

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY071922\
 Data File : VY009601.D
 Acq On : 19 Jul 2022 19:47
 Operator : KP/MD
 Sample : N3652-04
 Misc : 5.75g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 SED-4-(6-12)

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\82Y071822S.M

Title : SW846 8260

Signal : TIC: VY009601.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.686	456	470	486	rBV4	88924	331562	2.61%	0.738%
2	7.710	953	966	971	rBV	160615	380565	3.00%	0.847%
3	7.783	971	978	986	rVB	330112	745789	5.88%	1.660%
4	8.130	1028	1035	1049	rVB	152929	352130	2.77%	0.784%
5	8.319	1049	1066	1078	rBV	4153496	10055324	79.23%	22.384%
6	8.685	1116	1126	1137	rBV	507245	1036906	8.17%	2.308%
7	9.313	1223	1229	1236	rBV	488720	988409	7.79%	2.200%
8	9.612	1270	1278	1289	rVB	4130226	7806996	61.51%	17.379%
9	9.740	1289	1299	1312	rBV	6294952	12691314	100.00%	28.252%
10	9.862	1312	1319	1324	rVV3	95328	236042	1.86%	0.525%
11	9.929	1324	1330	1336	rVB3	108322	227481	1.79%	0.506%
12	10.002	1336	1342	1347	rBV	76571	138063	1.09%	0.307%
13	10.105	1353	1359	1365	rVB	587639	1040677	8.20%	2.317%
14	10.173	1365	1370	1384	rVB	683287	1241859	9.79%	2.764%
15	10.349	1390	1399	1409	rBV5	78641	310575	2.45%	0.691%
16	10.471	1414	1419	1424	rVB3	105838	192446	1.52%	0.428%
17	10.636	1436	1446	1450	rBV3	129864	330620	2.61%	0.736%
18	10.874	1479	1485	1490	rBV4	69311	140369	1.11%	0.312%
19	11.063	1510	1516	1524	rBV5	96934	252855	1.99%	0.563%
20	11.142	1524	1529	1543	rVB4	224724	543656	4.28%	1.210%
21	11.270	1545	1550	1553	rBV2	72288	136337	1.07%	0.303%
22	11.404	1564	1572	1575	rBV6	86594	210324	1.66%	0.468%
23	11.489	1580	1586	1592	rVV	775314	1261254	9.94%	2.808%
24	11.745	1624	1628	1637	rVB5	91619	188672	1.49%	0.420%
25	12.477	1741	1748	1756	rVB3	552800	1066719	8.41%	2.375%
26	12.623	1768	1772	1782	rVB6	56171	148600	1.17%	0.331%
27	12.928	1818	1822	1829	rVB	89812	179291	1.41%	0.399%
28	12.995	1829	1833	1838	rBV	96064	163880	1.29%	0.365%
29	13.233	1868	1872	1876	rVV	101997	140810	1.11%	0.313%
30	13.282	1877	1880	1888	rVB3	95647	154278	1.22%	0.343%
31	13.422	1889	1903	1909	rBV	746317	1381044	10.88%	3.074%
32	13.745	1953	1956	1972	rVB9	62730	156372	1.23%	0.348%
33	13.965	1987	1992	2002	rVB2	109968	213361	1.68%	0.475%
34	14.147	2015	2022	2028	rVB4	106146	254782	2.01%	0.567%
35	14.257	2036	2040	2054	rVB5	59077	222662	1.75%	0.496%

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Instrument :
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Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y071822S.M
Title : SW846 8260

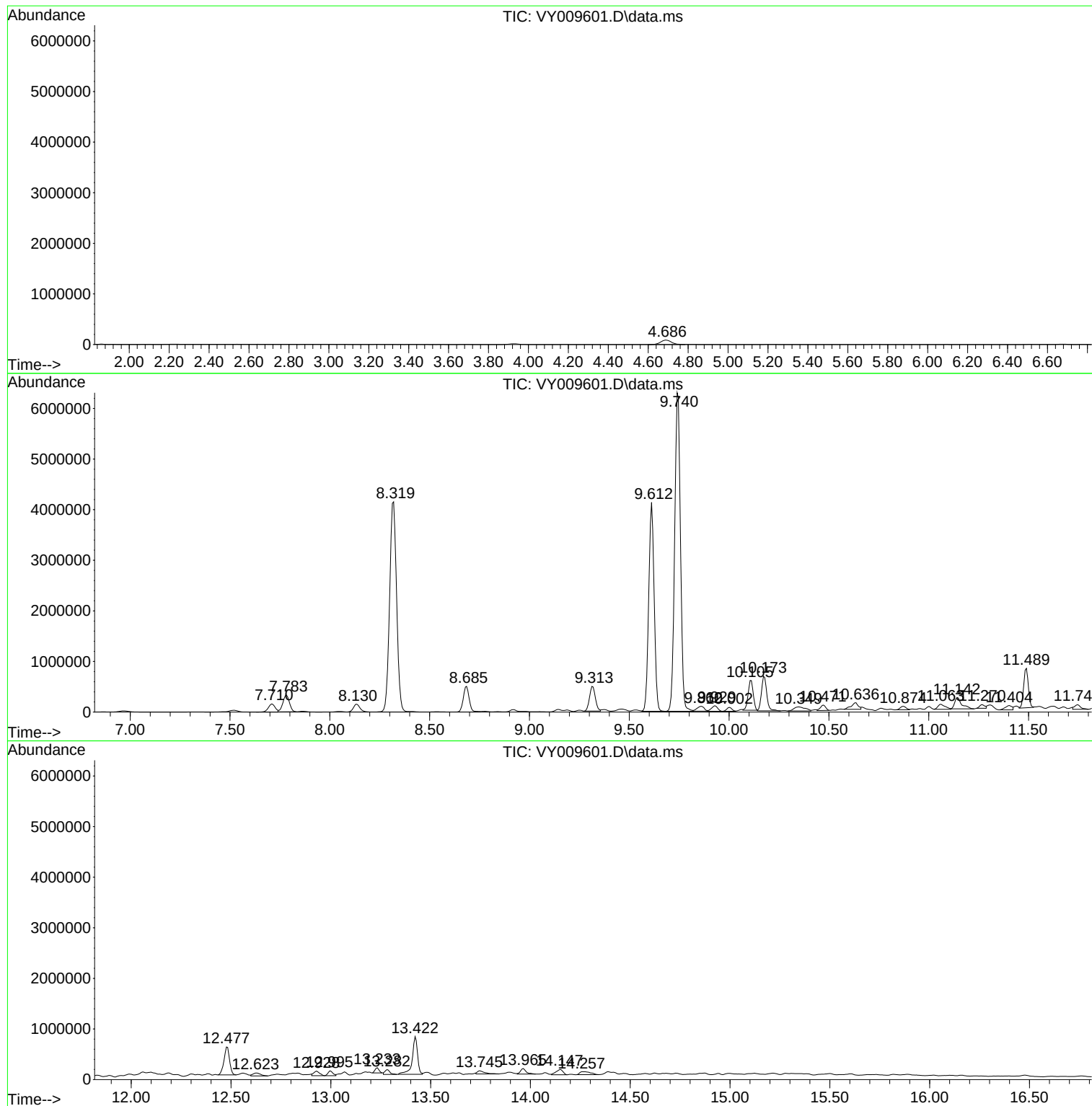
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y071822S.M
Quant Title : SW846 8260

