

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU081420\
 Data File : VU039874.D
 Acq On : 14 Aug 2020 21:44
 Operator : SY/MD
 Sample : L3517-15
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 F4L72

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.1

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_U\METHOD\SOMUTR080320WMA.M

Title : TRACE VOA SOM01.0

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.501	45	50	60	rBV	8456	11492	3.47%	0.303%
2	1.617	76	86	96	rBV3	31912	49020	14.79%	1.294%
3	1.951	182	190	199	rBV	28396	37286	11.25%	0.984%
4	2.594	380	390	402	rBV	57363	94097	28.39%	2.484%
5	2.822	457	461	468	rBV6	2781	3737	1.13%	0.099%
6	3.436	632	652	669	rVB2	43286	118004	35.60%	3.115%
7	4.732	1041	1055	1063	rBV4	15114	36894	11.13%	0.974%
8	4.803	1067	1077	1078	rBV4	9610	14437	4.36%	0.381%
9	5.128	1163	1178	1192	rBV2	50420	108110	32.62%	2.854%
10	5.243	1205	1214	1229	rVB3	11038	21570	6.51%	0.569%
11	5.774	1362	1379	1389	rBV2	104904	245406	74.04%	6.478%
12	6.292	1528	1540	1552	rVV	92079	170095	51.32%	4.490%
13	6.359	1554	1561	1574	rVB	11125	17836	5.38%	0.471%
14	6.578	1616	1629	1642	rBV7	45411	97593	29.44%	2.576%
15	6.729	1665	1676	1689	rVB	64518	120775	36.44%	3.188%
16	6.819	1695	1704	1725	rVB3	10979	30766	9.28%	0.812%
17	7.591	1934	1944	1959	rBV4	31059	54356	16.40%	1.435%
18	7.771	1990	2000	2007	rBV5	4180	7284	2.20%	0.192%
19	7.832	2010	2019	2033	rBV2	14868	26411	7.97%	0.697%
20	7.919	2035	2046	2056	rBV	111360	184278	55.60%	4.865%
21	7.989	2061	2068	2078	rVB2	33557	57965	17.49%	1.530%
22	8.198	2124	2133	2151	rVB2	23006	40488	12.22%	1.069%
23	8.285	2151	2160	2167	rBV6	5533	10575	3.19%	0.279%
24	8.449	2203	2211	2216	rBV7	2754	4129	1.25%	0.109%
25	8.565	2236	2247	2258	rVB2	27986	45511	13.73%	1.201%
26	8.661	2258	2277	2291	rBV	181983	331451	100.00%	8.750%
27	8.806	2312	2322	2352	rBV5	17969	46683	14.08%	1.232%
28	9.427	2503	2515	2538	rVB	140973	232987	70.29%	6.150%
29	9.578	2551	2562	2571	rBV	24122	36860	11.12%	0.973%
30	9.697	2587	2599	2612	rBV	79779	127223	38.38%	3.358%
31	10.105	2714	2726	2749	rVB2	23578	42845	12.93%	1.131%
32	10.240	2756	2768	2777	rBV2	15451	27269	8.23%	0.720%
33	10.375	2791	2810	2822	rVB7	6862	20419	6.16%	0.539%
34	10.571	2858	2871	2873	rBV4	4401	5541	1.67%	0.146%

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35	10.761	2918	2930	2945	rVB	61612	97525	29.42%	2.574%
36	10.996	2994	3003	3009	rBV4	5197	8309	2.51%	0.219%
37	11.054	3016	3021	3028	rBV4	3819	5043	1.52%	0.133%
38	11.208	3058	3069	3077	rBV8	6852	12379	3.73%	0.327%
39	11.256	3077	3084	3092	rVV3	12224	20415	6.16%	0.539%
40	11.288	3092	3094	3105	rVB5	4298	4731	1.43%	0.125%
41	11.417	3118	3134	3135	rBV4	3625	5130	1.55%	0.135%
42	11.468	3141	3150	3159	rBV2	51212	73619	22.21%	1.943%
43	11.578	3172	3184	3197	rBV3	98581	192515	58.08%	5.082%
44	11.648	3197	3206	3216	rVV2	24535	37364	11.27%	0.986%
45	11.748	3226	3237	3245	rBV2	5315	10542	3.18%	0.278%
46	11.809	3245	3256	3269	rVB	161254	250869	75.69%	6.622%
47	11.886	3272	3280	3283	rBV5	3509	4534	1.37%	0.120%
48	11.963	3295	3304	3317	rVB6	16290	27897	8.42%	0.736%
49	12.053	3322	3332	3340	rVV5	11440	19808	5.98%	0.523%
50	12.086	3340	3342	3353	rVB5	2680	3524	1.06%	0.093%
51	12.189	3362	3374	3391	rBV	138502	223461	67.42%	5.899%
52	12.343	3404	3422	3438	rVB7	24727	46611	14.06%	1.230%
53	12.491	3448	3468	3471	rBV10	5971	14112	4.26%	0.373%
54	12.671	3515	3524	3541	rVB9	6708	13299	4.01%	0.351%
55	12.767	3541	3554	3560	rBV	57127	88966	26.84%	2.349%
56	12.803	3560	3565	3580	rVV2	38469	60553	18.27%	1.598%
57	12.880	3582	3589	3598	rVB10	6745	10055	3.03%	0.265%
58	13.279	3708	3713	3722	rVB7	5228	7331	2.21%	0.194%
59	13.394	3739	3749	3756	rBV9	2770	4734	1.43%	0.125%
60	13.462	3763	3770	3777	rVB8	3216	4882	1.47%	0.129%
61	13.619	3806	3819	3827	rBV7	3385	7186	2.17%	0.190%
62	13.864	3882	3895	3907	rBV5	15028	30238	9.12%	0.798%
63	14.536	4094	4104	4112	rBV5	3182	5868	1.77%	0.155%
64	14.632	4121	4134	4143	rBV10	3104	6187	1.87%	0.163%
65	14.870	4198	4208	4217	rBV4	7392	11069	3.34%	0.292%

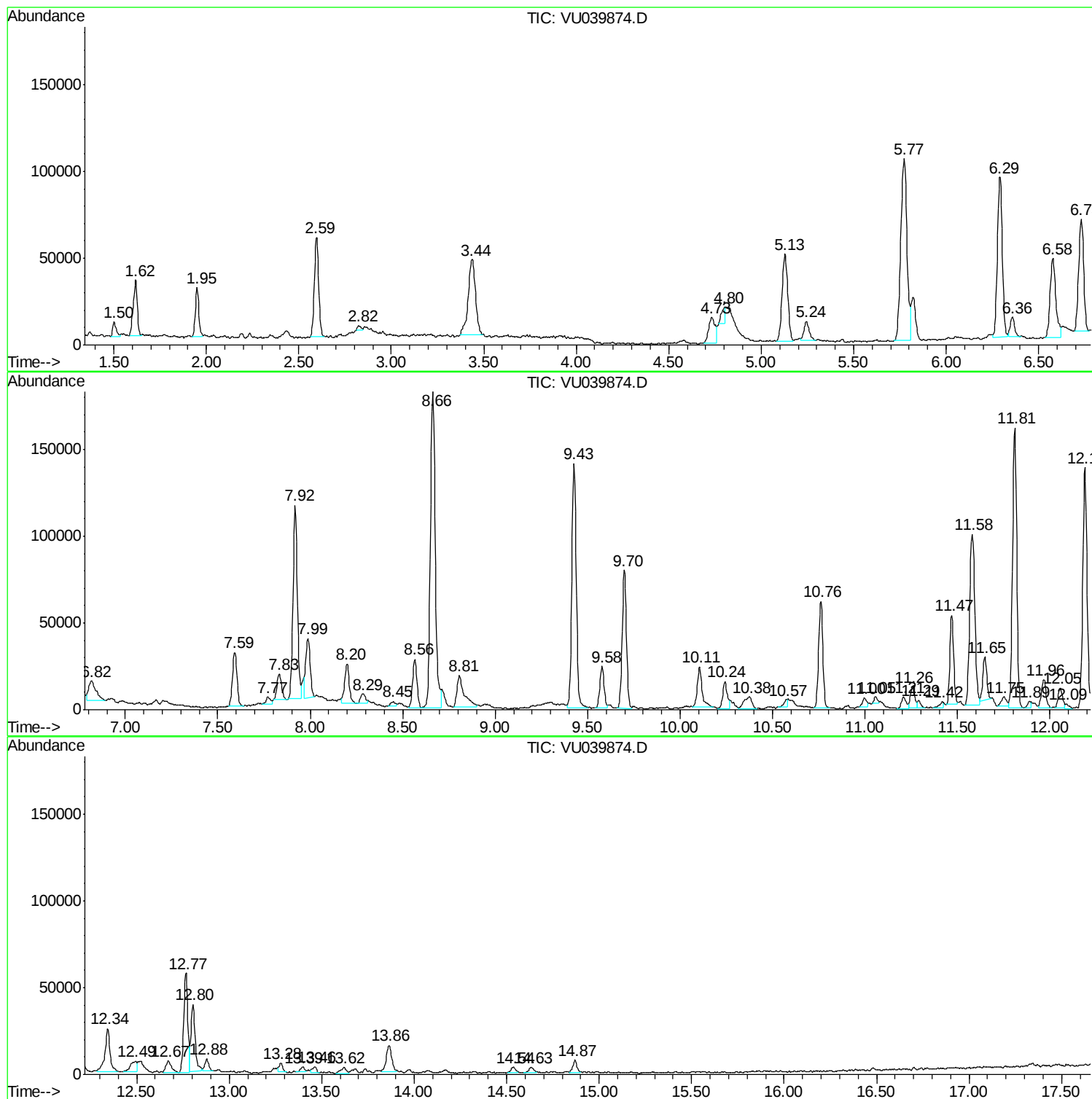
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR080320WMA.M
Quant Title : TRACE VOA SOM01.0

