

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY092821\
 Data File : VY006192.D
 Acq On : 28 Sep 2021 16:46
 Operator : SY/MD
 Sample : M3933-01
 Misc : 1.06G/5ML/MSVOA_Y/SOIL
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 191-0922-XYL-1

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\82Y092821S.M
 Title : SW846 8260

Signal : TIC: VY006192.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.954	340	350	368	rVB2	25623	71476	3.32%	0.381%
2	4.729	468	477	490	rBV3	8324	27169	1.26%	0.145%
3	5.594	606	619	631	rBV2	7559	26248	1.22%	0.140%
4	6.710	791	802	815	rBV3	24561	72635	3.37%	0.387%
5	6.996	838	849	858	rBV	18483	49834	2.31%	0.265%
6	7.728	957	969	974	rBV	110313	263275	12.22%	1.402%
7	7.795	974	980	991	rVB	201891	448675	20.82%	2.390%
8	8.149	1029	1038	1049	rBV	132323	298900	13.87%	1.592%
9	8.557	1096	1105	1113	rBV2	32402	68199	3.17%	0.363%
10	8.697	1120	1128	1137	rVB	317202	615905	28.59%	3.281%
11	8.899	1150	1161	1186	rBV	106802	332092	15.41%	1.769%
12	9.142	1195	1201	1210	rVB2	35079	66725	3.10%	0.355%
13	9.264	1214	1221	1234	rVB	65377	126927	5.89%	0.676%
14	9.612	1273	1278	1288	rVB	18039	32786	1.52%	0.175%
15	10.075	1347	1354	1364	rVB	923967	1542984	71.62%	8.220%
16	10.185	1364	1372	1378	rBV	523764	900196	41.78%	4.795%
17	10.246	1378	1382	1386	rVV	15541	21867	1.01%	0.116%
18	10.368	1395	1402	1407	rBV	126335	212679	9.87%	1.133%
19	10.441	1407	1414	1433	rBV	908677	1639226	76.08%	8.732%
20	10.807	1467	1474	1486	rVB2	125996	271624	12.61%	1.447%
21	10.935	1488	1495	1505	rBV	468910	740952	34.39%	3.947%
22	11.093	1515	1521	1526	rVB2	20012	35858	1.66%	0.191%
23	11.294	1548	1554	1563	rVB5	15946	40426	1.88%	0.215%
24	11.386	1563	1569	1574	rBV	10379	21958	1.02%	0.117%
25	11.490	1579	1586	1596	rVB	443540	733521	34.05%	3.908%
26	11.593	1596	1603	1611	rBV	279488	464162	21.54%	2.473%
27	11.697	1611	1620	1636	rVB2	1117258	2154524	100.00%	11.477%
28	11.953	1657	1662	1666	rVV4	17491	36220	1.68%	0.193%
29	12.032	1666	1675	1682	rVV	907516	1427629	66.26%	7.605%
30	12.105	1682	1687	1698	rVB	197879	326693	15.16%	1.740%
31	12.209	1698	1704	1709	rBV	132928	207324	9.62%	1.104%
32	12.331	1718	1724	1729	rBV2	31488	54871	2.55%	0.292%
33	12.483	1739	1749	1754	rBV2	359158	592929	27.52%	3.159%
34	12.538	1754	1758	1762	rVB2	17141	25696	1.19%	0.137%
35	12.672	1776	1780	1785	rVB	66105	98526	4.57%	0.525%
36	12.733	1785	1790	1794	rBV	241932	384251	17.83%	2.047%

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37	12.770	1794	1796	1799	rVV	75791	98521	4.57%	0.525%
38	12.812	1799	1803	1816	rVB	249943	409906	19.03%	2.184%
39	12.971	1821	1829	1836	rBV	200815	364658	16.93%	1.943%
40	13.075	1843	1846	1848	rBV2	15799	22914	1.06%	0.122%
41	13.123	1848	1854	1858	rVV	749233	1105544	51.31%	5.889%
42	13.166	1858	1861	1867	rVB2	96103	138393	6.42%	0.737%
43	13.270	1868	1878	1884	rBV4	119457	228753	10.62%	1.219%
44	13.343	1884	1890	1897	rVB3	30679	74726	3.47%	0.398%
45	13.428	1897	1904	1907	rBV	430077	714440	33.16%	3.806%
46	13.459	1907	1909	1913	rVB	189744	217632	10.10%	1.159%
47	13.532	1918	1921	1932	rVV	116673	207915	9.65%	1.108%
48	13.629	1932	1937	1940	rVV2	37409	62969	2.92%	0.335%
49	13.666	1940	1943	1951	rVB3	22356	40161	1.86%	0.214%
50	13.751	1952	1957	1961	rBV5	26917	43453	2.02%	0.231%
51	13.904	1976	1982	1987	rBV4	26966	65051	3.02%	0.347%
52	13.971	1989	1993	1999	rVB4	29800	55137	2.56%	0.294%
53	14.105	2011	2015	2019	rVB	51780	68514	3.18%	0.365%
54	14.209	2028	2032	2037	rVB2	55187	76820	3.57%	0.409%
55	14.410	2062	2065	2072	rVB4	14170	25628	1.19%	0.137%
56	14.672	2103	2108	2114	rBV6	12592	23327	1.08%	0.124%
57	14.739	2114	2119	2126	rVB5	20336	39415	1.83%	0.210%
58	14.958	2151	2155	2160	rVB2	23515	31588	1.47%	0.168%
59	15.239	2196	2201	2210	rVB2	13335	29071	1.35%	0.155%
60	16.275	2365	2371	2376	rVV4	17197	38550	1.79%	0.205%
61	16.336	2376	2381	2391	rVB	78723	153827	7.14%	0.819%