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



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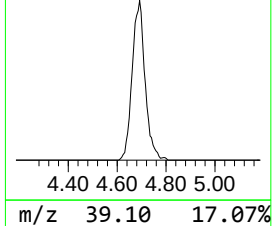
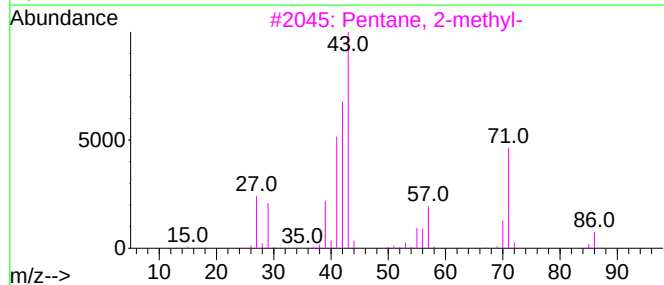
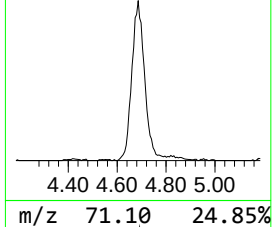
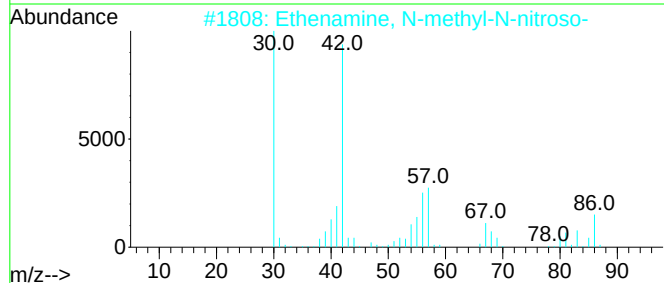
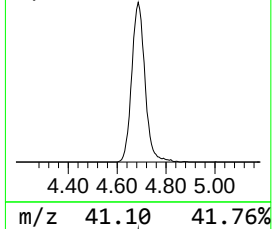
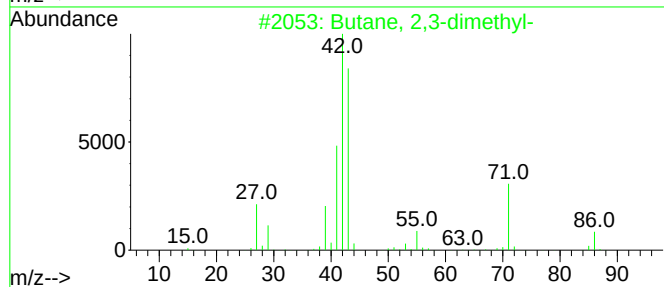
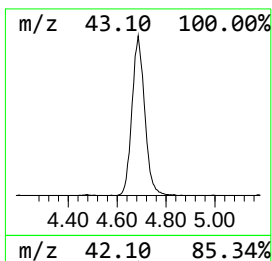
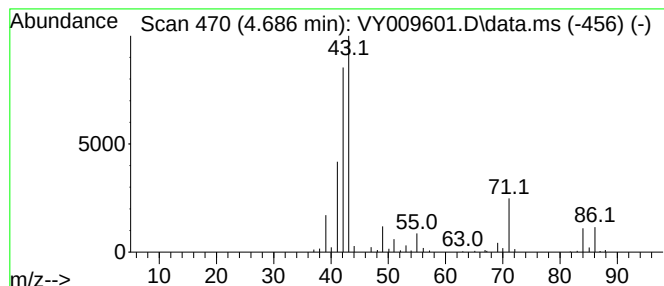
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y071822S.M
Quant Title : SW846 8260

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

Peak Number 1 Butane, 2,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.686	22.23 ug/l	331562	Pentafluorobenzene	7.783

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	87
2			Ethenamine, N-methyl-N-nitroso-	86	C3H6N2O	004549-40-0	47
3			Pentane, 2-methyl-	86	C6H14	000107-83-5	36
4			Tetrahydrofuran	72	C4H8O	000109-99-9	33
5			Propanoyl chloride, 2-methyl-	106	C4H7ClO	000079-30-1	10



