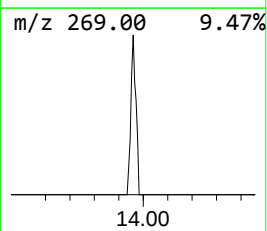
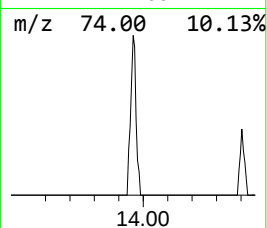
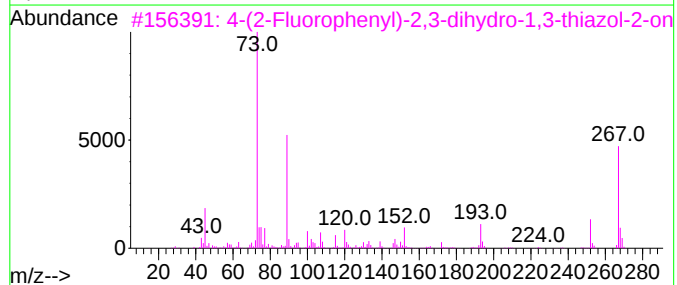
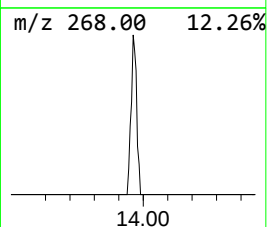
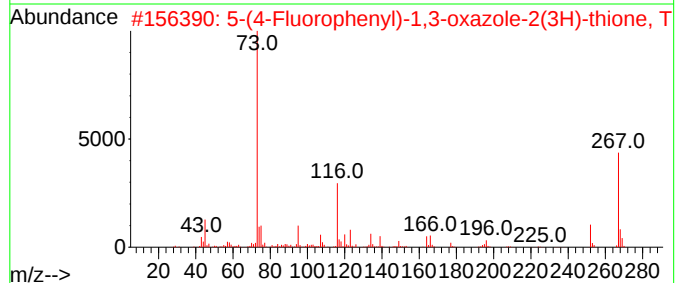
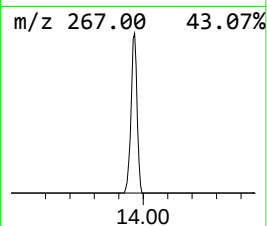
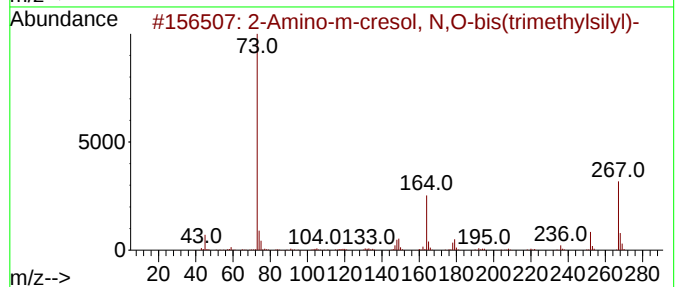
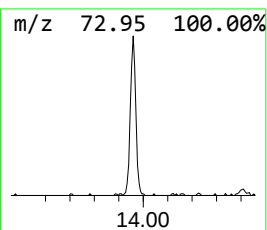
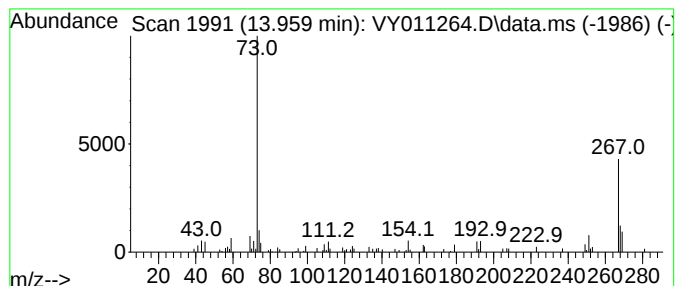
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y110122S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 2-Amino-m-cresol, N,O-bis(t... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.959	12.51 ug/l	45390	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Amino-m-cresol, N,O-bis(trimet...	267	C13H25NOSi2	1000449-77-1	53
2			5-(4-Fluorophenyl)-1,3-oxazole-2...	267	C12H14FNOSSi	1000497-07-6	50
3			4-(2-Fluorophenyl)-2,3-dihydro-1...	267	C12H14FNOSSi	1000504-37-6	38
4			Benzeneethanamine, N-[(pentafluo...	475	C21H26F5NO2Si2	055429-85-1	36
5			4-(Methylamino)phenol, N,O-bis(t...	267	C13H25NOSi2	1000471-13-0	33