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



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Instrument :
MSVOA_U
ClientSampleId :
F4L72

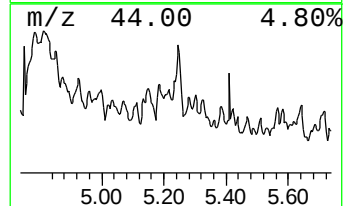
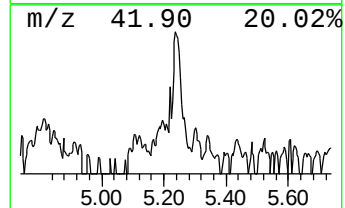
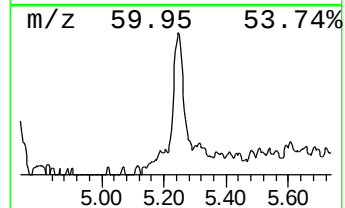
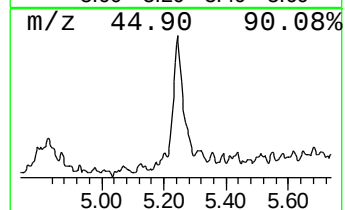
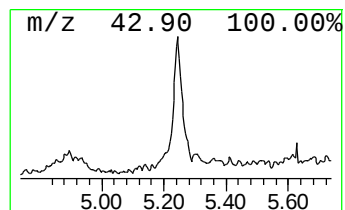
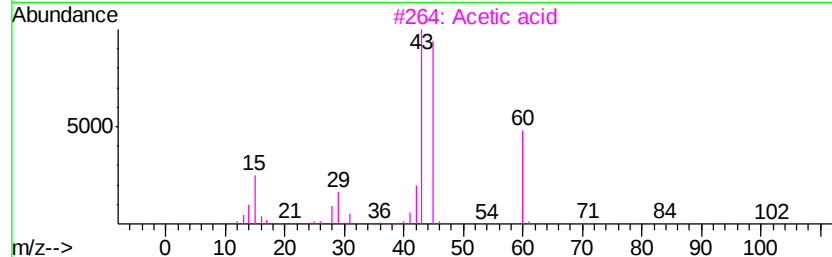
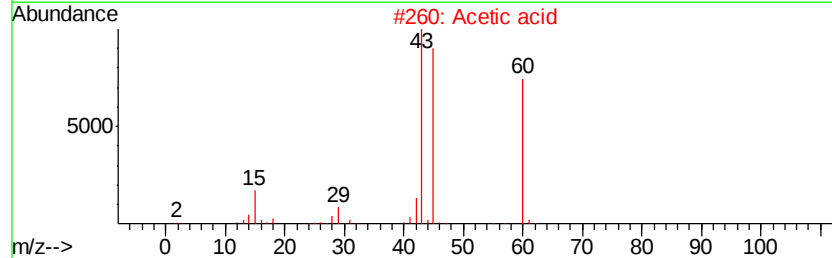
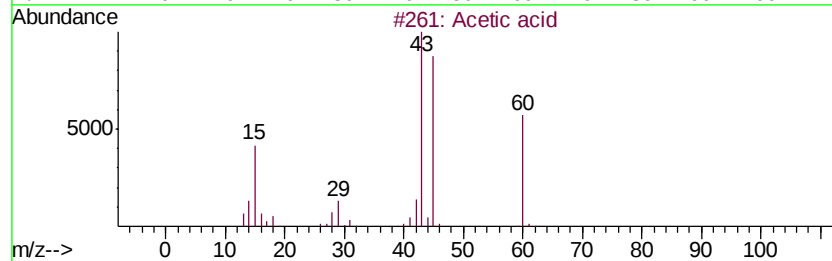
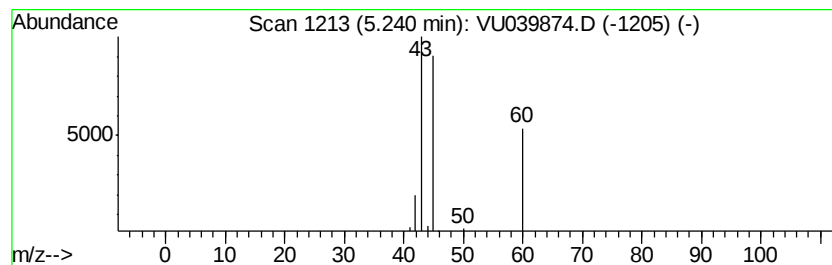
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR080320WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 1 Acetic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.24	0.63 ug/L	21570	1,4-Difluorobenzene	6.29

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetic acid			60	C2H4O2	000064-19-7	90
2	Acetic acid			60	C2H4O2	000064-19-7	83
3	Acetic acid			60	C2H4O2	000064-19-7	64
4	Acetic acid			60	C2H4O2	000064-19-7	9
5	Hydrazine, 1,2-dimethyl-			60	C2H8N2	000540-73-8	5



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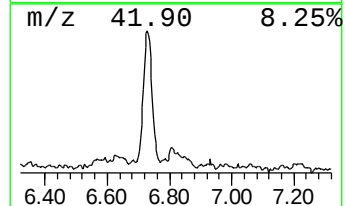
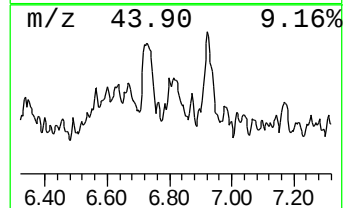
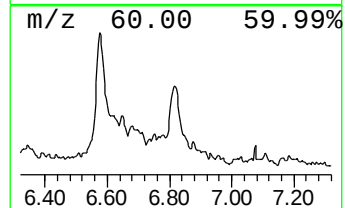
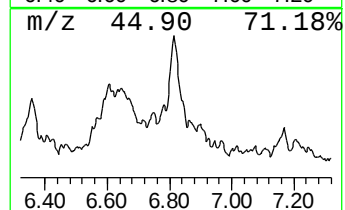
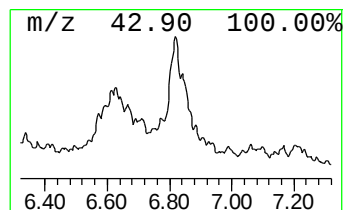
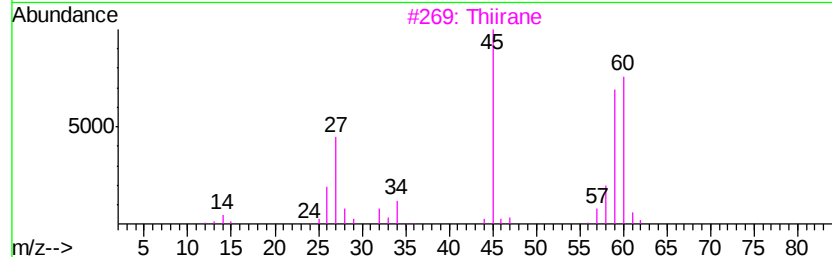
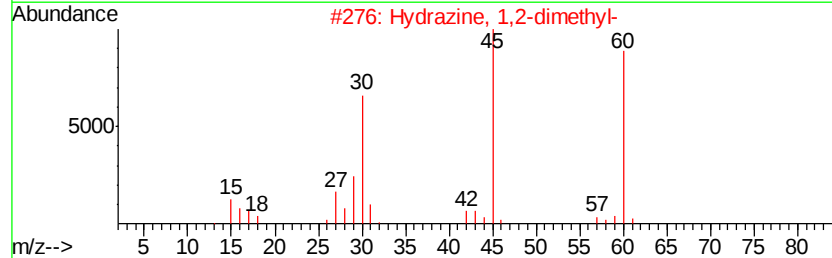
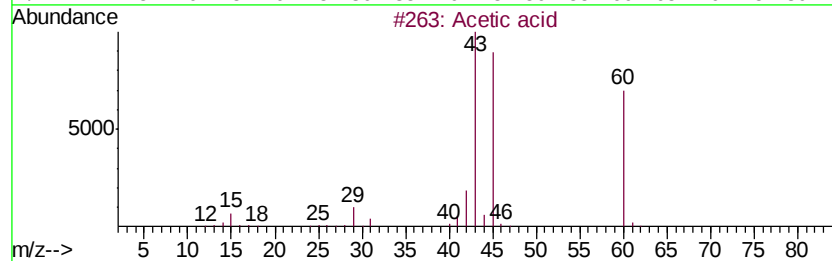
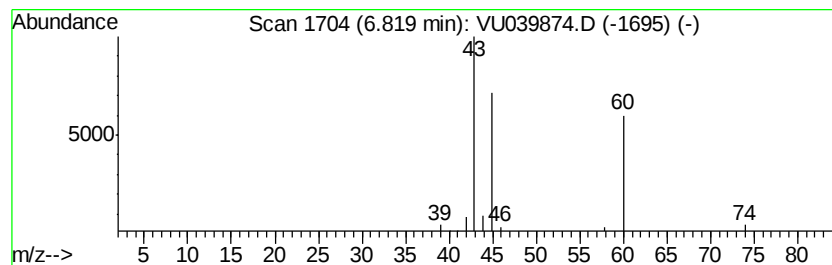
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR080320WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 3 Hydrazine, 1,2-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.82	0.90 ug/L	30766	1,4-Difluorobenzene	6.29

