

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV060922\
 Data File : VV026082.D
 Acq On : 10 Jun 2022 02:55
 Operator : SY/MD
 Sample : N3272-19
 Misc : 25mL/MSVOA_V/WATER
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 C0B66

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.1

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_V\Method\SFAMVTR060822WMA.M
 Title : TRACE VOA SFAM1.0

Signal : TIC: VV026082.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.320	79	87	98	rVB	62217	66293	3.08%	0.908%
2	1.584	161	169	182	rVB	70967	82720	3.84%	1.133%
3	2.124	326	337	352	rVB	117070	175864	8.17%	2.409%
4	3.195	654	670	685	rBV2	15899	34404	1.60%	0.471%
5	3.921	886	896	926	rBV	33407	129188	6.00%	1.770%
6	4.362	1018	1033	1054	rBV	101819	246911	11.48%	3.382%
7	5.056	1234	1249	1276	rBV3	192924	490589	22.80%	6.720%
8	5.625	1412	1426	1452	rBV	184581	383074	17.80%	5.248%
9	5.918	1503	1517	1551	rBV	1101955	2151724	100.00%	29.476%
10	6.075	1554	1566	1594	rVB	125784	264498	12.29%	3.623%
11	6.995	1839	1852	1873	rBV	40822	80637	3.75%	1.105%
12	7.320	1941	1953	1973	rBV	200233	357611	16.62%	4.899%
13	7.629	2040	2049	2062	rBV2	22703	41848	1.94%	0.573%
14	7.979	2148	2158	2178	rVB2	65086	119393	5.55%	1.636%
15	8.091	2182	2193	2219	rBV2	190757	510721	23.74%	6.996%
16	8.381	2271	2283	2294	rBV6	22367	48839	2.27%	0.669%
17	8.513	2307	2324	2340	rBV6	29849	76877	3.57%	1.053%
18	8.799	2401	2413	2421	rBV4	32429	62416	2.90%	0.855%
19	8.854	2421	2430	2451	rVB	315185	516100	23.99%	7.070%
20	10.008	2774	2789	2808	rBV	61879	111602	5.19%	1.529%
21	10.217	2844	2854	2871	rVB	91836	151453	7.04%	2.075%
22	10.407	2897	2913	2926	rBV4	17036	35483	1.65%	0.486%
23	10.857	3041	3053	3067	rBV	40554	68678	3.19%	0.941%
24	10.950	3073	3082	3089	rBV	17438	26111	1.21%	0.358%
25	11.098	3115	3128	3153	rBV3	70400	150175	6.98%	2.057%
26	11.249	3163	3175	3186	rBV	291788	464040	21.57%	6.357%
27	11.622	3280	3291	3305	rBV	230024	364487	16.94%	4.993%
28	12.082	3425	3434	3446	rBV	38587	57641	2.68%	0.790%
29	16.178	4696	4708	4719	rBV5	20395	30565	1.42%	0.419%