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MSVOA_Y
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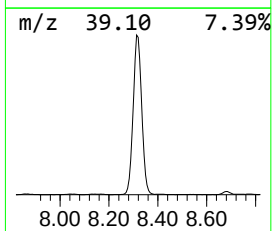
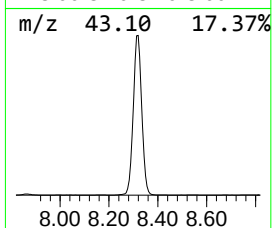
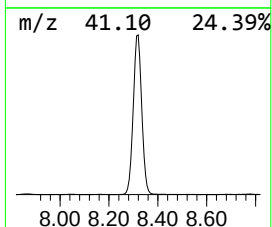
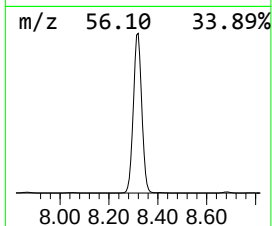
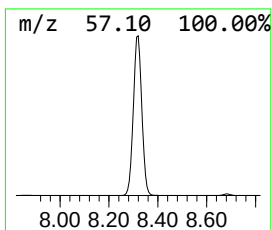
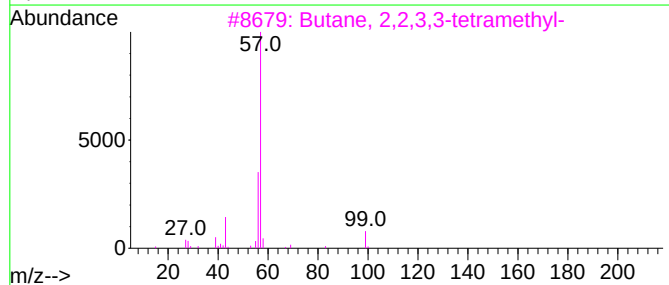
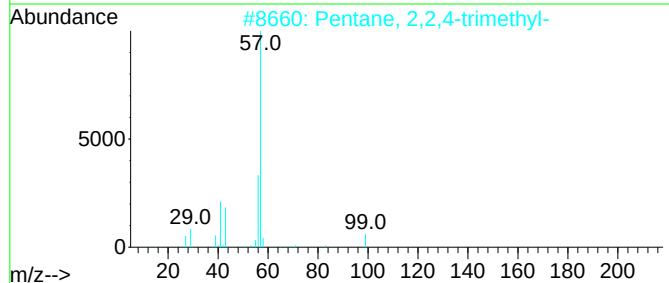
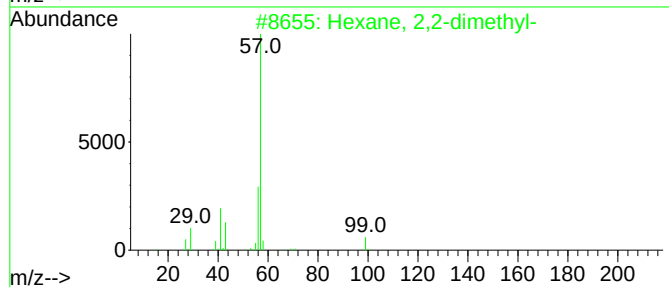
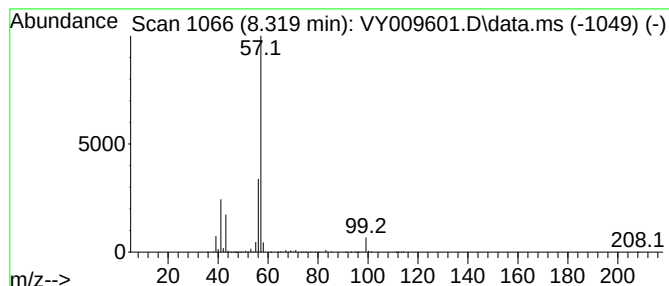
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y071822S.M
Quant Title : SW846 8260

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

Peak Number 2 Hexane, 2,2-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.319	484.87 ug/l	10055300	1,4-Difluorobenzene	8.685

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	83
2			Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	83
3			Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	83
4			Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	78
5			Pentane, 2,2,3-trimethyl-	114	C8H18	000564-02-3	64



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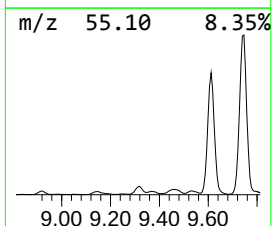
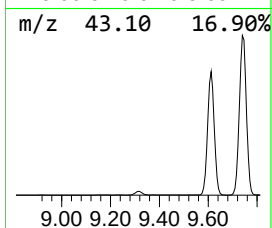
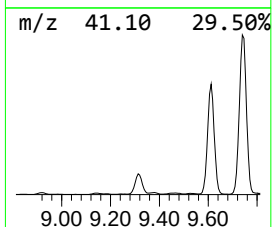
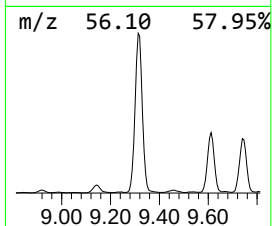
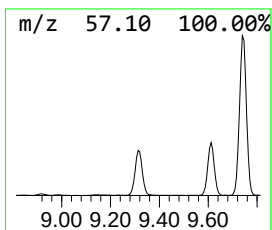
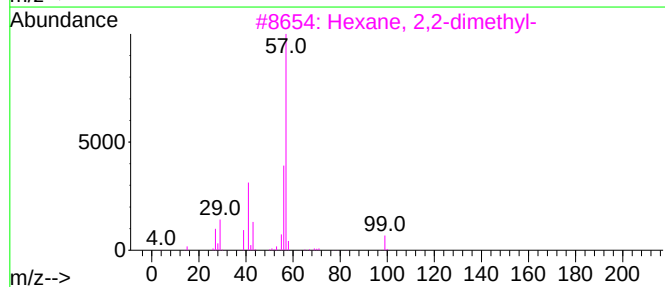
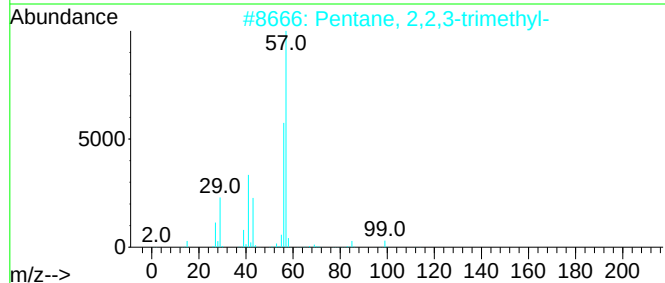
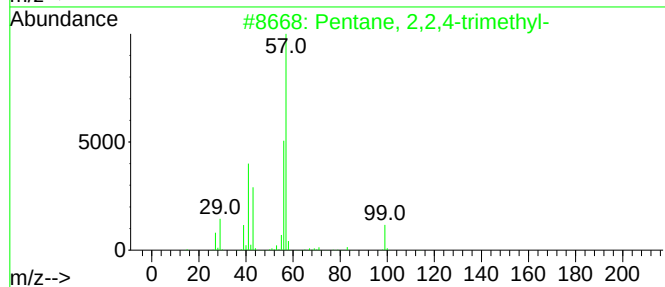
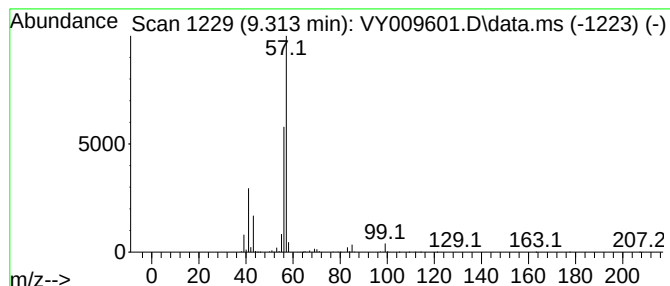
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y071822S.M
Quant Title : SW846 8260

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

Peak Number 3 Pentane, 2,2,4-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.313	47.66 ug/l	988409	1,4-Difluorobenzene	8.685

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	72
2			Pentane, 2,2,3-trimethyl-	114	C8H18	000564-02-3	72
3			Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	64
4			Pentane, 3-methyl-	86	C6H14	000096-14-0	59
5			Heptane, 2,2,4-trimethyl-	142	C10H22	014720-74-2	50