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetic acid		60	C2H4O2	000064-19-7	64
2	Hydrazine, 1,2-dimethyl-		60	C2H8N2	000540-73-8	53
3	Thiirane		60	C2H4S	000420-12-2	9
4	Hydrazine, 1,2-dimethyl-		60	C2H8N2	000540-73-8	9
5	Acetic acid		60	C2H4O2	000064-19-7	9



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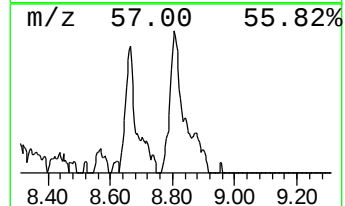
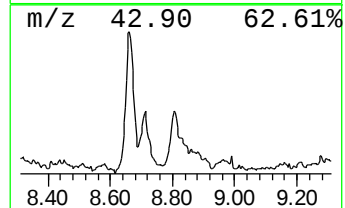
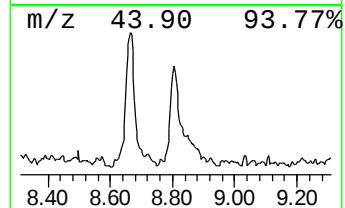
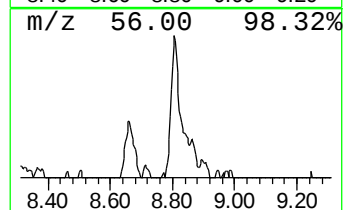
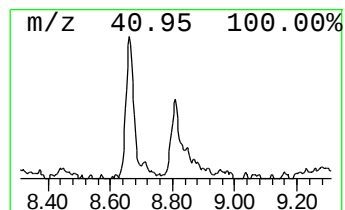
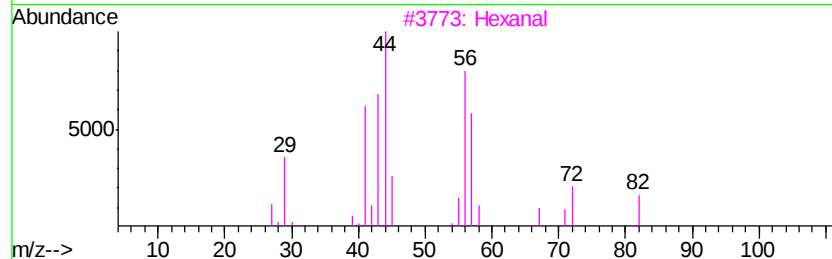
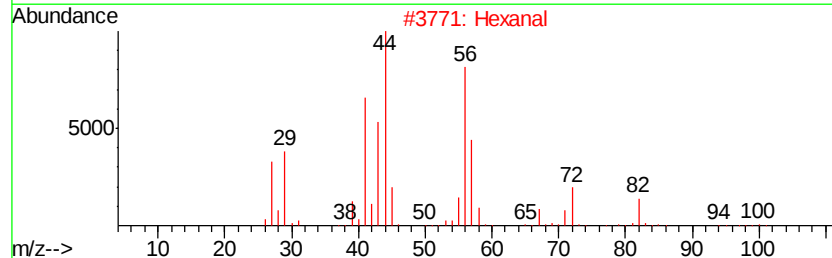
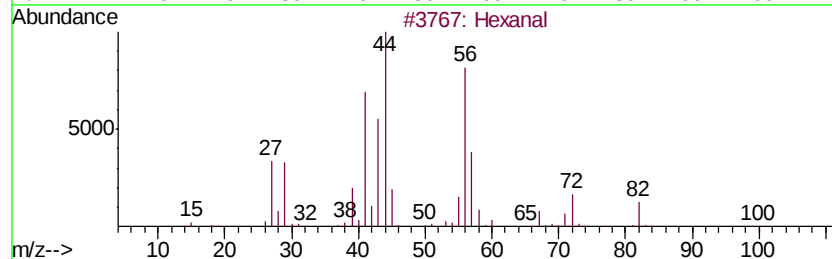
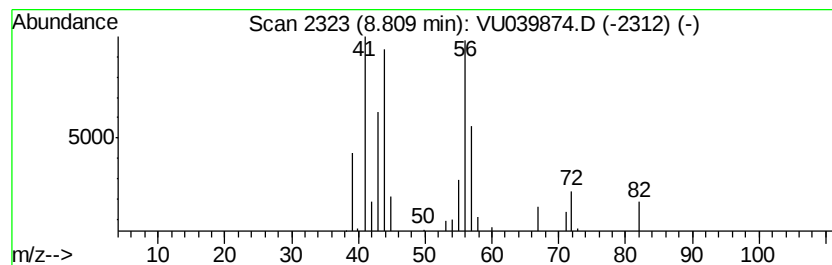
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Hexanal Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.81	1.00 ug/L	46683	Chlorobenzene-d5	9.43

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanal		100	C6H12O	000066-25-1	80
2	Hexanal		100	C6H12O	000066-25-1	72
3	Hexanal		100	C6H12O	000066-25-1	72
4	Hexanal		100	C6H12O	000066-25-1	59
5	Cyclobutanol, 2-ethyl-		100	C6H12O	035301-43-0	37