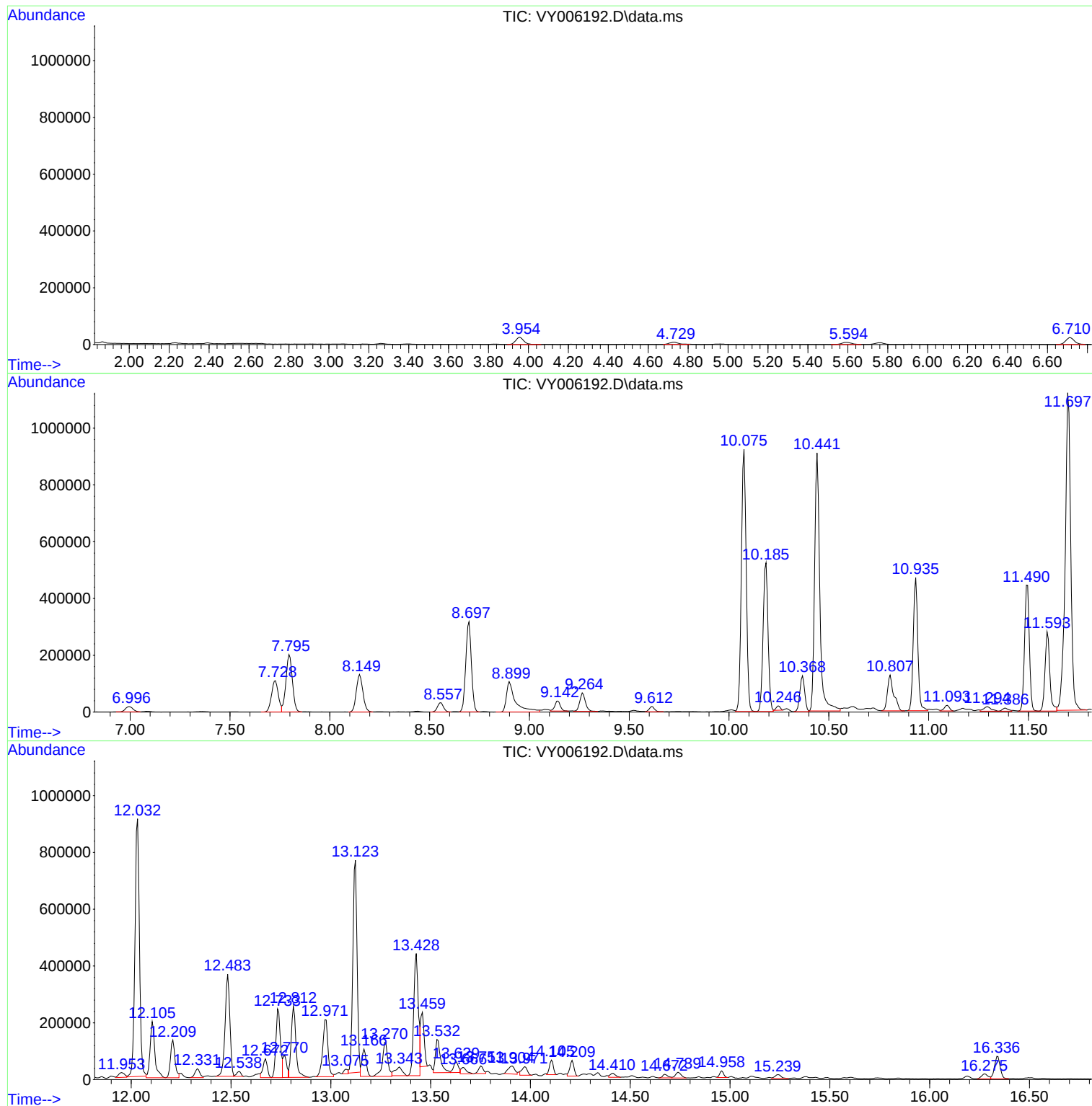
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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Instrument :
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ClientSampleId :
191-0922-XYL-1

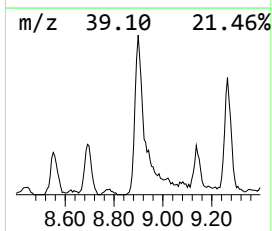
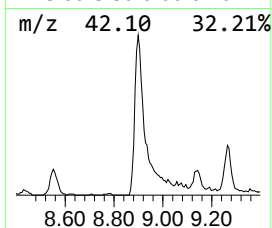
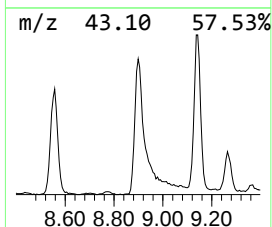
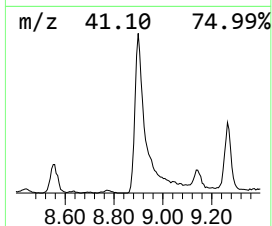
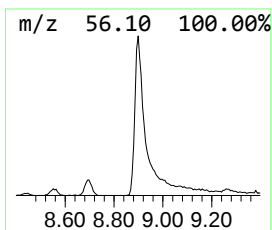
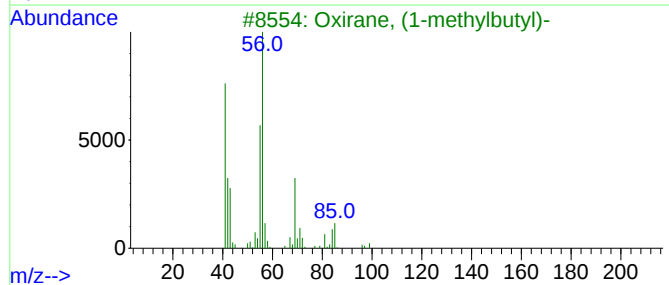
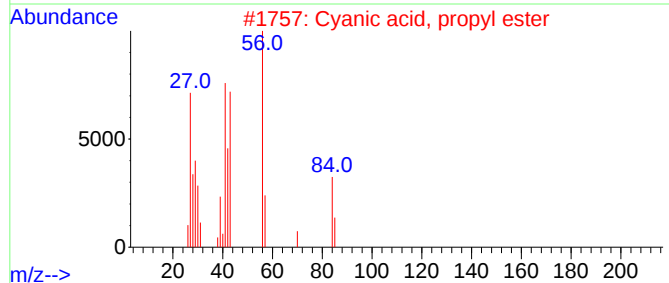
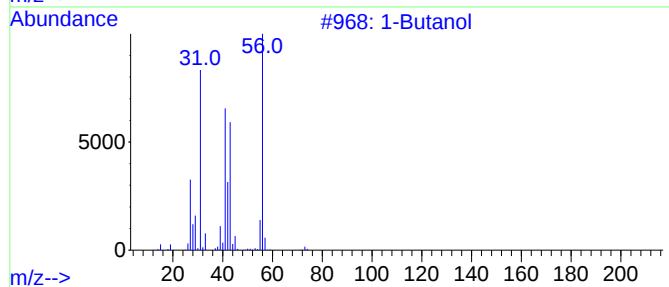
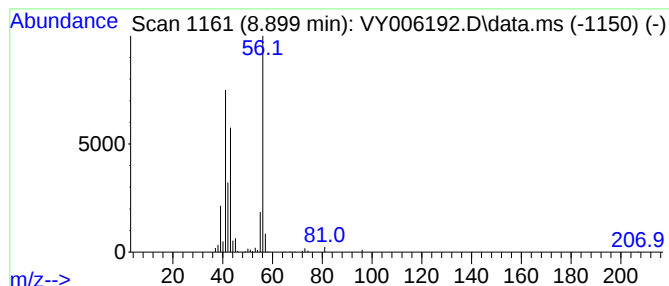
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y092821S.M
Quant Title : SW846 8260

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

Peak Number 1 1-Butanol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.899	26.96 ug/l	332092	1,4-Difluorobenzene	8.697