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Misc : 5.75g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SED-4-(6-12)

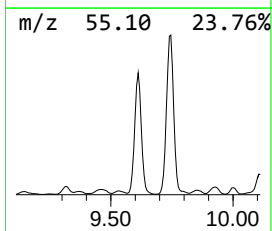
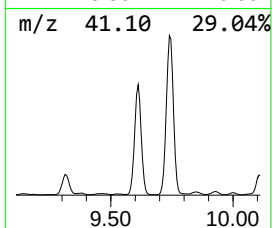
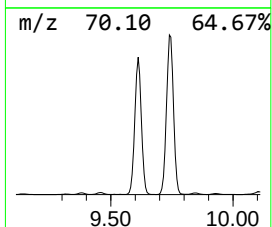
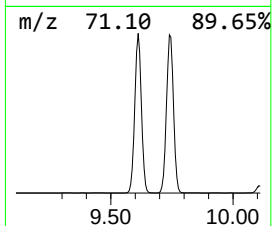
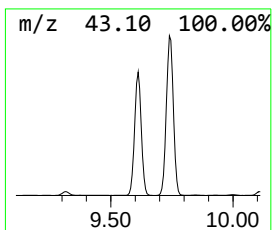
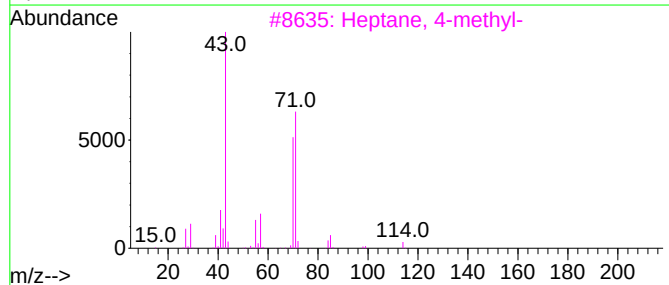
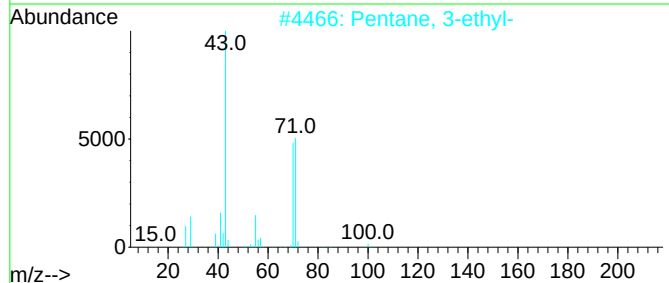
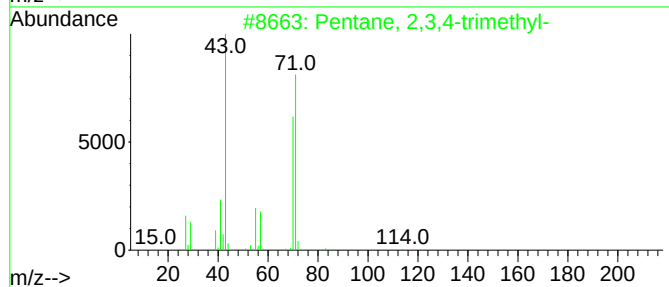
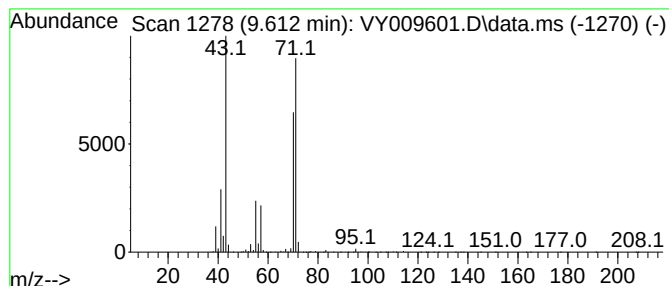
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y071822S.M
Quant Title : SW846 8260

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

Peak Number 4 Pentane, 2,3,4-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.612	376.46 ug/l	7807000	1,4-Difluorobenzene	8.685

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	91
2			Pentane, 3-ethyl-	100	C7H16	000617-78-7	83
3			Heptane, 4-methyl-	114	C8H18	000589-53-7	83
4			Heptane, 3,3,4-trimethyl-	142	C10H22	020278-87-9	78
5			Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	78



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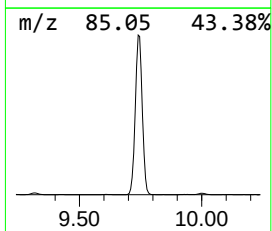
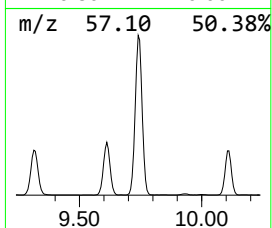
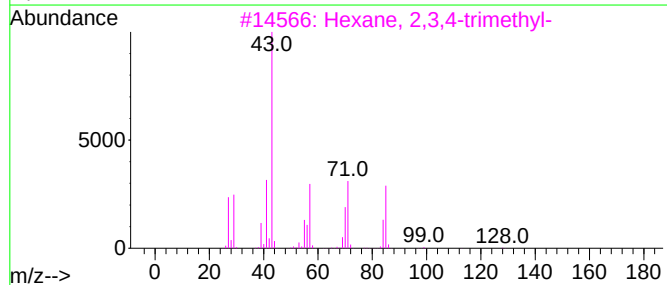
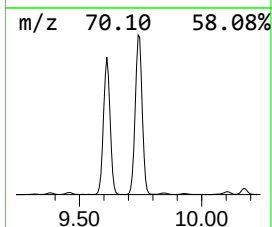
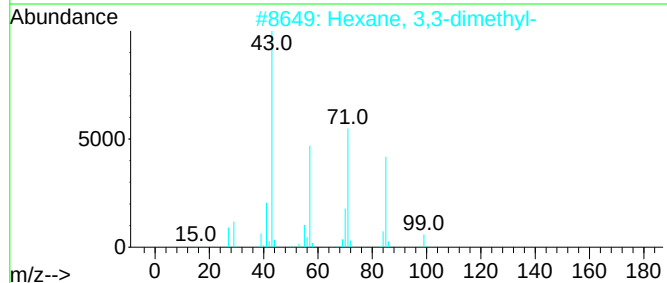
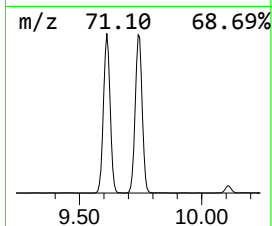
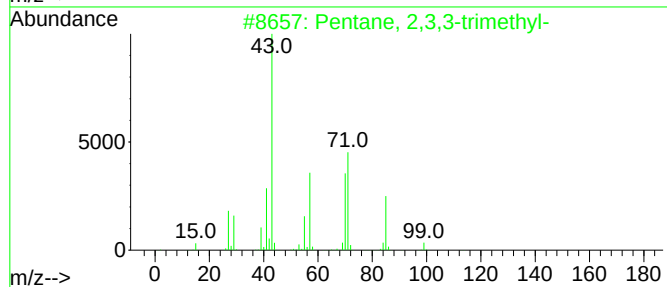
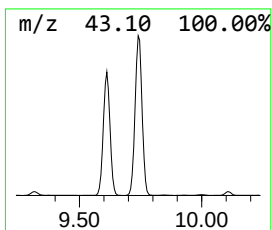
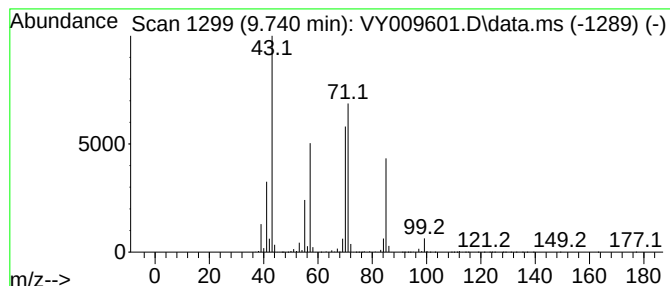
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y071822S.M
Quant Title : SW846 8260