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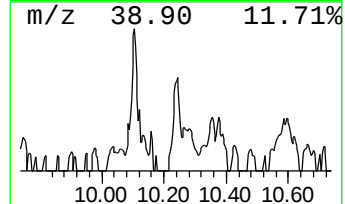
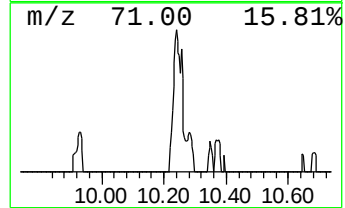
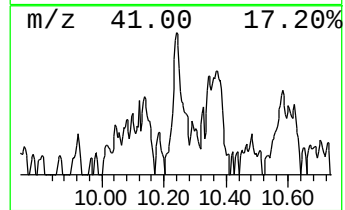
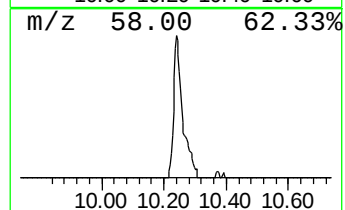
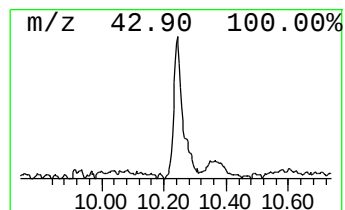
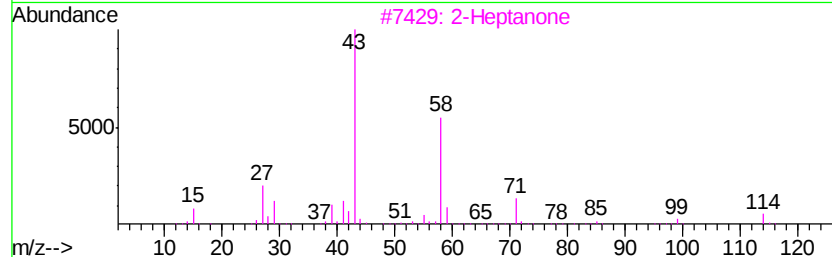
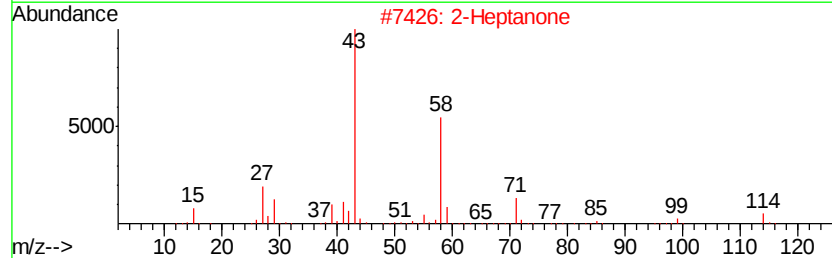
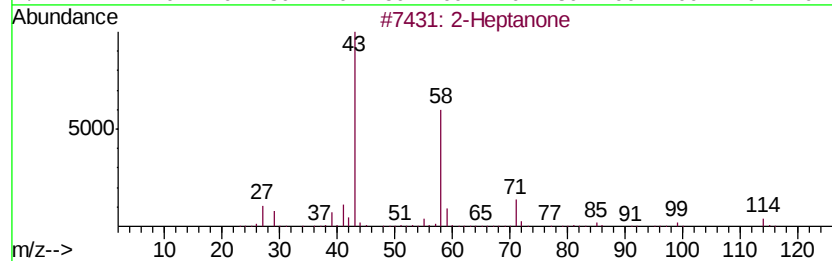
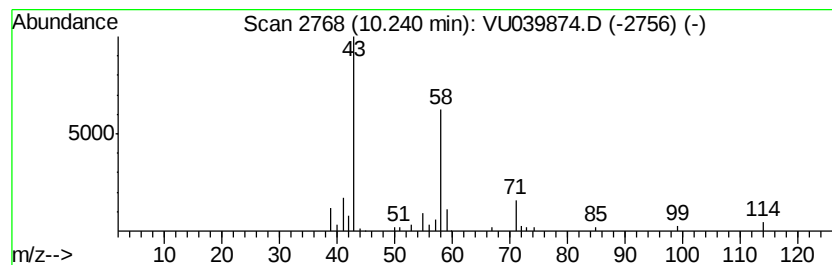
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 6 2-Heptanone Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.24	0.59 ug/L	27269	Chlorobenzene-d5	9.43	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Heptanone	114	C7H14O	000110-43-0	91
2	2-Heptanone	114	C7H14O	000110-43-0	90
3	2-Heptanone	114	C7H14O	000110-43-0	87
4	2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	86
5	2-Heptanone	114	C7H14O	000110-43-0	83



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ClientSampleId :
F4L72

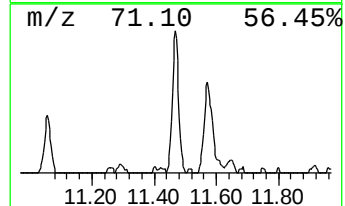
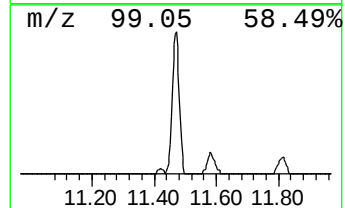
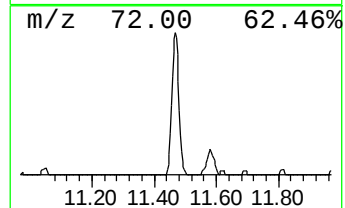
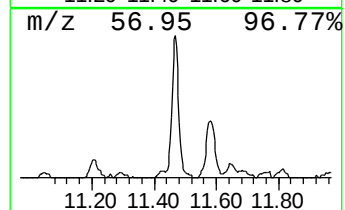
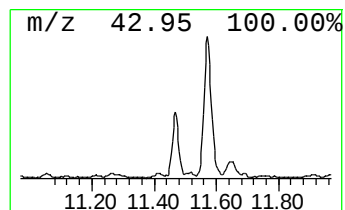
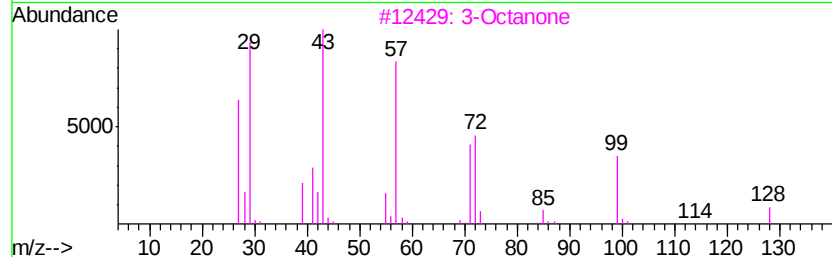
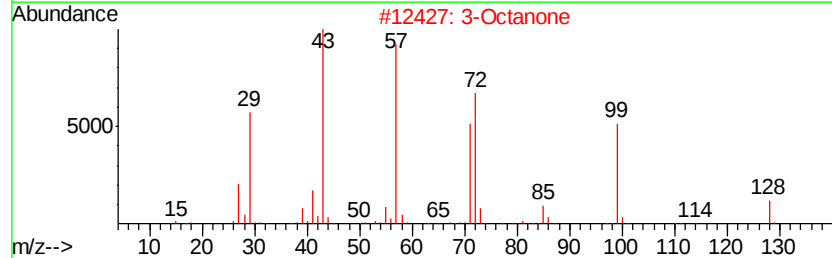
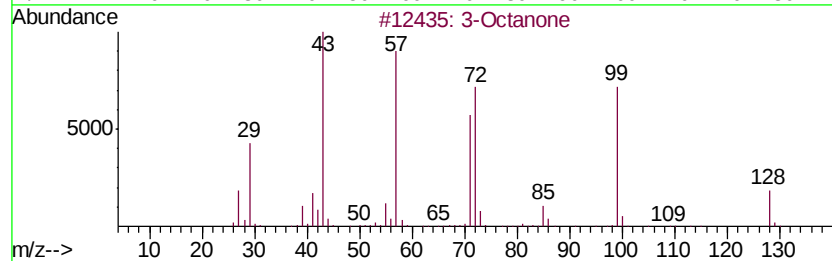
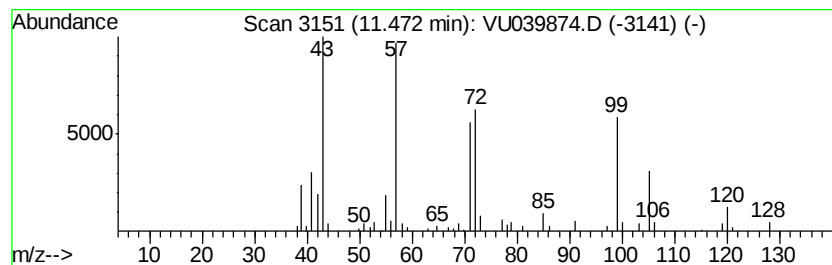
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 7 3-Octanone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.47	1.47 ug/L	73619	1,4-Dichlorobenzene-d4	11.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Octanone	128	C8H16O	000106-68-3	81
2			3-Octanone	128	C8H16O	000106-68-3	81
3			3-Octanone	128	C8H16O	000106-68-3	59
4			3-Heptanone, 5-methyl-	128	C8H16O	000541-85-5	53
5			3-Heptanone, 5-methyl-	128	C8H16O	000541-85-5	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU081420\
Data File : VU039874.D
Acq On : 14 Aug 2020 21:44
Operator : SY/MD
Sample : L3517-15
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
F4L72