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Butanol			74	C4H10O	000071-36-3	91
2	Cyanoic acid, propyl ester			85	C4H7NO	001768-36-1	50
3	Oxirane, (1-methylbutyl)-			114	C7H14O	053229-39-3	40
4	1-Pentanol, 4-methyl-			102	C6H14O	000626-89-1	40
5	Propane, 2-cyclopropyl-			84	C6H12	003638-35-5	39



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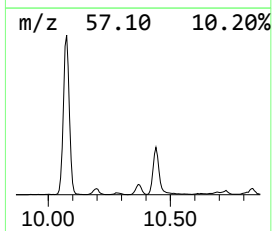
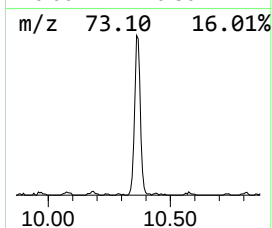
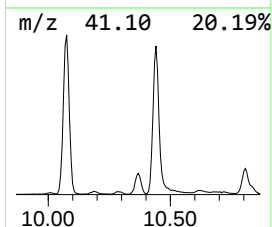
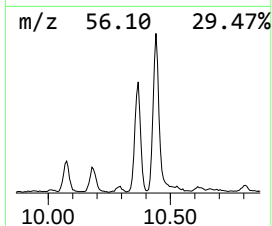
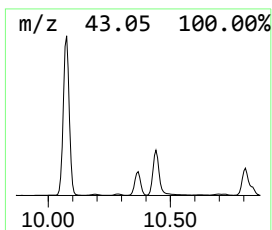
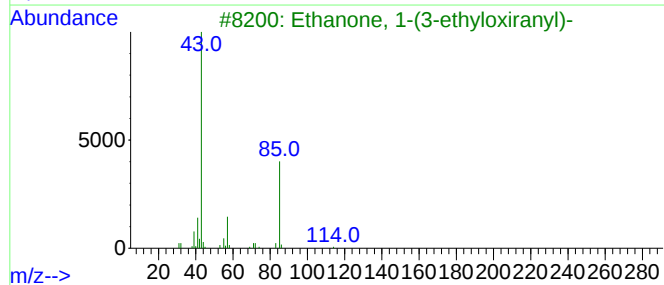
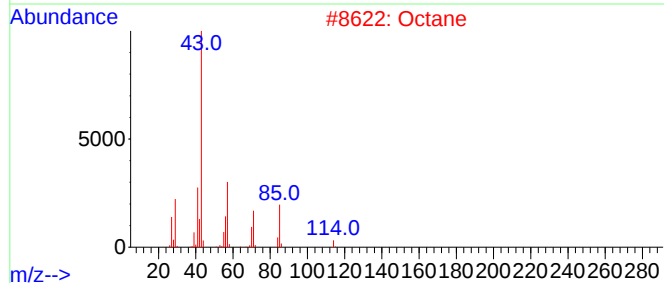
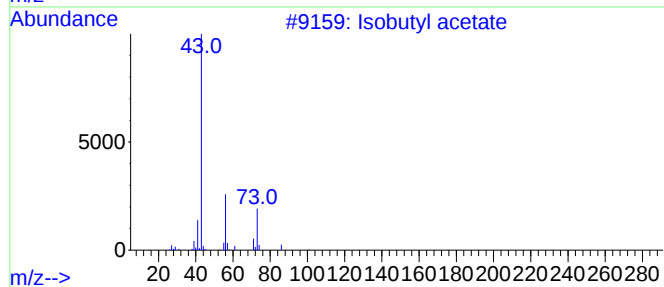
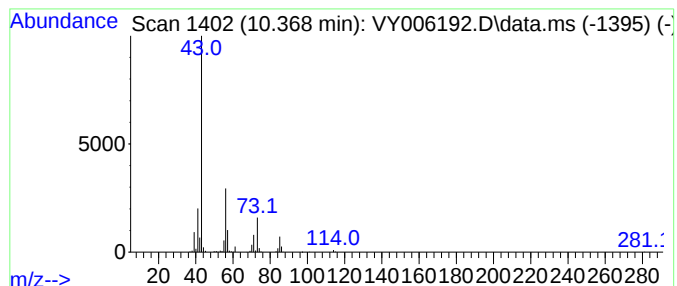
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y092821S.M
Quant Title : SW846 8260

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

Peak Number 2 Isobutyl acetate Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.368	14.50 ug/l	212679	Chlorobenzene-d5	11.496

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isobutyl acetate	116	C6H12O2	000110-19-0	53
2			Octane	114	C8H18	000111-65-9	35
3			Ethanone, 1-(3-ethyloxiranyl)-	114	C6H10O2	017257-81-7	35
4			Acetic acid, butyl ester	116	C6H12O2	000123-86-4	33
5			Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	32



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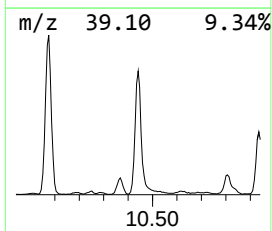
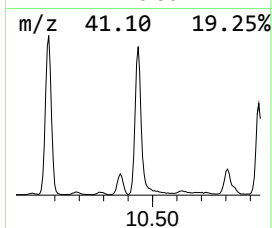
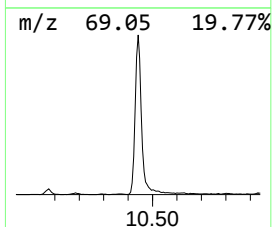
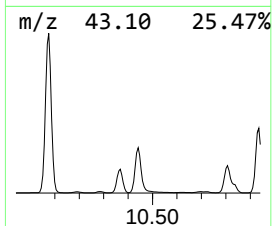
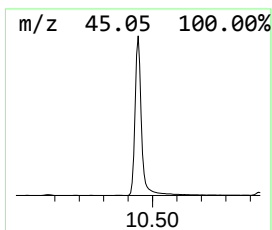
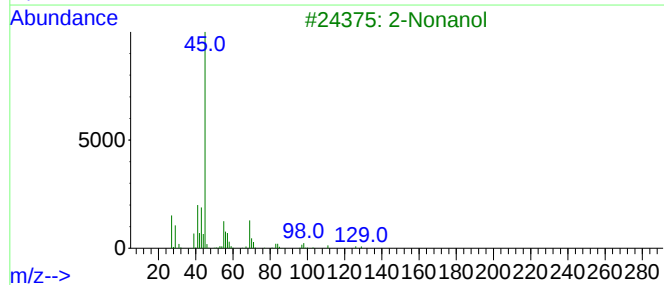
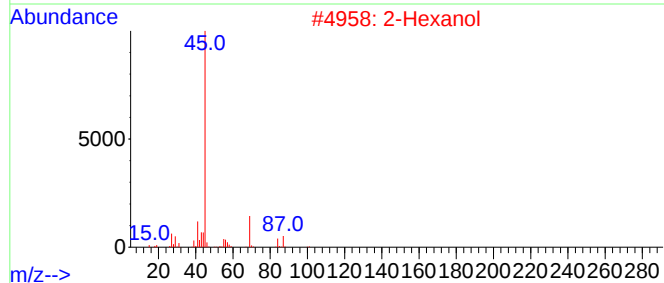
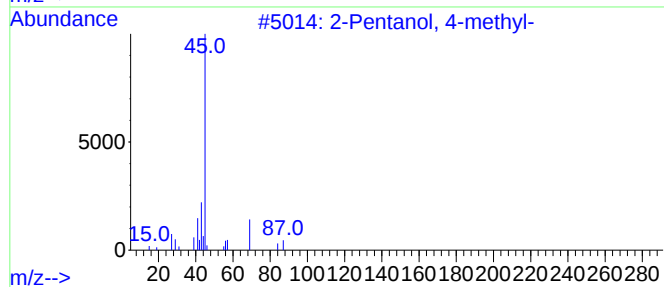
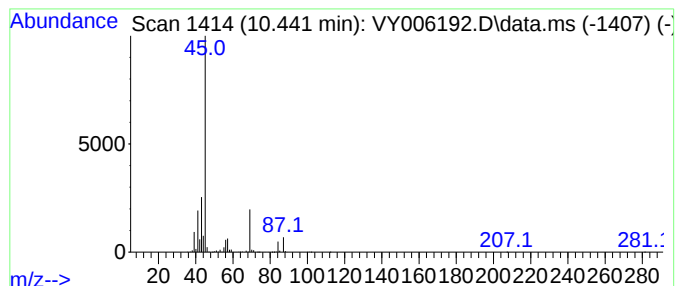
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanol, 4-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.441	111.74 ug/l	1639230	Chlorobenzene-d5	11.496