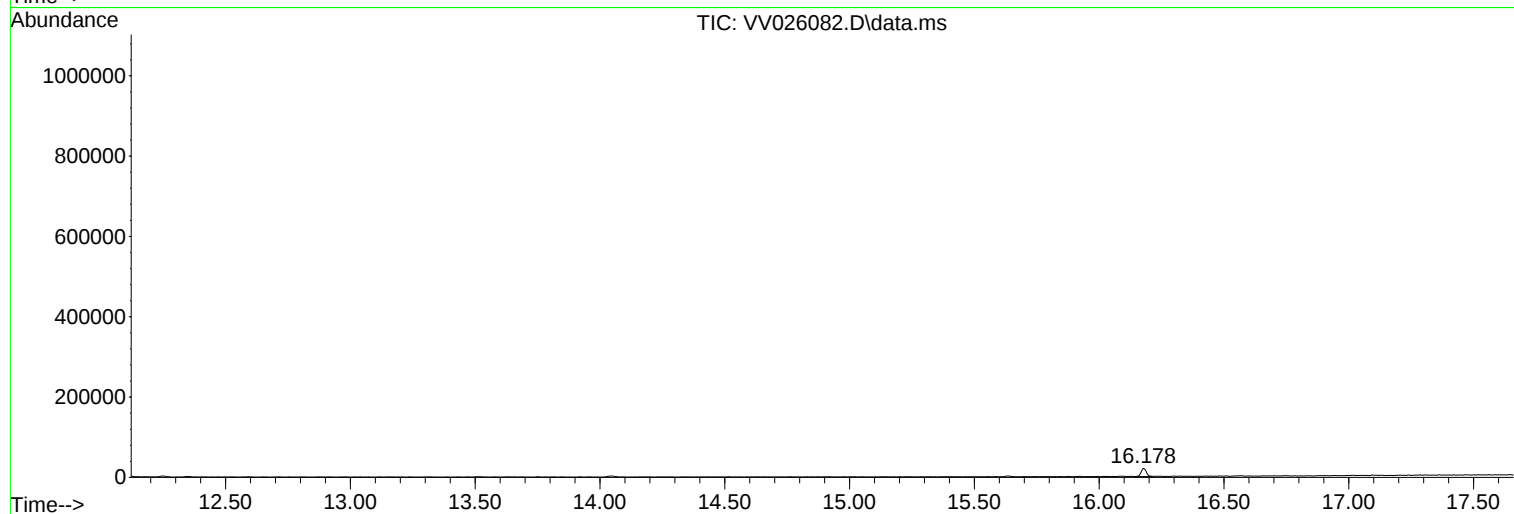
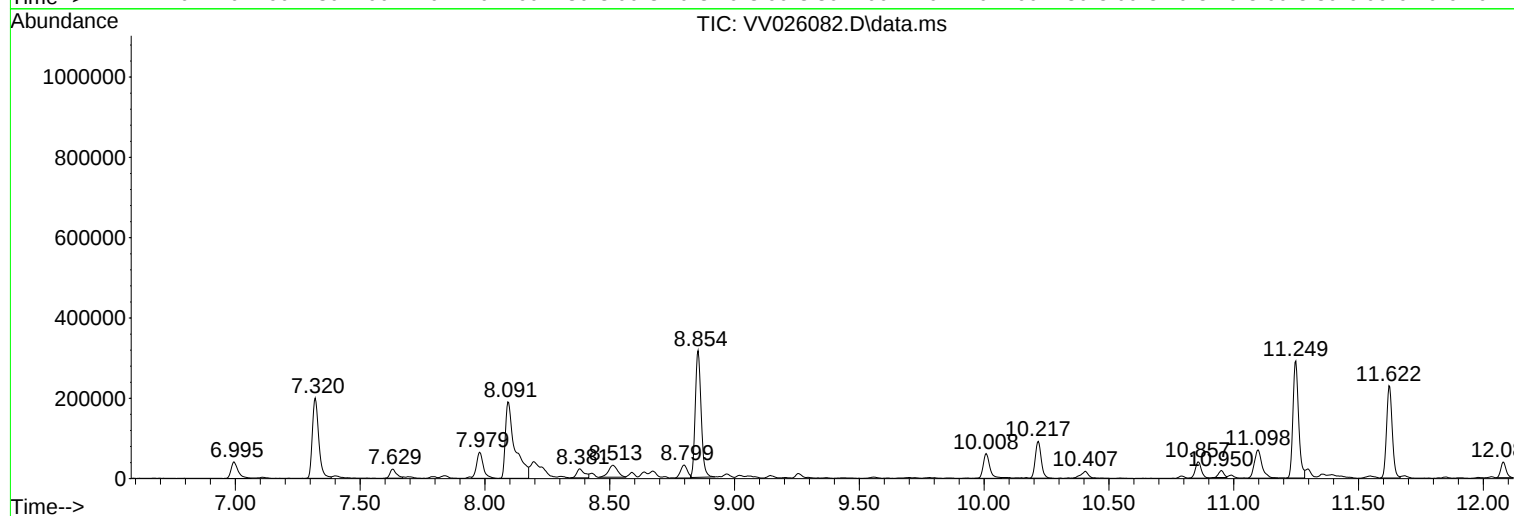
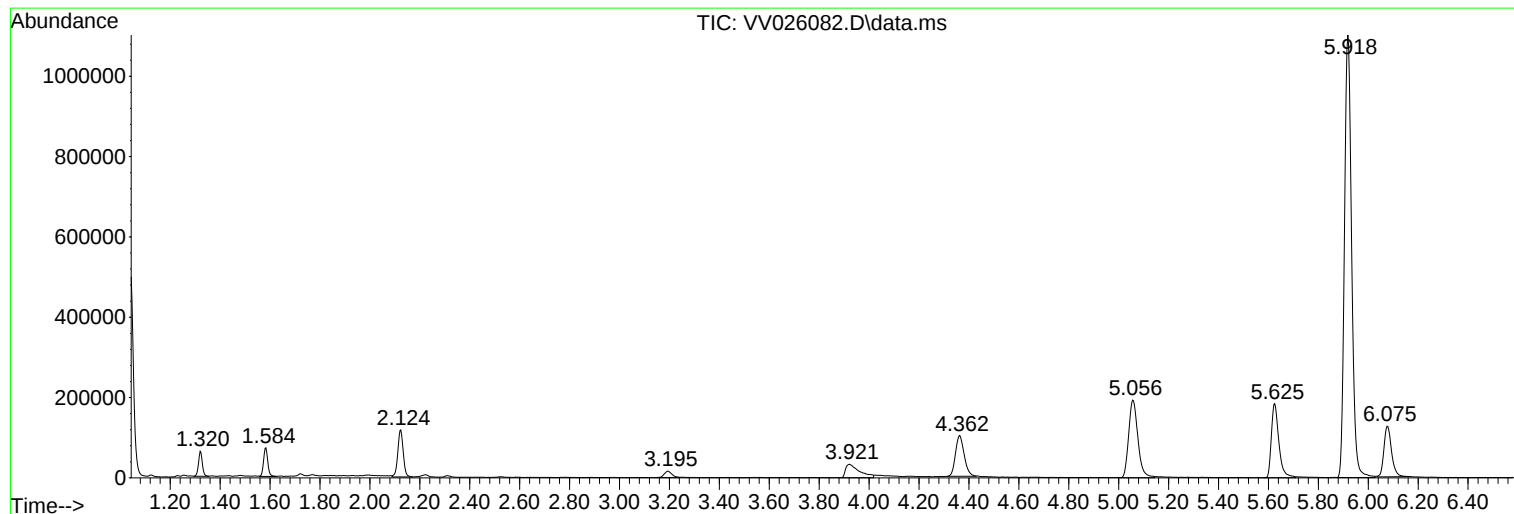
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR060822WMA.M
Quant Title : TRACE VOA SFAM1.0