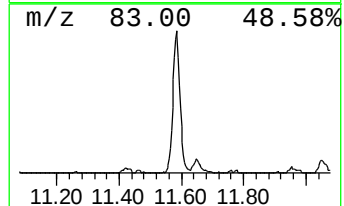
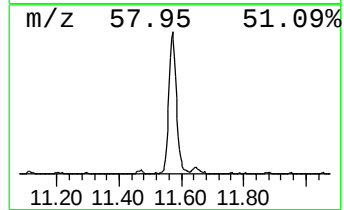
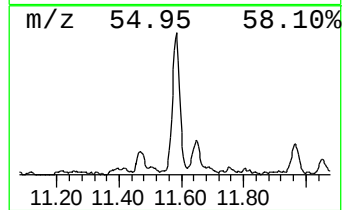
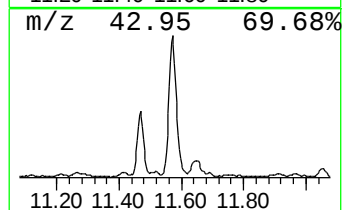
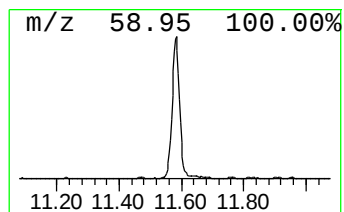
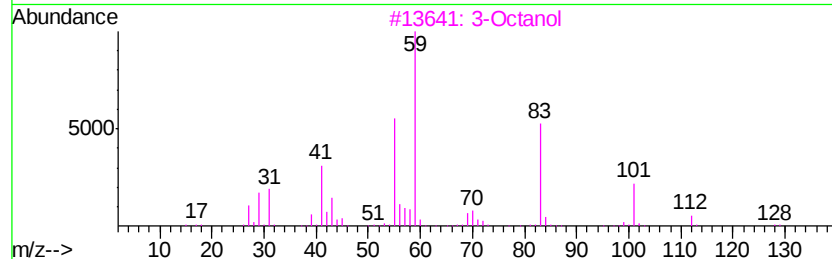
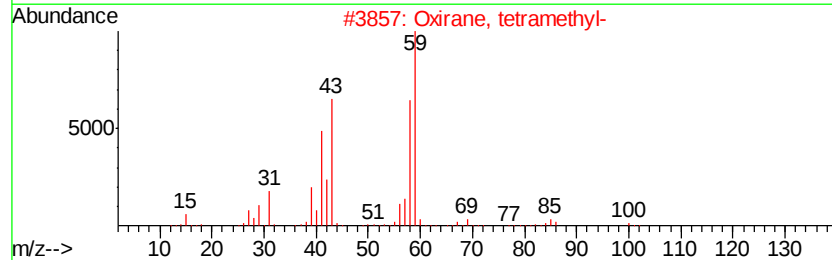
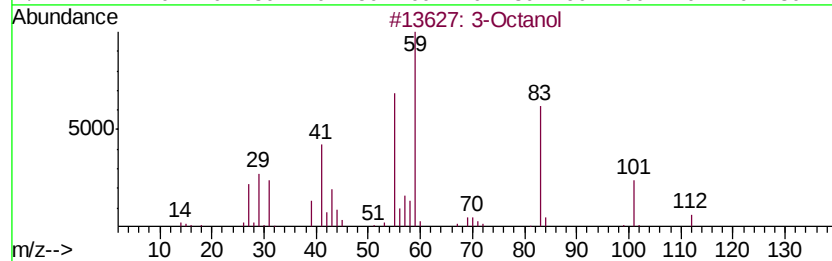
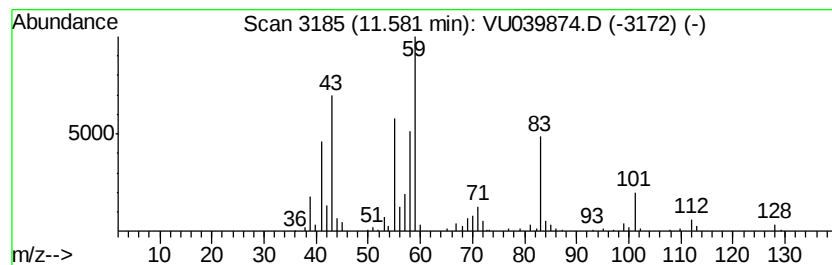
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 8 3-Octanol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.58	3.84 ug/L	192515	1,4-Dichlorobenzene-d4	11.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Octanol	130	C8H18O	000589-98-0	59
2		Oxirane, tetramethyl-	100	C6H12O	005076-20-0	53
3		3-Octanol	130	C8H18O	000589-98-0	53
4		5-Ethyl-3-methylhept-1-en-4-ol	156	C10H20O	286424-80-4	53
5		3-Octanol	130	C8H18O	000589-98-0	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU081420\
Data File : VU039874.D
Acq On : 14 Aug 2020 21:44
Operator : SY/MD
Sample : L3517-15
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
F4L72

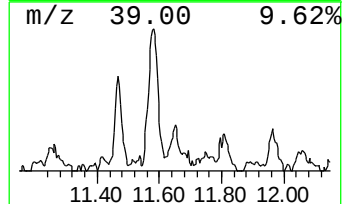
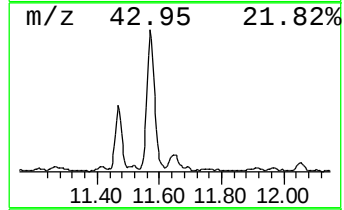
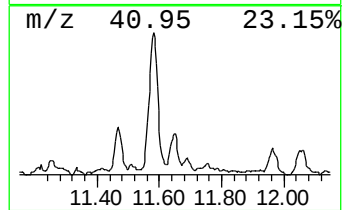
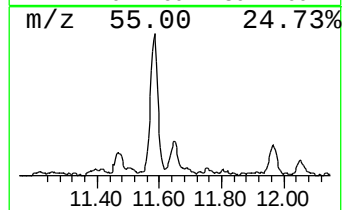
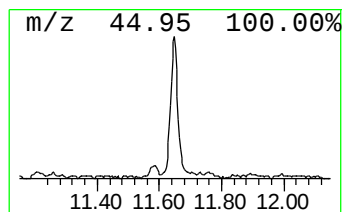
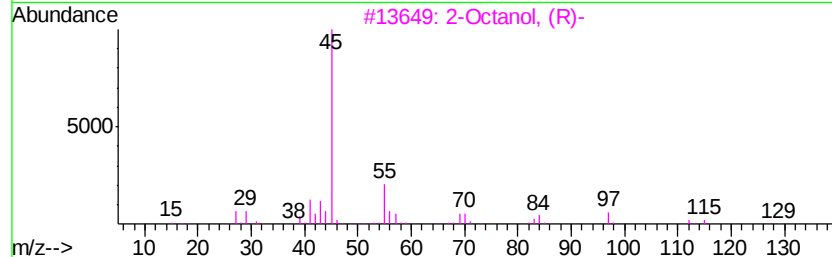
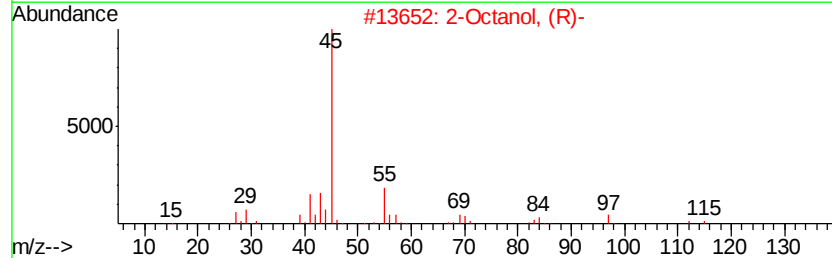
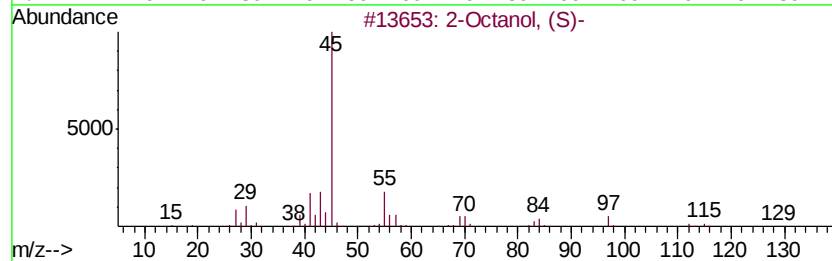
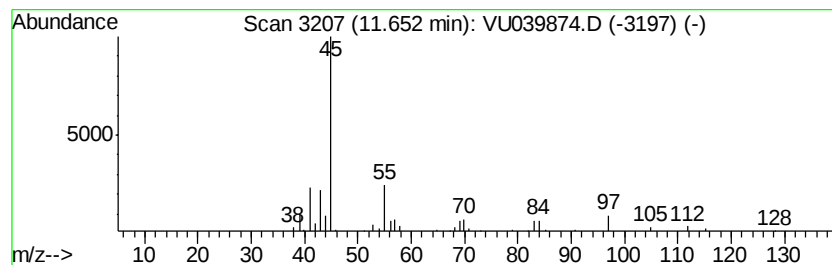
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 9 2-Octanol, (S)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.65	0.74 ug/L	37364	1,4-Dichlorobenzene-d4	11.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Octanol, (S)-	130	C8H18O	006169-06-8	83
2		2-Octanol, (R)-	130	C8H18O	005978-70-1	83
3		2-Octanol, (R)-	130	C8H18O	005978-70-1	83
4		2-Octanol	130	C8H18O	000123-96-6	83
5		2-Octanol, (S)-	130	C8H18O	006169-06-8	78



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU081420\
Data File : VU039874.D
Acq On : 14 Aug 2020 21:44
Operator : SY/MD
Sample : L3517-15
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