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanol, 4-methyl-			102	C6H14O	000108-11-2	83
2	2-Hexanol			102	C6H14O	000626-93-7	83
3	2-Nonanol			144	C9H20O	000628-99-9	43
4	2-Heptanol			116	C7H16O	000543-49-7	40
5	2-Butanol, 3-methyl-			88	C5H12O	000598-75-4	9



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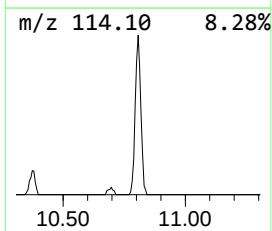
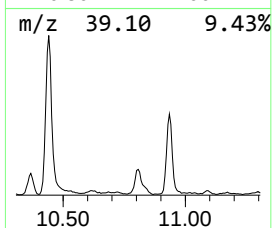
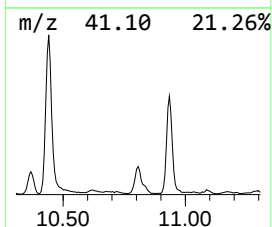
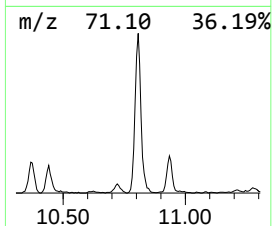
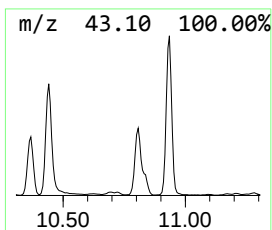
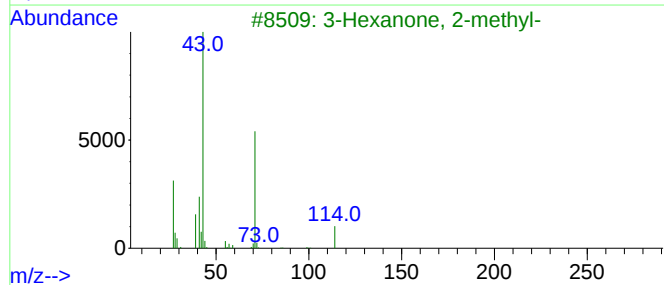
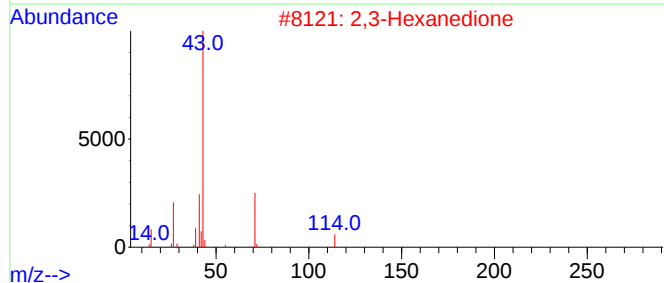
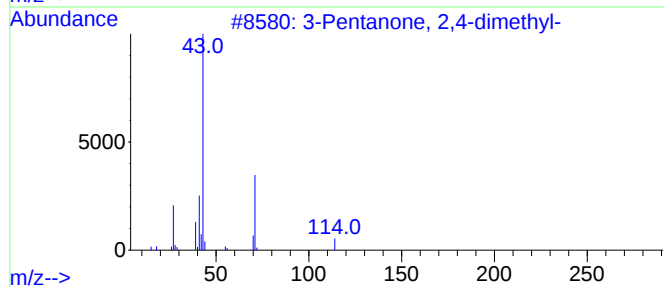
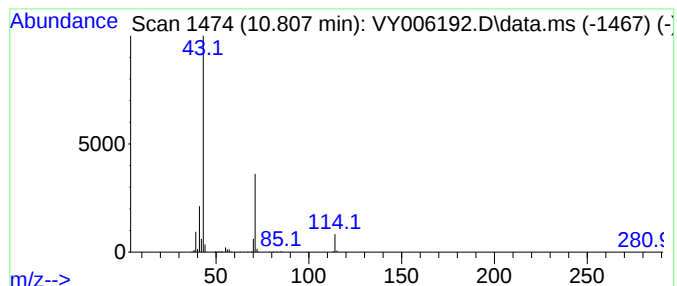
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y092821S.M
Quant Title : SW846 8260

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

Peak Number 4 3-Pentanone, 2,4-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.807	18.52 ug/l	271624	Chlorobenzene-d5	11.496

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Pentanone, 2,4-dimethyl-	114	C7H14O	000565-80-0	86
2			2,3-Hexanedione	114	C6H10O2	003848-24-6	64
3			3-Hexanone, 2-methyl-	114	C7H14O	007379-12-6	59
4			Pyrrolidine	71	C4H9N	000123-75-1	40
5			4-Heptanone	114	C7H14O	000123-19-3	40