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



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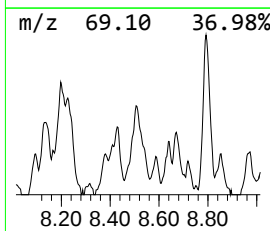
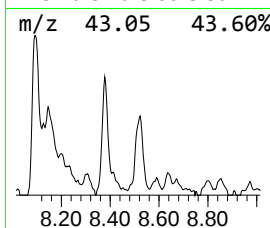
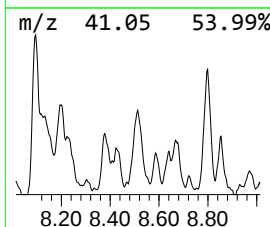
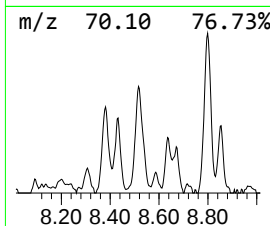
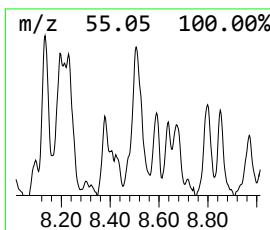
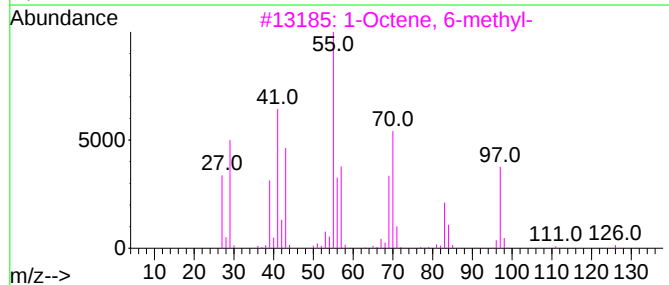
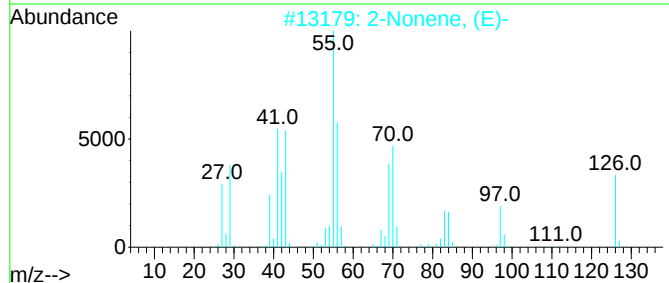
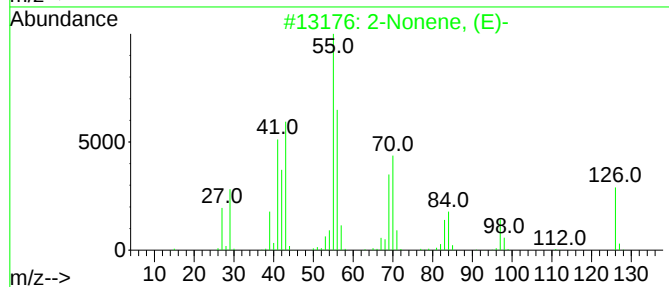
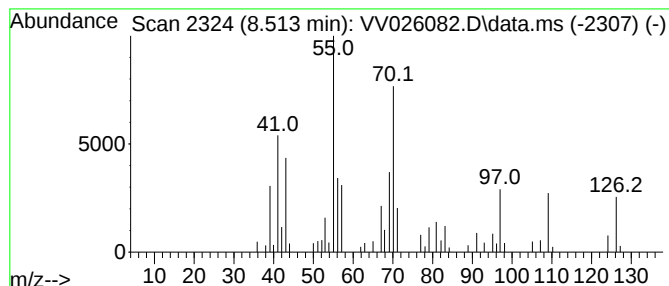
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR060822WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.513	0.74 ug/L	76877	Chlorobenzene-d5	8.854