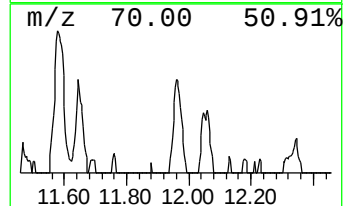
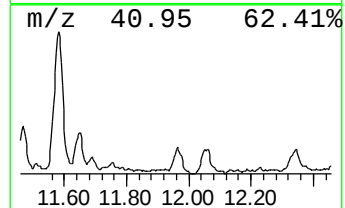
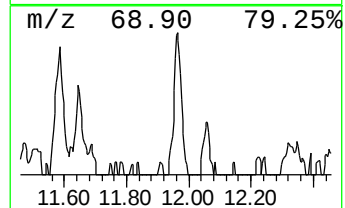
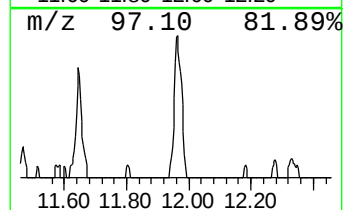
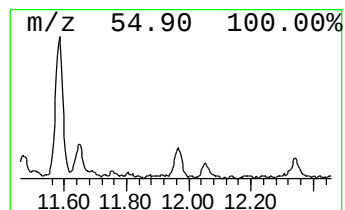
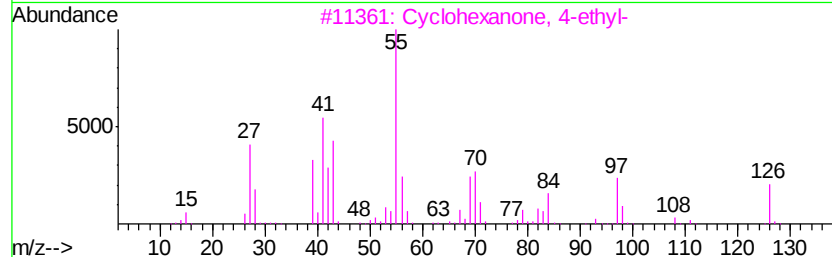
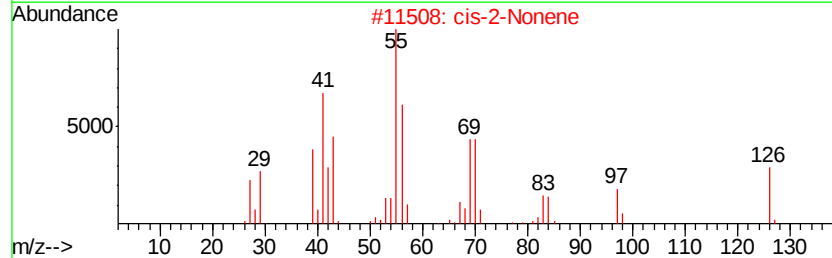
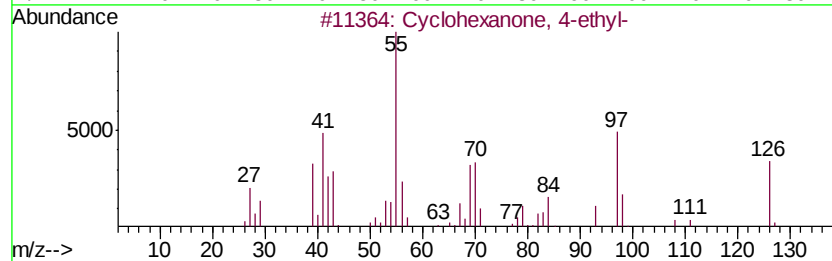
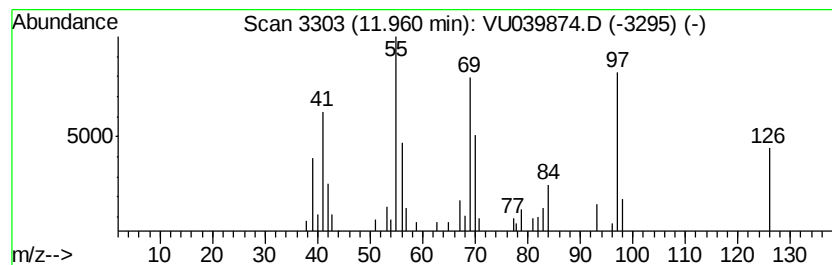
Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR080320WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 10 Cyclohexanone, 4-ethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.96	0.56 ug/L	27897	1,4-Dichlorobenzene-d4	11.81
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclohexanone, 4-ethyl-	126 C8H14O	005441-51-0 70
2		cis-2-Nonene	126 C9H18	006434-77-1 53
3		Cyclohexanone, 4-ethyl-	126 C8H14O	005441-51-0 49
4		1-Pentene, 3-methyl-	84 C6H12	000760-20-3 46
5		2-Heptenal, 2-methyl-	126 C8H14O	030567-26-1 43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU081420\
Data File : VU039874.D
Acq On : 14 Aug 2020 21:44
Operator : SY/MD
Sample : L3517-15
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

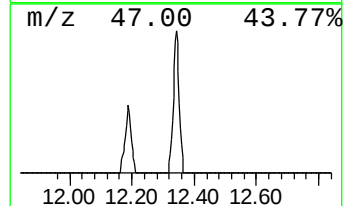
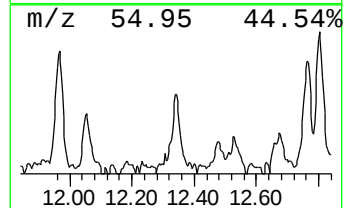
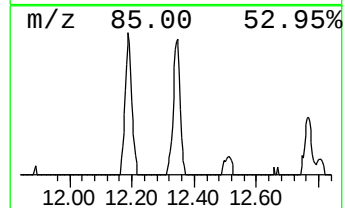
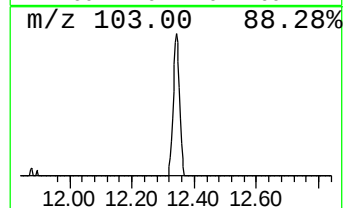
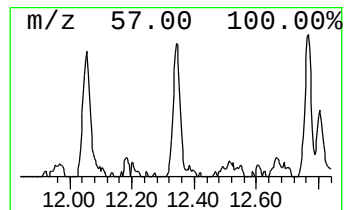
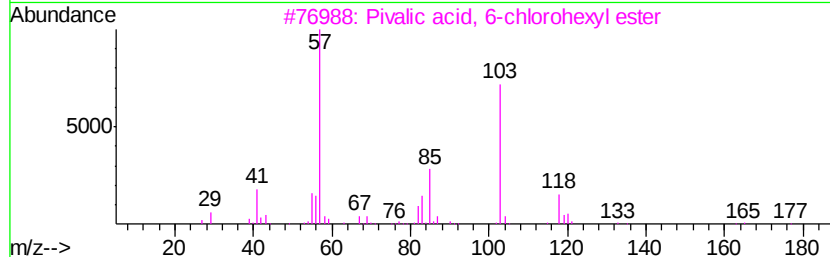
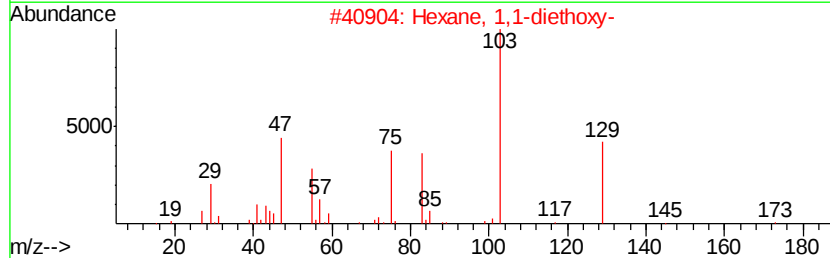
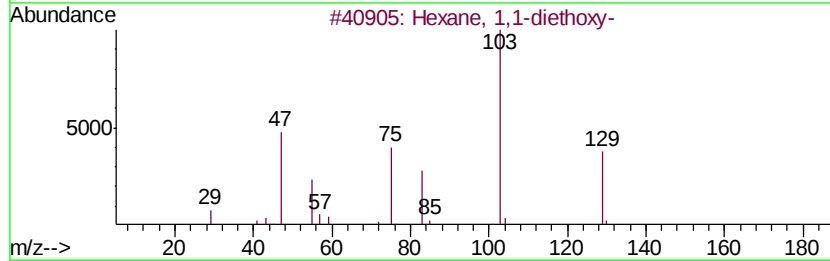
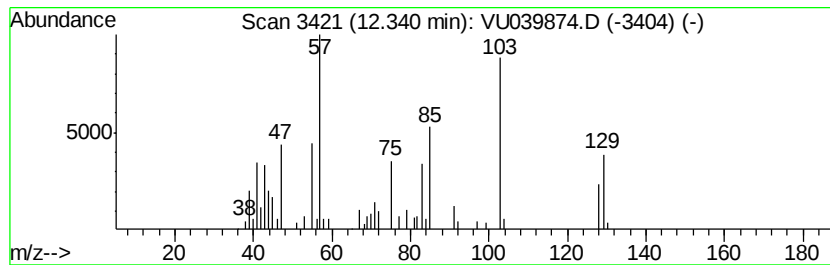
Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR080320WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 11 Hexane, 1,1-diethoxy- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.34	0.93 ug/L	46611	1,4-Dichlorobenzene-d4	11.81
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexane, 1,1-diethoxy-	174 C10H22O2	003658-93-3	53
2	Hexane, 1,1-diethoxy-	174 C10H22O2	003658-93-3	53
3	Pivalic acid, 6-chlorohexyl ester	220 C11H21ClO2	1000330-91-8	37
4	Butanoic acid, 2-methyl-, hexyl ...	186 C11H22O2	010032-15-2	35
5	Butanoic acid, 2-methyl-, pentyl...	172 C10H20O2	068039-26-9	35