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

Peak Number 5 Pentane, 2,3,3-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.740	611.98 ug/l	12691300	1,4-Difluorobenzene	8.685

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	90
2			Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	64
3			Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	64
4			Heptane, 4-methyl-	114	C8H18	000589-53-7	53
5			Pentane, 2,2,3,3-tetramethyl-	128	C9H20	007154-79-2	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY071922\
Data File : VY009601.D
Acq On : 19 Jul 2022 19:47
Operator : KP/MD
Sample : N3652-04
Misc : 5.75g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SED-4-(6-12)

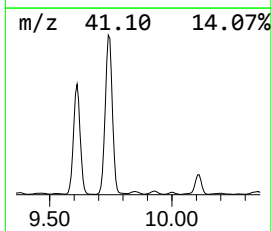
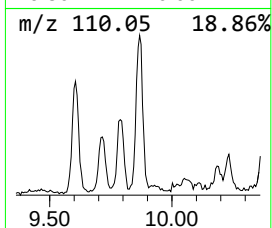
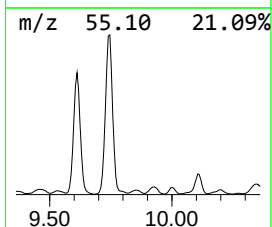
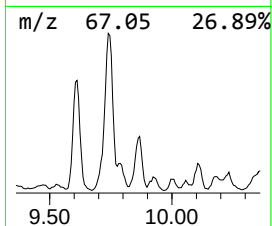
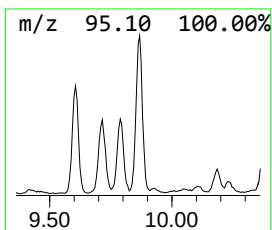
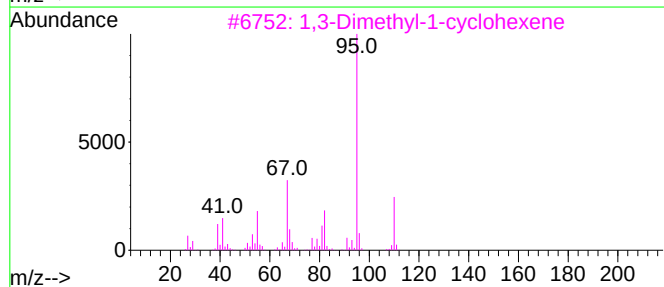
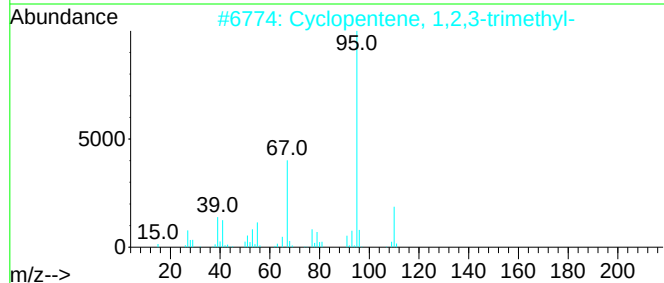
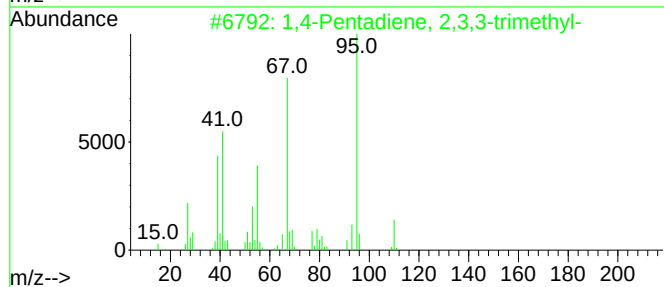
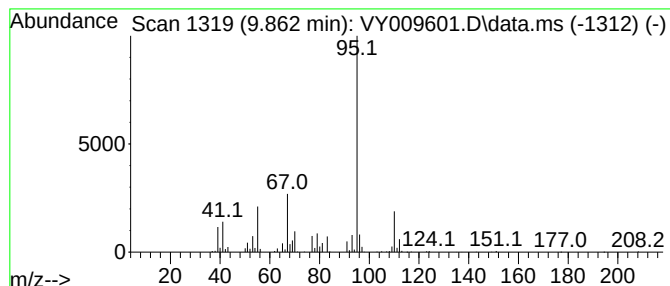
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Quant Title : SW846 8260

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

Peak Number 6 1,4-Pentadiene, 2,3,3-trime... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.862	11.38 ug/l	236042	1,4-Difluorobenzene	8.685

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1,4-Pentadiene, 2,3,3-trimethyl-	110	C8H14	000756-02-5	87
2		Cyclopentene, 1,2,3-trimethyl-	110	C8H14	000473-91-6	87
3		1,3-Dimethyl-1-cyclohexene	110	C8H14	002808-76-6	83
4		1,4-Pentadiene, 2,3,4-trimethyl-	110	C8H14	072014-90-5	59
5		Bicyclo[3.1.0]hexane, 1,5-dimethyl-	110	C8H14	1010142-17-5	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY071922\
Data File : VY009601.D
Acq On : 19 Jul 2022 19:47
Operator : KP/MD
Sample : N3652-04
Misc : 5.75g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SED-4-(6-12)