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191-0922-XYL-1

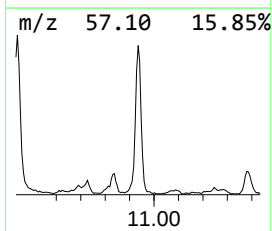
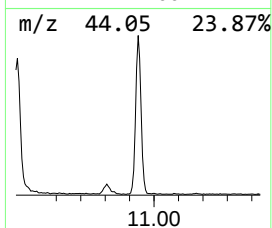
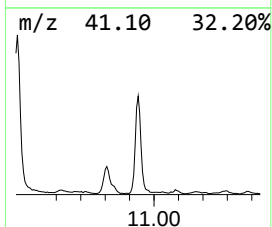
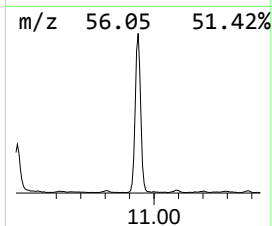
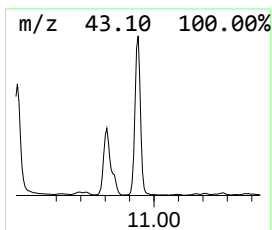
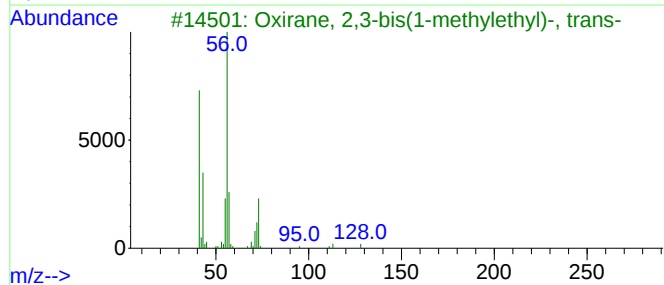
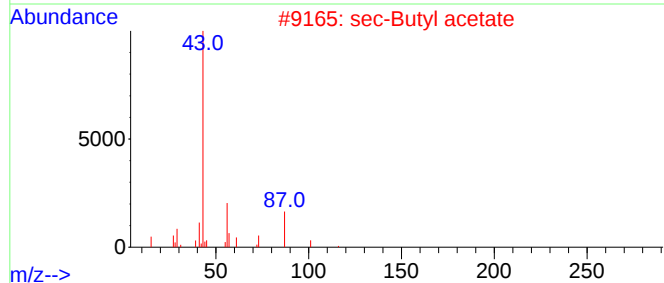
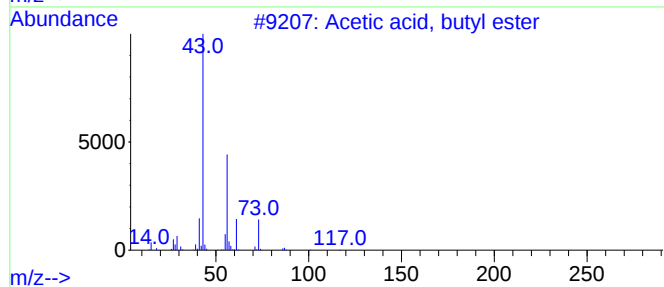
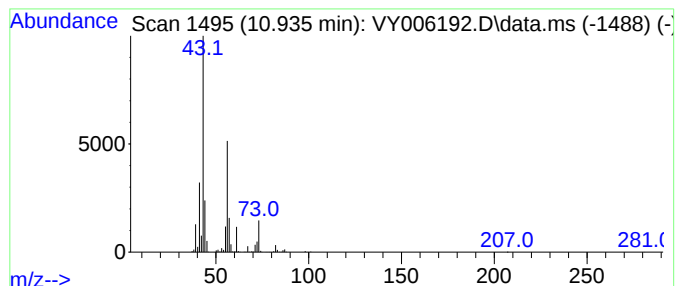
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Quant Title : SW846 8260

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

Peak Number 5 Acetic acid, butyl ester Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.935	50.51 ug/l	740952	Chlorobenzene-d5	11.496

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetic acid, butyl ester	116	C6H12O2	000123-86-4	53
2			sec-Butyl acetate	116	C6H12O2	000105-46-4	38
3			Oxirane, 2,3-bis(1-methylethyl)-...	128	C8H16O	054644-32-5	25
4			4-Methyloxazolidine	87	C4H9NO	155799-98-7	17
5			1-Butanol	74	C4H10O	000071-36-3	12



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY092821\
Data File : VY006192.D
Acq On : 28 Sep 2021 16:46
Operator : SY/MD
Sample : M3933-01
Misc : 1.06G/5ML/MSVOA_Y/SOIL
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
191-0922-XYL-1

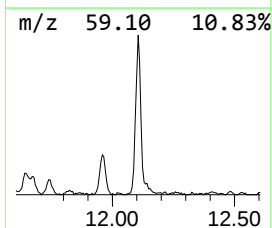
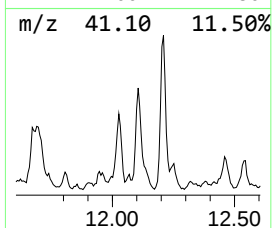
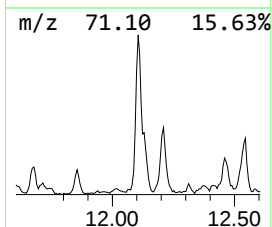
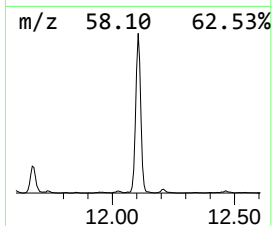
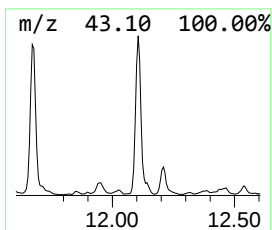
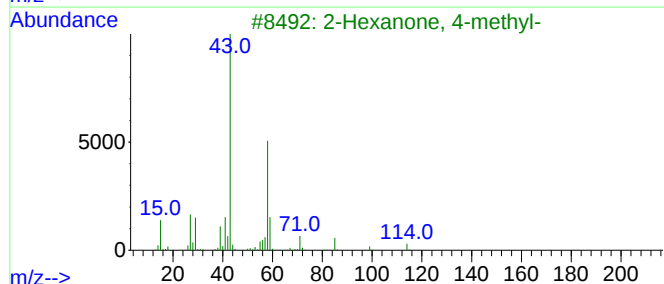
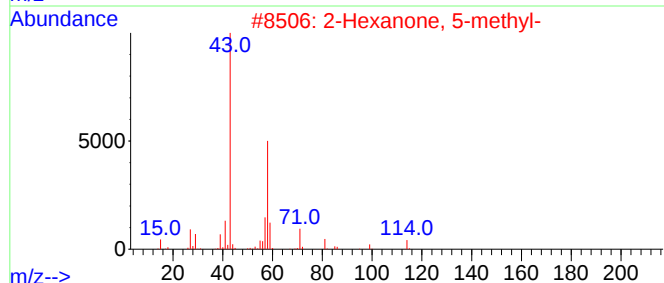
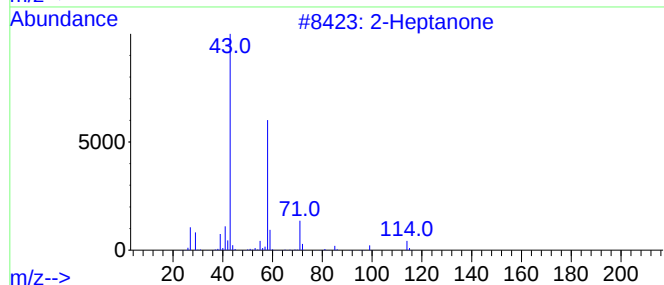
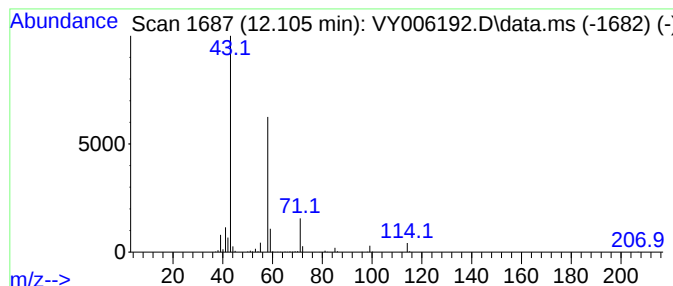
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Quant Title : SW846 8260