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU081420\
Data File : VU039874.D
Acq On : 14 Aug 2020 21:44
Operator : SY/MD
Sample : L3517-15
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
F4L72

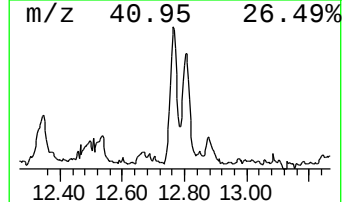
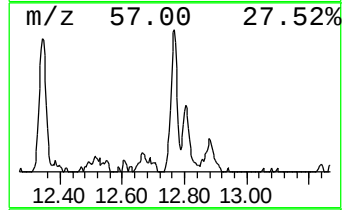
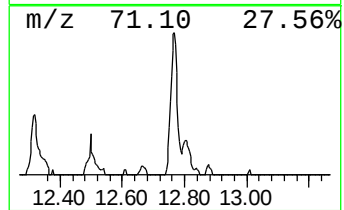
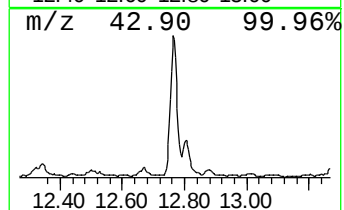
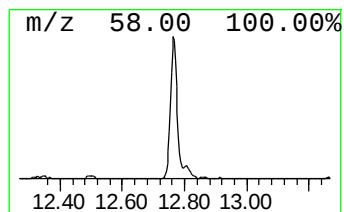
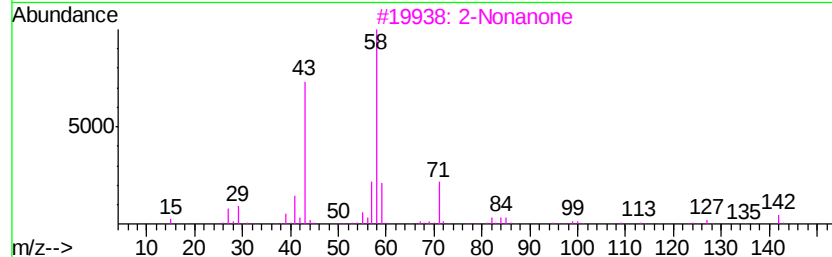
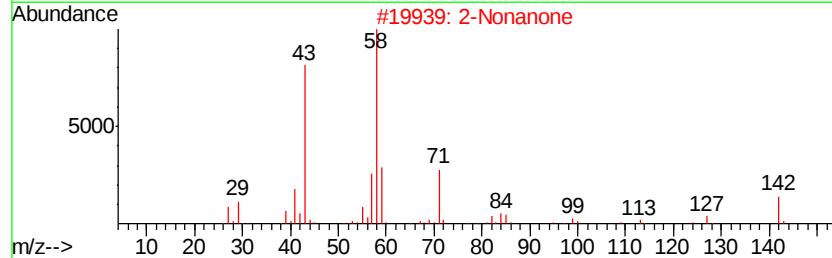
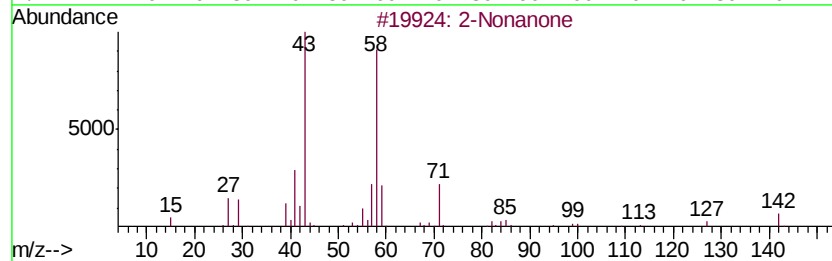
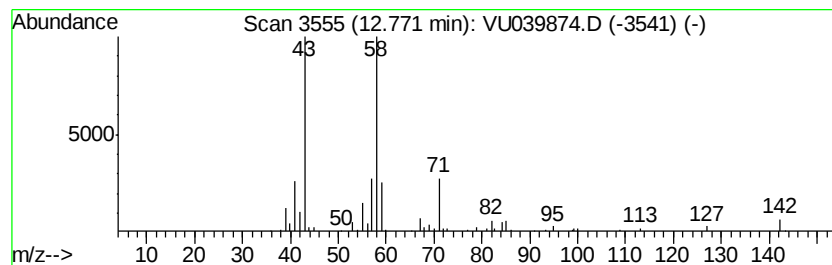
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 12 2-Nonanone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.77	1.77 ug/L	88966	1,4-Dichlorobenzene-d4	11.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Nonanone			142	C9H18O	000821-55-6	94
2	2-Nonanone			142	C9H18O	000821-55-6	91
3	2-Nonanone			142	C9H18O	000821-55-6	87
4	2-Nonanone			142	C9H18O	000821-55-6	80
5	2-Hexanone, 5-methyl-			114	C7H14O	000110-12-3	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU081420\
Data File : VU039874.D
Acq On : 14 Aug 2020 21:44
Operator : SY/MD
Sample : L3517-15
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

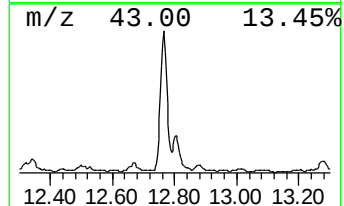
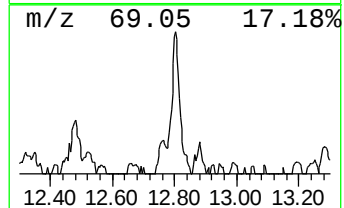
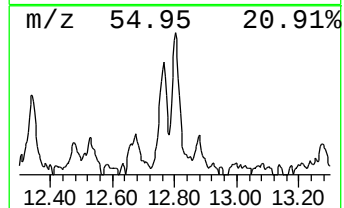
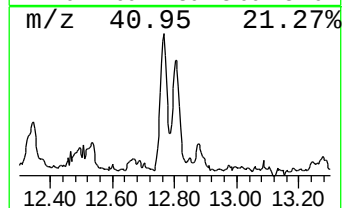
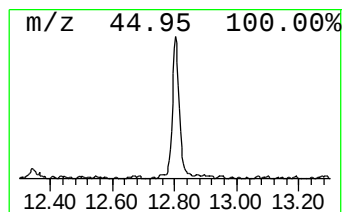
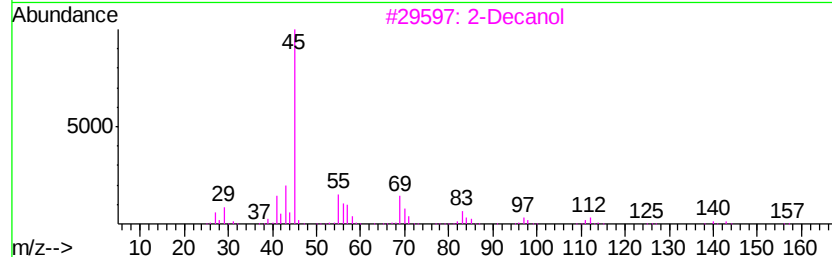
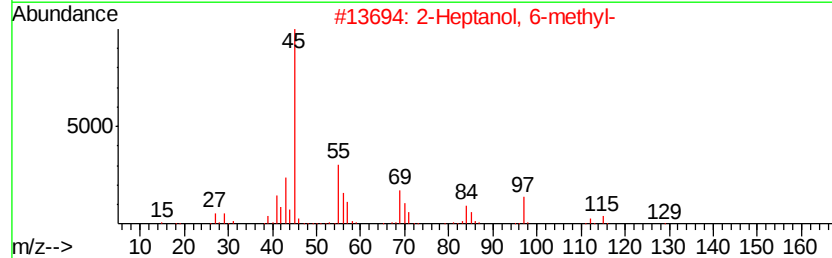
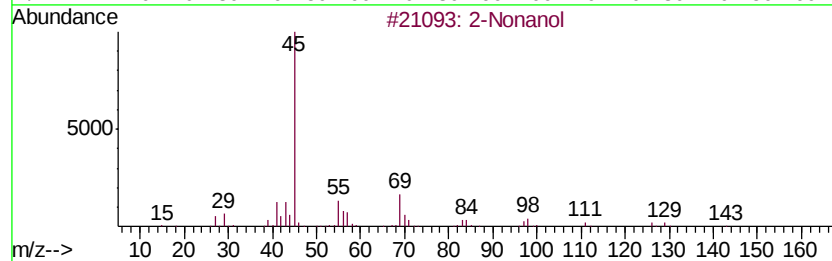
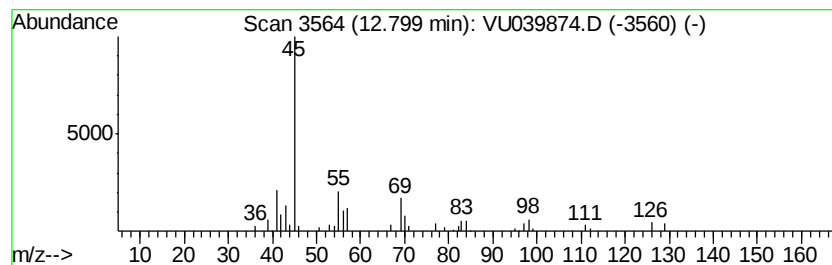
Instrument :
MSVOA_U
ClientSampleId :
F4L72