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Nonene, (E)-	126	C9H18	006434-78-2	50	
2	2-Nonene, (E)-	126	C9H18	006434-78-2	50	
3	1-Octene, 6-methyl-	126	C9H18	013151-10-5	50	
4	Butane, 2-cyclopropyl-	98	C7H14	005750-02-7	43	
5	Cycloheptanone, 2-methyl-	126	C8H14O	000932-56-9	35	



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Instrument :
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ClientSampleId :
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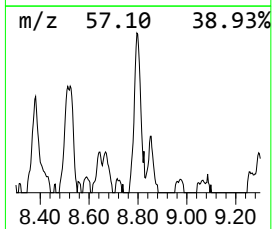
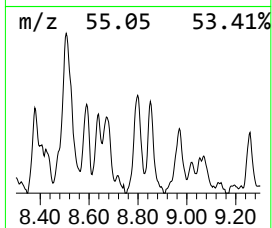
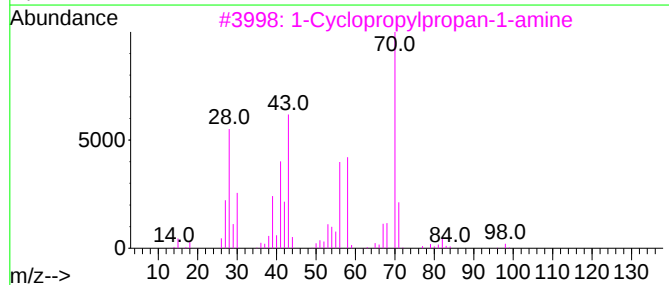
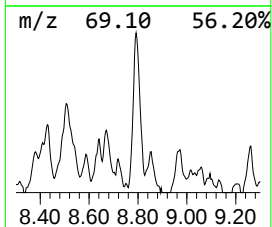
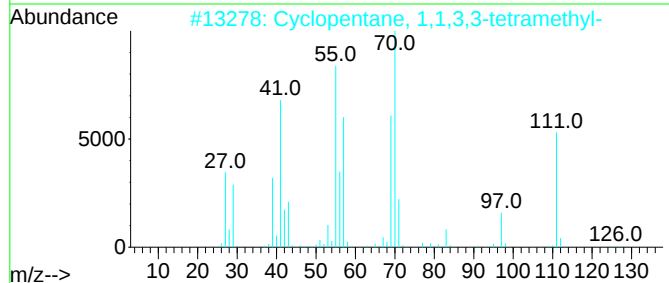
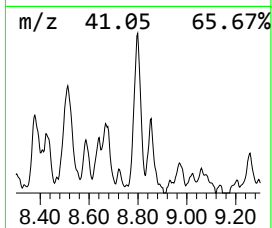
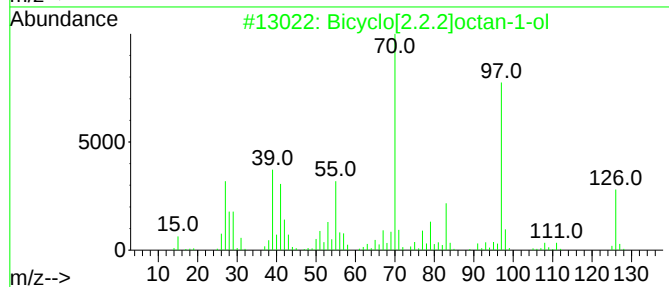
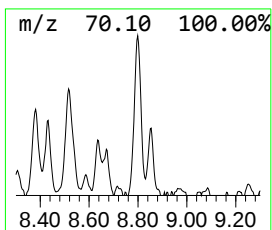
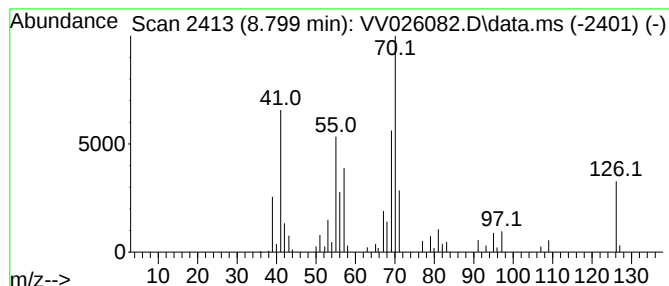
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR060822WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 3 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.799	0.60 ug/L	62416	Chlorobenzene-d5	8.854