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

Peak Number 6 2-Heptanone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.105	22.27 ug/l	326693	Chlorobenzene-d5	11.496

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Heptanone	114	C7H14O	000110-43-0	91
2			2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	86
3			2-Hexanone, 4-methyl-	114	C7H14O	000105-42-0	64
4			2-Pentanone, 4,4-dimethyl-	114	C7H14O	000590-50-1	50
5			1-Butoxy-2-propanol acetate	174	C9H18O3	085409-76-3	39



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY092821\
Data File : VY006192.D
Acq On : 28 Sep 2021 16:46
Operator : SY/MD
Sample : M3933-01
Misc : 1.06G/5ML/MSVOA_Y/SOIL
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
191-0922-XYL-1

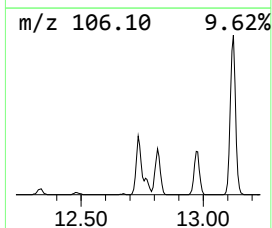
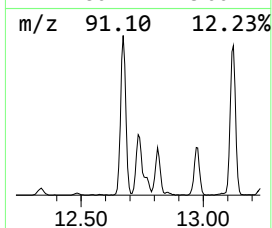
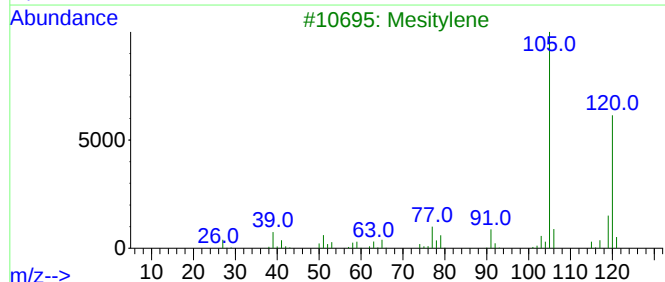
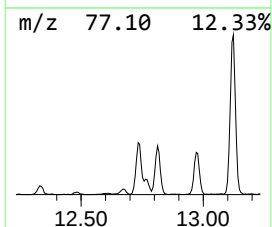
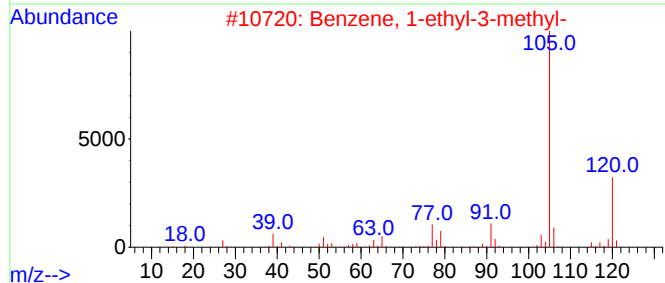
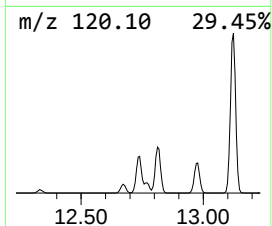
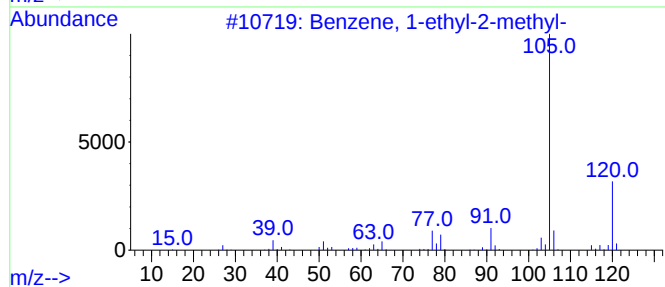
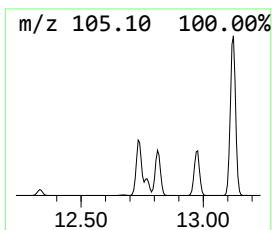
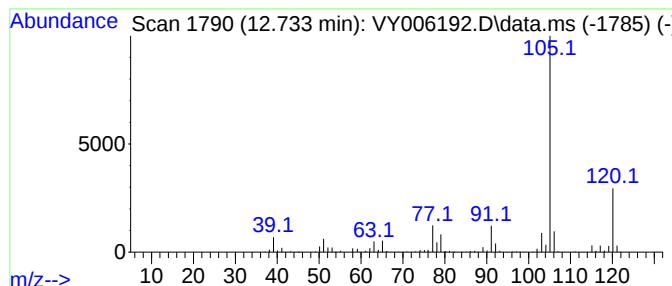
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.733	26.89 ug/l	384251	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY092821\
Data File : VY006192.D
Acq On : 28 Sep 2021 16:46
Operator : SY/MD
Sample : M3933-01
Misc : 1.06G/5ML/MSVOA_Y/SOIL
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
191-0922-XYL-1