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR080320WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 13 2-Nonanol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.80	1.21 ug/L	60553	1,4-Dichlorobenzene-d4		11.81
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Nonanol	144	C9H20O	000628-99-9	83
2	2-Heptanol, 6-methyl-	130	C8H18O	004730-22-7	78
3	2-Decanol	158	C10H22O	001120-06-5	78
4	2-Nonanol	144	C9H20O	000628-99-9	78
5	2-Decanol	158	C10H22O	001120-06-5	78



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU081420\
Data File : VU039874.D
Acq On : 14 Aug 2020 21:44
Operator : SY/MD
Sample : L3517-15
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
F4L72

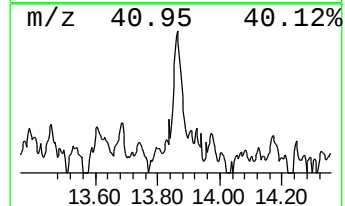
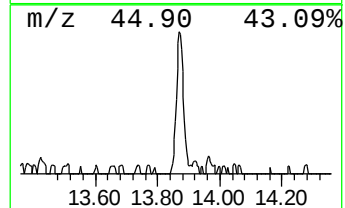
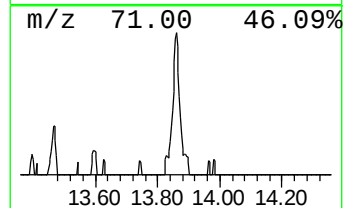
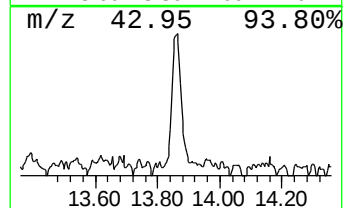
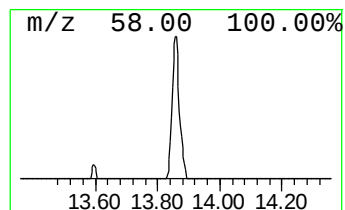
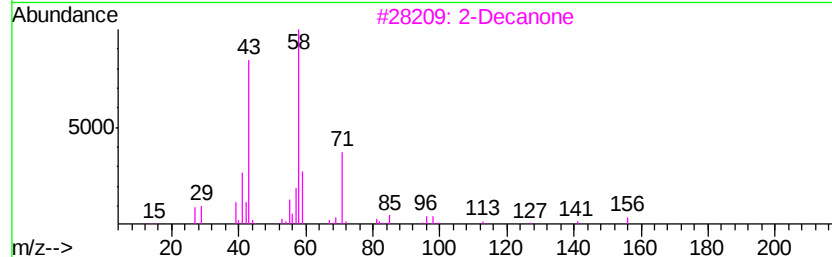
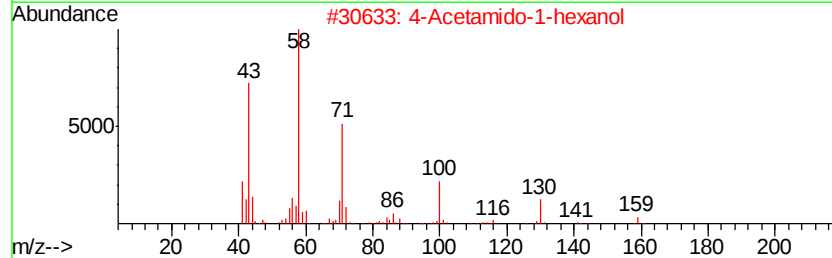
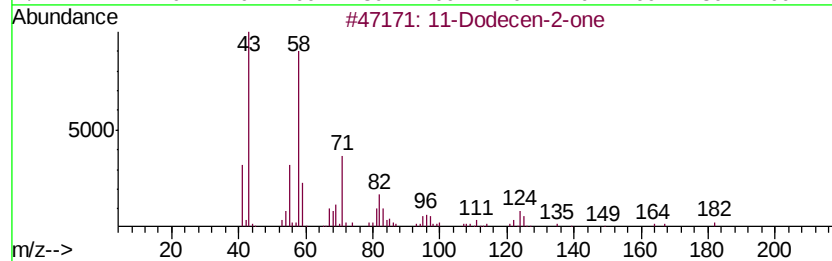
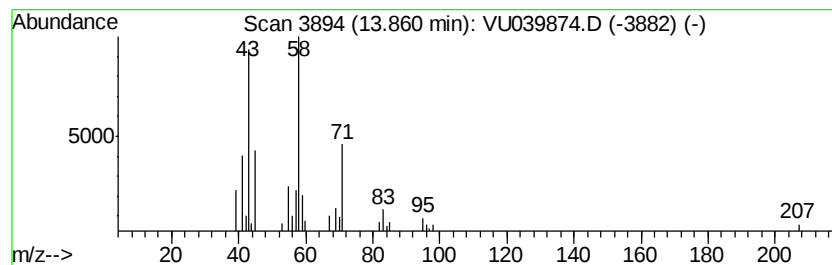
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 14 11-Dodecen-2-one Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.86	0.60 ug/L	30238	1,4-Dichlorobenzene-d4	11.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11-Dodecen-2-one	182	C12H22O	005009-33-6	53
2		4-Acetamido-1-hexanol	159	C8H17NO2	1000213-28-5	47
3		2-Decanone	156	C10H20O	000693-54-9	38
4		3-Oxetanol, 2,2,3-trimethyl-	116	C6H12O2	025910-96-7	38
5		2-Nonadecanol	284	C19H40O	026533-36-8	22



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU081420\
 Data File : VU039874.D
 Acq On : 14 Aug 2020 21:44
 Operator : SY/MD
 Sample : L3517-15
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 F4L72

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR080320WMA.M
 Quant Title : TRACE VOA SOM01.0

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Acetic acid	5.24	0.6	ug/L	21570	1	6.29	170095	5.0
Hydrazine, 1,2-di...	6.82	0.9	ug/L	30766	1	6.29	170095	5.0
Hexanal	8.81	1.0	ug/L	46683	2	9.43	232987	5.0
2-Heptanone	10.24	0.6	ug/L	27269	2	9.43	232987	5.0
3-Octanone	11.47	1.5	ug/L	73619	3	11.81	250869	5.0
3-Octanol	11.58	3.8	ug/L	192515	3	11.81	250869	5.0
2-Octanol, (S)-	11.65	0.7	ug/L	37364	3	11.81	250869	5.0
Cyclohexanone, 4-...	11.96	0.6	ug/L	27897	3	11.81	250869	5.0
Hexane, 1,1-dieth...	12.34	0.9	ug/L	46611	3	11.81	250869	5.0
2-Nonanone	12.77	1.8	ug/L	88966	3	11.81	250869	5.0
2-Nonanol	12.80	1.2	ug/L	60553	3	11.81	250869	5.0
11-Dodecen-2-one	13.86	0.6	ug/L	30238	3	11.81	250869	5.0