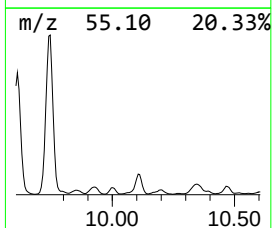
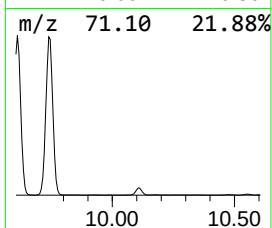
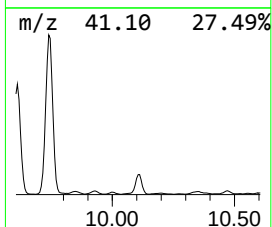
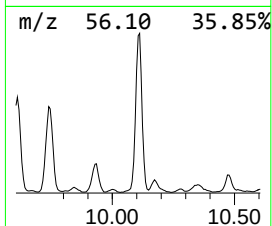
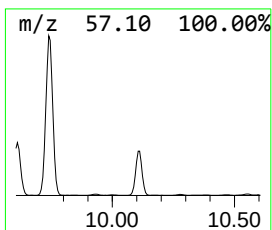
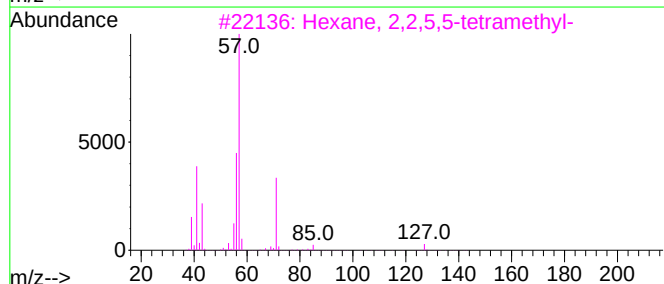
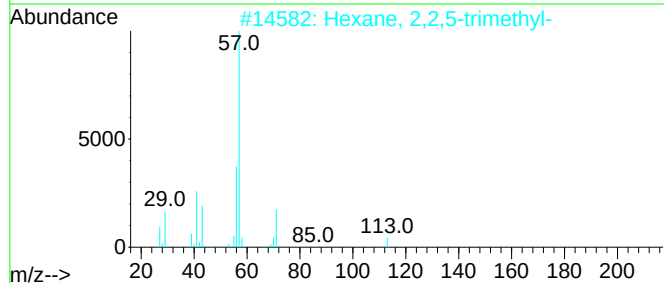
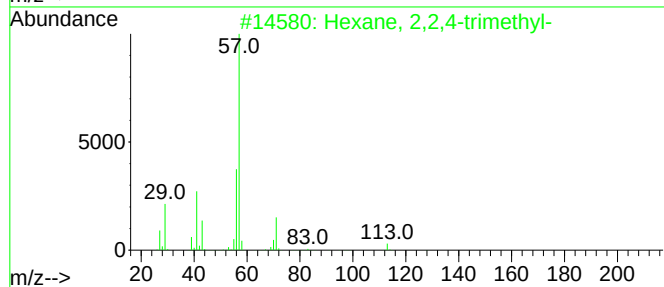
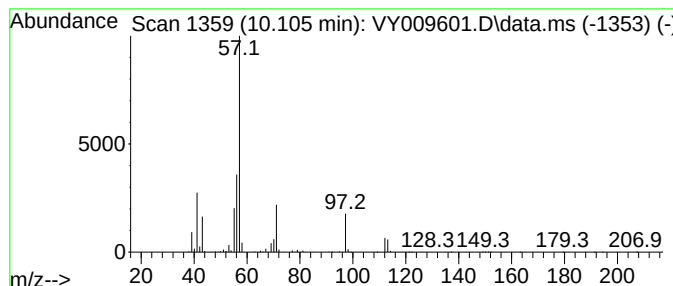
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Quant Title : SW846 8260

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

Peak Number 7 Hexane, 2,2,4-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.105	41.26 ug/l	1040680	Chlorobenzene-d5	11.489

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	59
2			Hexane, 2,2,5-trimethyl-	128	C9H20	003522-94-9	59
3			Hexane, 2,2,5,5-tetramethyl-	142	C10H22	001071-81-4	53
4			Heptane, 2,2-dimethyl-	128	C9H20	001071-26-7	50
5			1-Hexene, 5,5-dimethyl-	112	C8H16	007116-86-1	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY071922\
Data File : VY009601.D
Acq On : 19 Jul 2022 19:47
Operator : KP/MD
Sample : N3652-04
Misc : 5.75g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SED-4-(6-12)