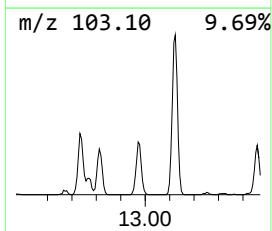
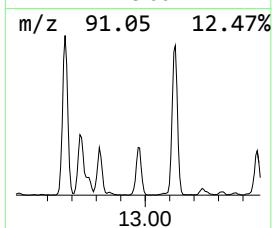
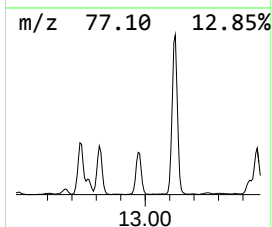
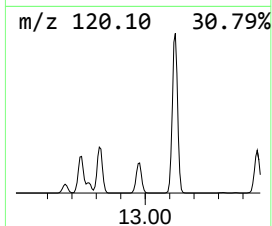
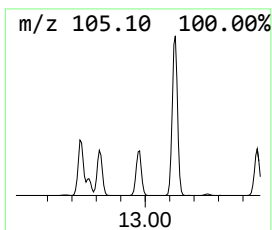
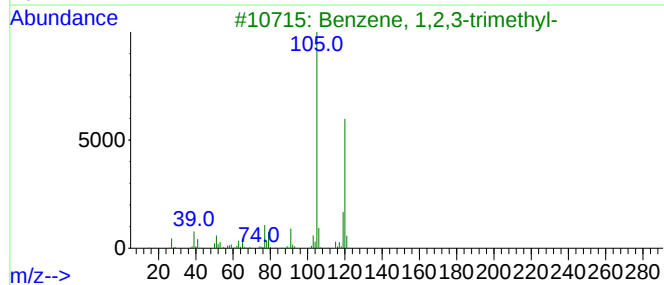
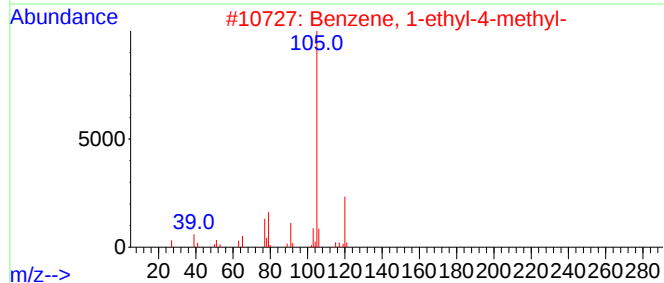
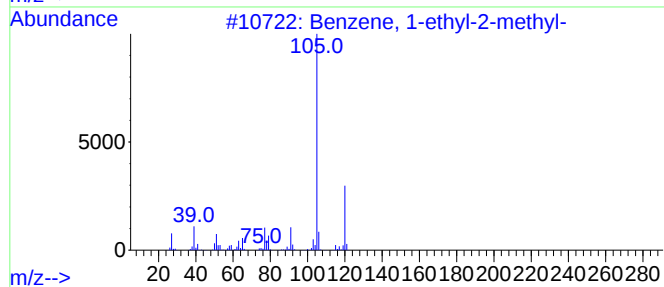
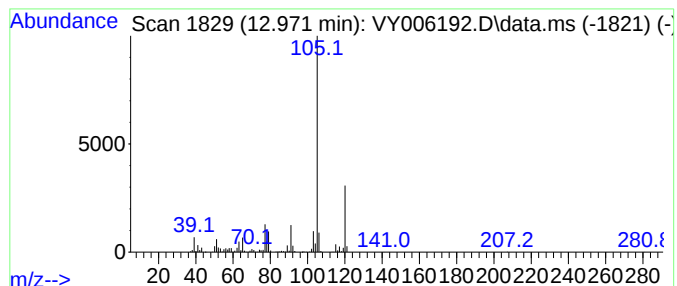
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1-ethyl-4-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.971	25.52 ug/l	364658	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY092821\
Data File : VY006192.D
Acq On : 28 Sep 2021 16:46
Operator : SY/MD
Sample : M3933-01
Misc : 1.06G/5ML/MSVOA_Y/SOIL
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
191-0922-XYL-1

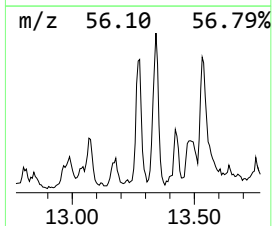
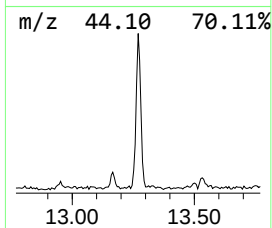
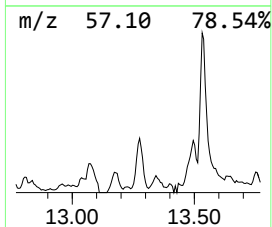
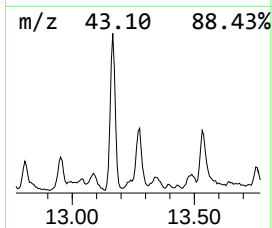
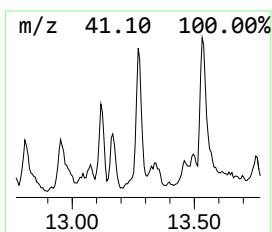
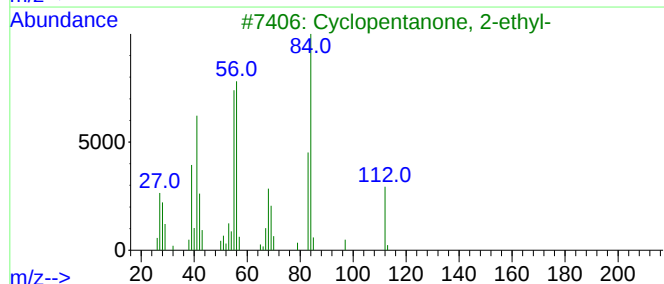
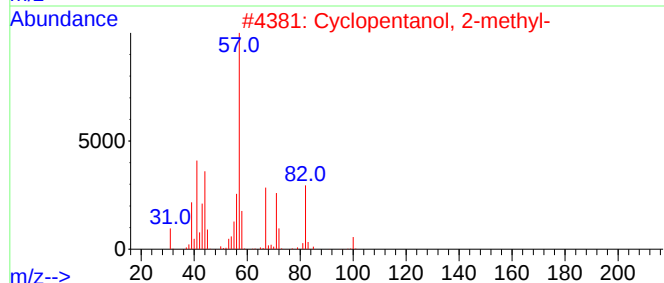
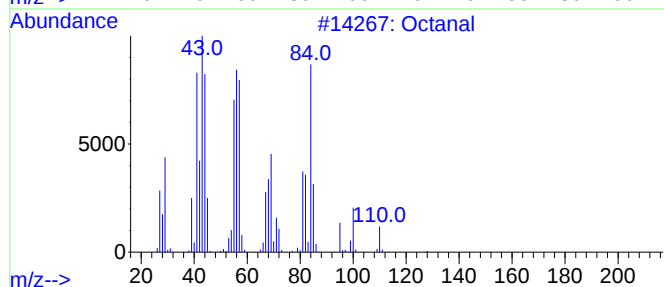
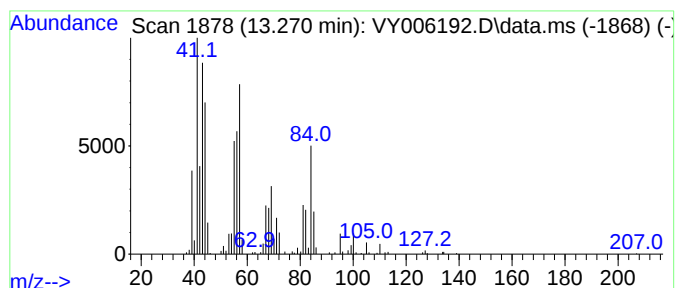
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Quant Title : SW846 8260