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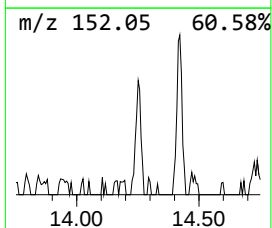
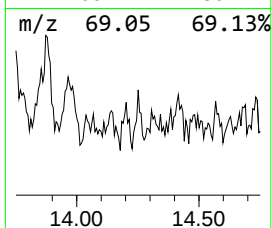
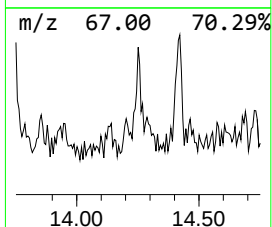
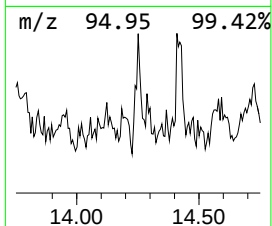
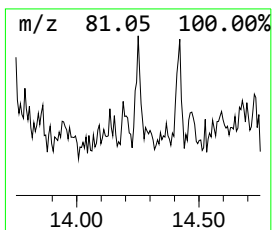
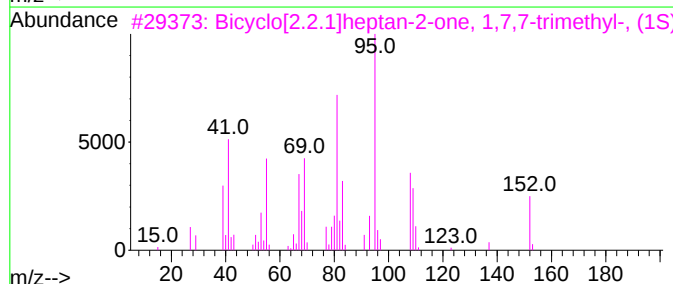
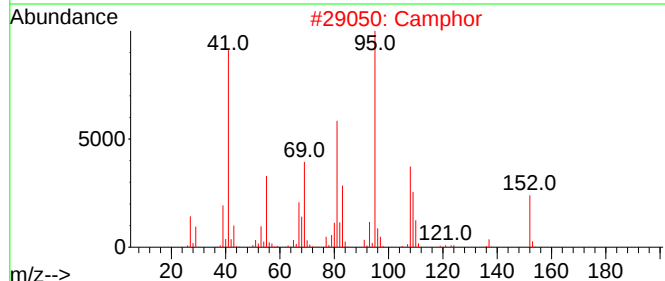
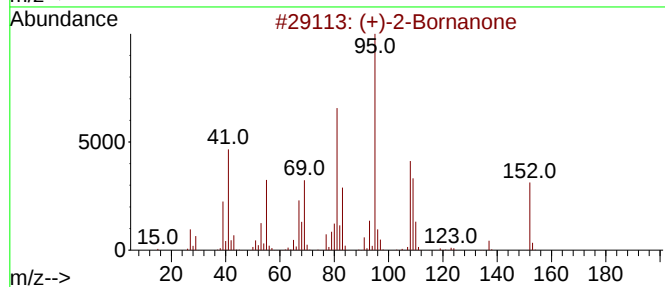
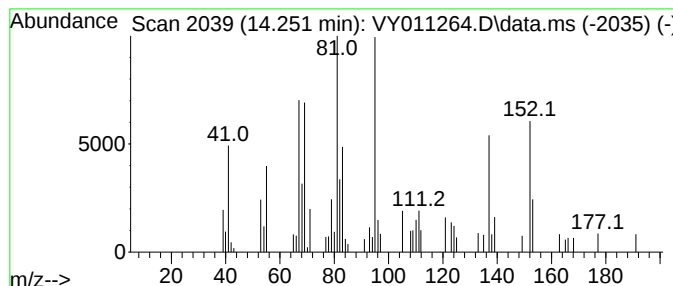
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y110122S.M
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TIC Integration Parameters: LSCINT.P

Peak Number 5 (+)-2-Bornanone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.251	5.54 ug/l	20104	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(+)-2-Bornanone	152	C10H16O	000464-49-3	76
2			Camphor	152	C10H16O	000076-22-2	76
3			Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	68
4			trans,trans-3-Ethylbicyclo[4.4.0]...	166	C12H22	058462-32-1	60
5			Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110222\
Data File : VY011264.D
Acq On : 02 Nov 2022 15:49
Operator : KP/MD
Sample : N5395-08
Misc : 5.12g/5.0mL/MSVOA_Y/soil
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
S-2-103122