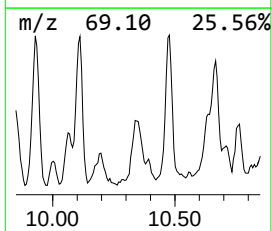
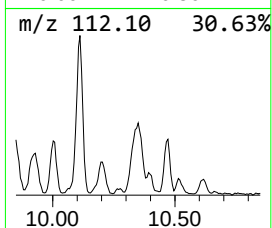
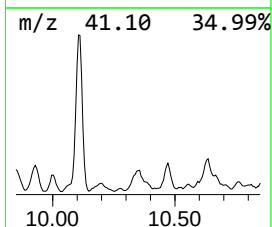
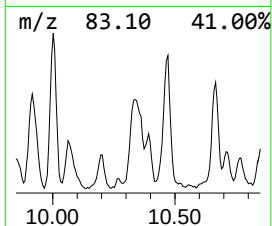
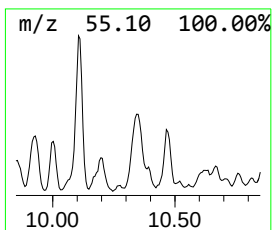
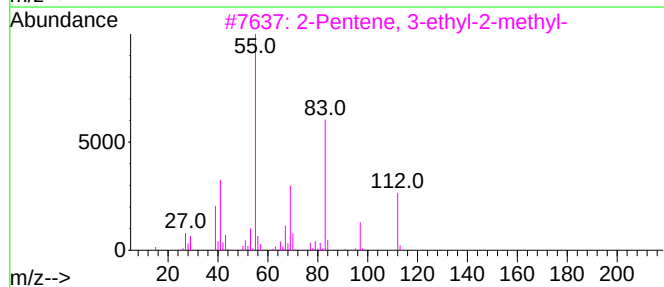
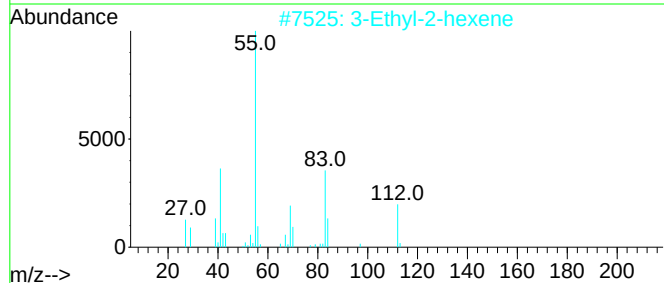
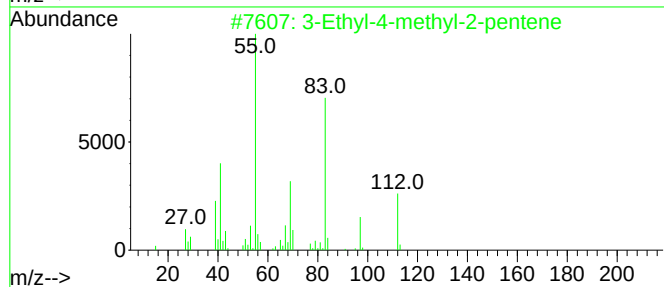
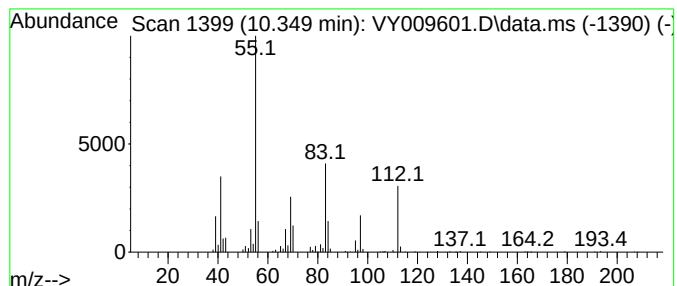
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y071822S.M
Quant Title : SW846 8260

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

Peak Number 8 3-Ethyl-4-methyl-2-pentene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.349	12.31 ug/l	310575	Chlorobenzene-d5	11.489

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Ethyl-4-methyl-2-pentene	112	C8H16	019780-68-8	87
2			3-Ethyl-2-hexene	112	C8H16	000620-00-8	87
3			2-Pentene, 3-ethyl-2-methyl-	112	C8H16	019780-67-7	80
4			Cyclohexane, 1,2-dimethyl-, cis-	112	C8H16	002207-01-4	64
5			Hexane, 3-methyl-4-methylene-	112	C8H16	003404-67-9	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY071922\
Data File : VY009601.D
Acq On : 19 Jul 2022 19:47
Operator : KP/MD
Sample : N3652-04
Misc : 5.75g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SED-4-(6-12)

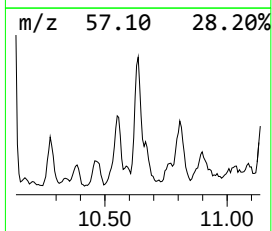
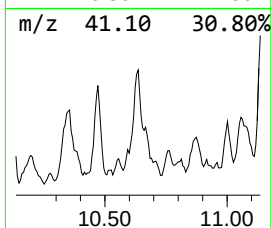
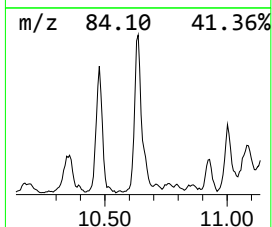
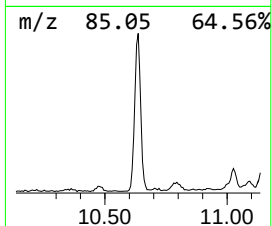
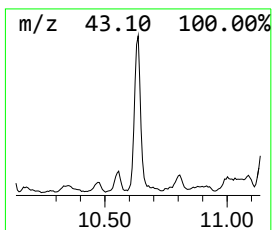
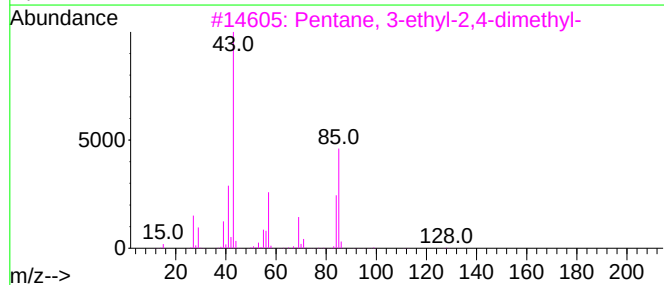
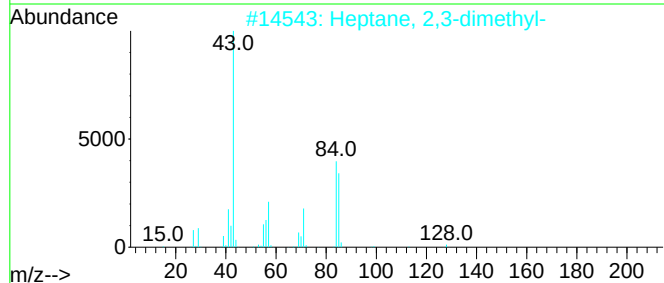
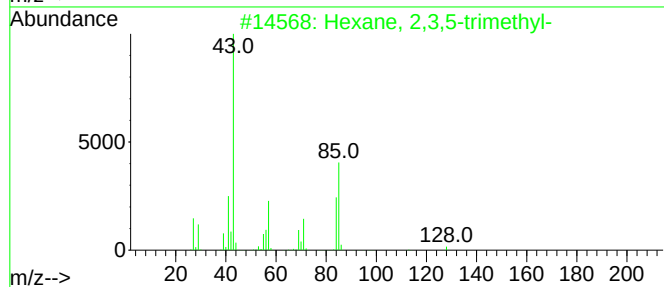
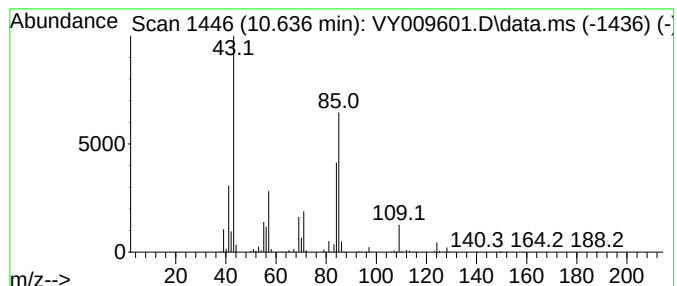
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y071822S.M
Quant Title : SW846 8260