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

Peak Number 9 Octanal Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.270	16.01 ug/l	228753	1,4-Dichlorobenzene-d4	13.428

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octanal	128	C8H16O	000124-13-0	91
2		Cyclopentanol, 2-methyl-	100	C6H12O	024070-77-7	38
3		Cyclopentanone, 2-ethyl-	112	C7H12O	004971-18-0	35
4		1-Hexene	84	C6H12	000592-41-6	27
5		Cyclopentanol, 2-methyl-, cis-	100	C6H12O	025144-05-2	27



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY092821\
Data File : VY006192.D
Acq On : 28 Sep 2021 16:46
Operator : SY/MD
Sample : M3933-01
Misc : 1.06G/5ML/MSVOA_Y/SOIL
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
191-0922-XYL-1

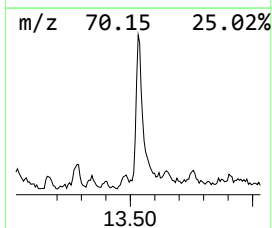
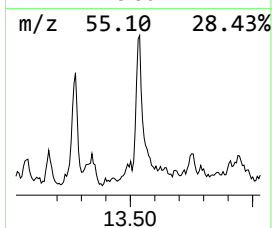
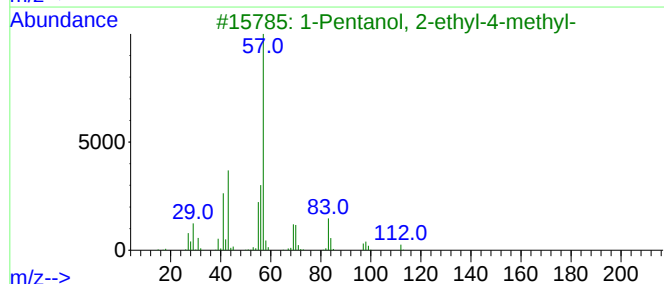
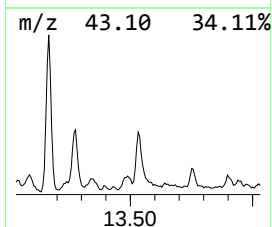
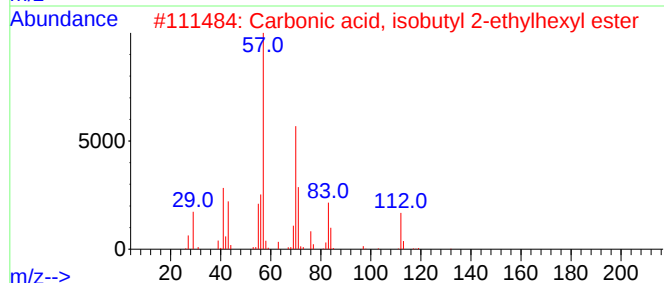
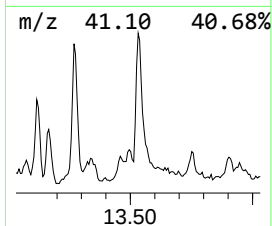
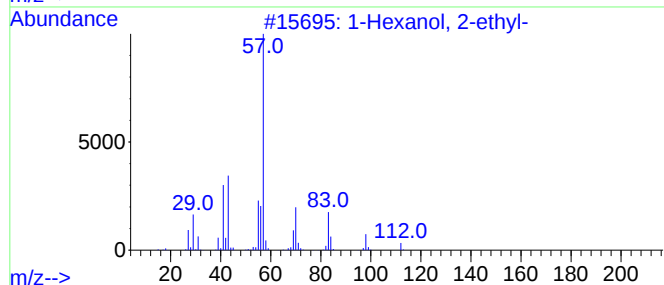
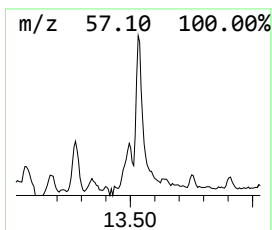
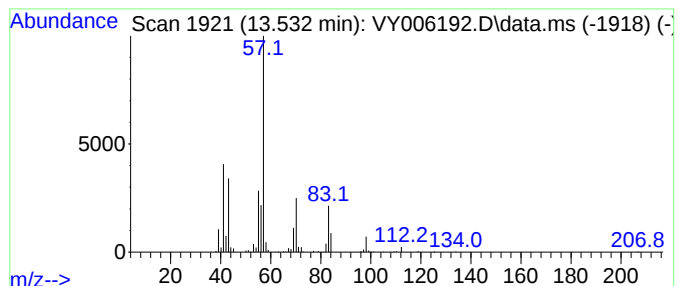
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Quant Title : SW846 8260

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

Peak Number 10 1-Hexanol, 2-ethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.532	14.55 ug/l	207915	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	83
2			Carbonic acid, isobutyl 2-ethylh...	230	C13H26O3	1000383-13-3	59
3			1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	56
4			Heptane, 1,1'-oxybis-	214	C14H30O	000629-64-1	53
5			2-Ethyl-1-hexanol, pentafluoropr...	276	C11H17F5O2	959012-82-9	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY092821\
Data File : VY006192.D
Acq On : 28 Sep 2021 16:46
Operator : SY/MD
Sample : M3933-01
Misc : 1.06G/5ML/MSVOA_Y/SOIL
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
191-0922-XYL-1

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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
1-Butanol	8.899	27.0	ug/l	332092	2	8.697	615905	50.0
Isobutyl acetate	10.368	14.5	ug/l	212679	3	11.496	733521	50.0
2-Pentanol, 4-m...	10.441	111.7	ug/l	1639230	3	11.496	733521	50.0
3-Pentanone, 2,...	10.807	18.5	ug/l	271624	3	11.496	733521	50.0
Acetic acid, bu...	10.935	50.5	ug/l	740952	3	11.496	733521	50.0
2-Heptanone	12.105	22.3	ug/l	326693	3	11.496	733521	50.0
Benzene, 1-ethy...	12.733	26.9	ug/l	384251	4	13.428	714440	50.0
Benzene, 1-ethy...	12.971	25.5	ug/l	364658	4	13.428	714440	50.0
Octanal	13.270	16.0	ug/l	228753	4	13.428	714440	50.0
1-Hexanol, 2-et...	13.532	14.6	ug/l	207915	4	13.428	714440	50.0