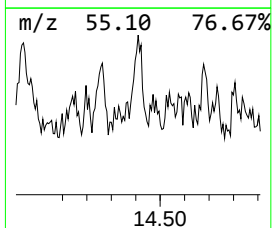
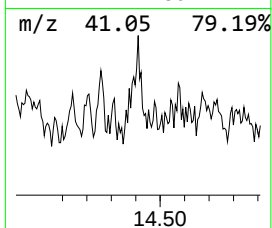
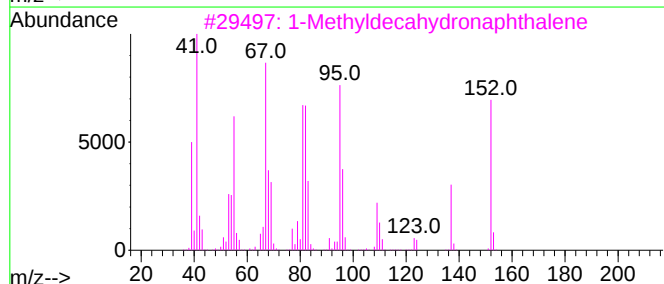
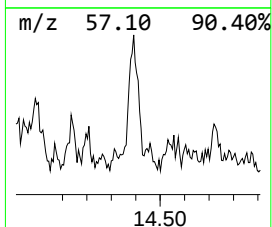
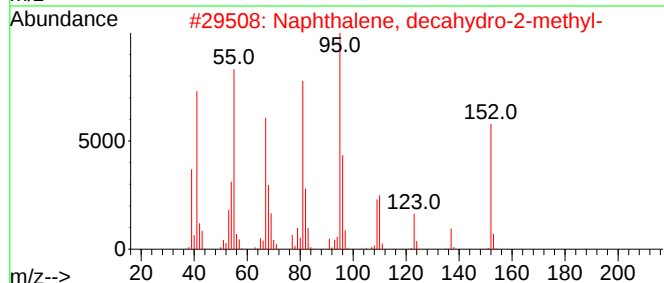
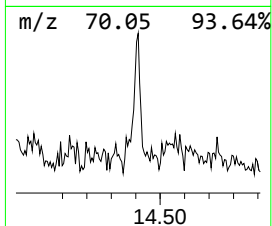
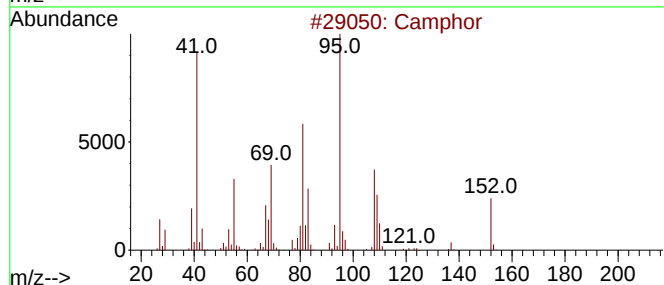
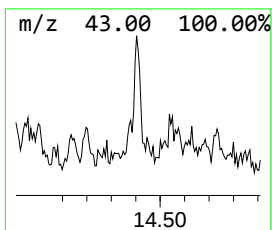
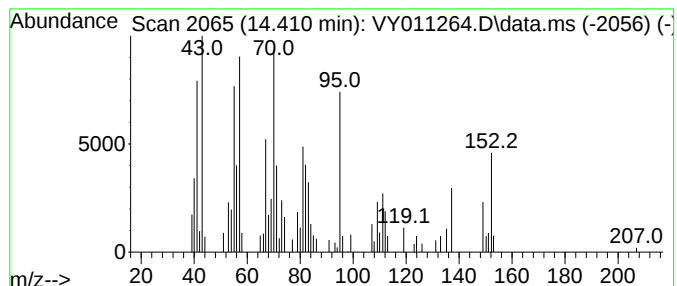
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown14.410 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.410	9.28 ug/l	33686	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Camphor	152	C10H16O	000076-22-2	30
2			Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	30
3			1-Methyldecahydronaphthalene	152	C11H20	002958-75-0	30
4			Diisopropylaminoacrylonitril	152	C9H16N2	038691-28-0	25
5			2-Butanone, 4-(5-methyl-2-furanyl)-	152	C9H12O2	013679-56-6	25



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110222\
Data File : VY011264.D
Acq On : 02 Nov 2022 15:49
Operator : KP/MD
Sample : N5395-08
Misc : 5.12g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
S-2-103122

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y110122S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
cis-1-Butyl-2-m...	12.642	6.6	ug/l	23840	4	13.422	181472	50.0
Cyclohexane, 1,...	12.812	7.2	ug/l	25939	4	13.422	181472	50.0
Bicyclo[2.2.1]h...	13.745	6.0	ug/l	21857	4	13.422	181472	50.0
2-Amino-m-creso...	13.959	12.5	ug/l	45390	4	13.422	181472	50.0
(+)-2-Bornanone	14.251	5.5	ug/l	20104	4	13.422	181472	50.0
unknown14.410	14.410	9.3	ug/l	33686	4	13.422	181472	50.0