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

Peak Number 9 Hexane, 2,3,5-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.636	13.11 ug/l	330620	Chlorobenzene-d5	11.489

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 2,3,5-trimethyl-	128	C9H20	001069-53-0	68
2			Heptane, 2,3-dimethyl-	128	C9H20	003074-71-3	62
3			Pentane, 3-ethyl-2,4-dimethyl-	128	C9H20	001068-87-7	59
4			Heptane, 3,4,5-trimethyl-	142	C10H22	020278-89-1	53
5			Heptane, 4,4-dimethyl-	128	C9H20	001068-19-5	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY071922\
Data File : VY009601.D
Acq On : 19 Jul 2022 19:47
Operator : KP/MD
Sample : N3652-04
Misc : 5.75g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SED-4-(6-12)

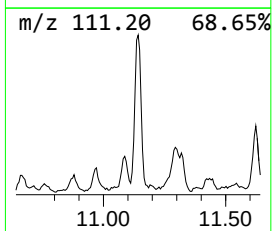
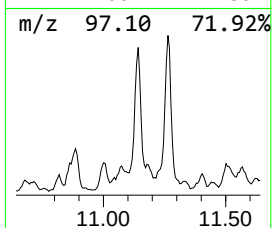
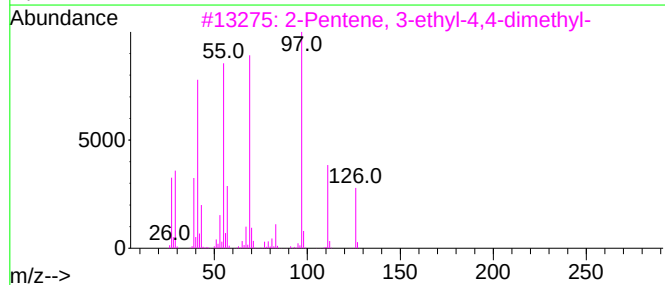
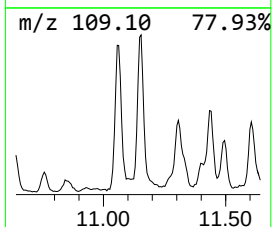
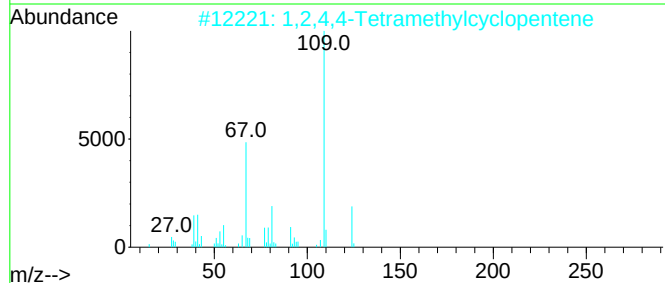
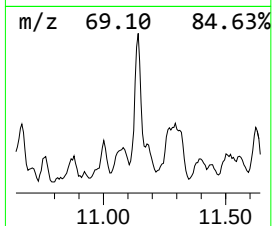
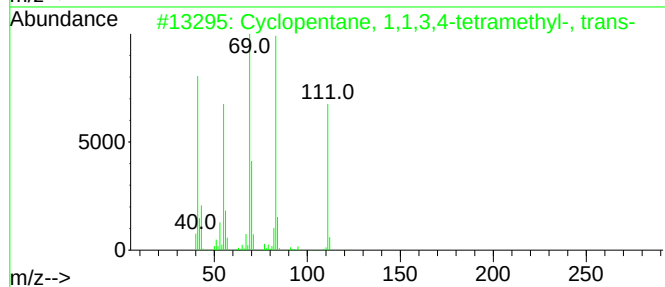
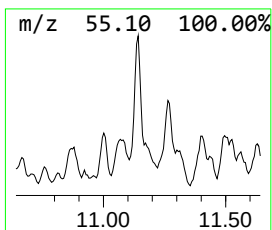
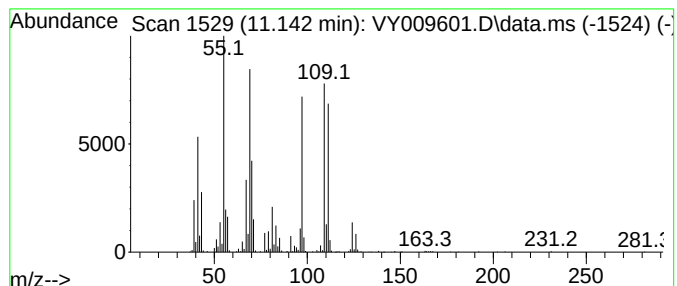
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y071822S.M
Quant Title : SW846 8260

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

Peak Number 10 unknown11.142 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.142	21.55 ug/l	543656	Chlorobenzene-d5	11.489

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, 1,1,3,4-tetramethyl...	126	C9H18	020309-77-7	38
2			1,2,4,4-Tetramethylcyclopentene	124	C9H16	065378-76-9	35
3			2-Pentene, 3-ethyl-4,4-dimethyl-	126	C9H18	053907-59-8	30
4			1-Hexene, 3,3-dimethyl-	112	C8H16	003404-77-1	30
5			6,6-Dimethylhepta-2,4-diene	124	C9H16	1000195-03-3	30



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY071922\
Data File : VY009601.D
Acq On : 19 Jul 2022 19:47
Operator : KP/MD
Sample : N3652-04
Misc : 5.75g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SED-4-(6-12)

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y071822S.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
Butane, 2,3-dim...	4.686	22.2	ug/l	331562	1	7.783	745789	50.0
Hexane, 2,2-dim...	8.319	484.9	ug/l	10055300	2	8.685	1036910	50.0
Pentane, 2,2,4-...	9.313	47.7	ug/l	988409	2	8.685	1036910	50.0
Pentane, 2,3,4-...	9.612	376.5	ug/l	7807000	2	8.685	1036910	50.0
Pentane, 2,3,3-...	9.740	612.0	ug/l	12691300	2	8.685	1036910	50.0
1,4-Pentadiene,...	9.862	11.4	ug/l	236042	2	8.685	1036910	50.0
Hexane, 2,2,4-t...	10.105	41.3	ug/l	1040680	3	11.489	1261250	50.0
3-Ethyl-4-methy...	10.349	12.3	ug/l	310575	3	11.489	1261250	50.0
Hexane, 2,3,5-t...	10.636	13.1	ug/l	330620	3	11.489	1261250	50.0
unknown11.142	11.142	21.6	ug/l	543656	3	11.489	1261250	50.0