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[2.2.2]octan-1-ol	126	C8H14O	020534-58-1	43
2			Cyclopentane, 1,1,3,3-tetramethyl-	126	C9H18	050876-33-0	40
3			1-Cyclopropylpropan-1-amine	99	C6H13N	1000436-93-8	40
4			1-Hexanol, 4-methyl-, acetate	158	C9H18O2	091367-59-8	38
5			Butane, 2-cyclopropyl-	98	C7H14	005750-02-7	38



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ClientSampleId :
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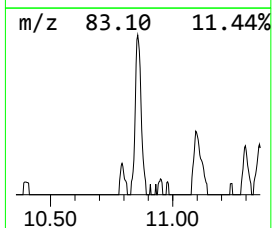
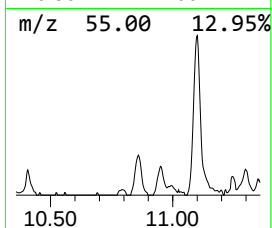
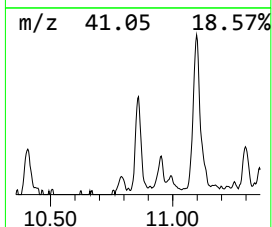
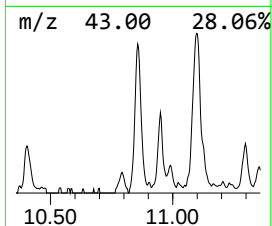
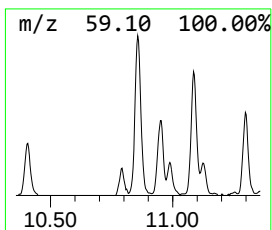
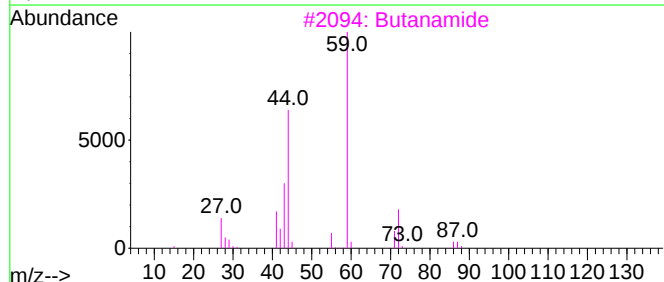
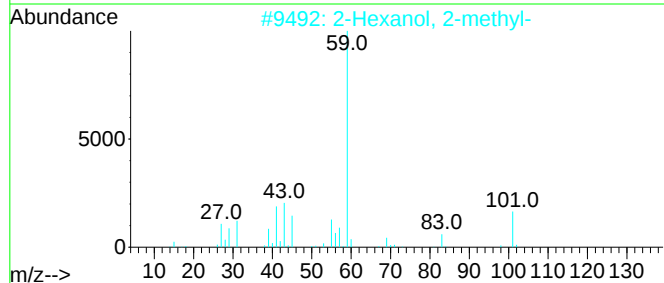
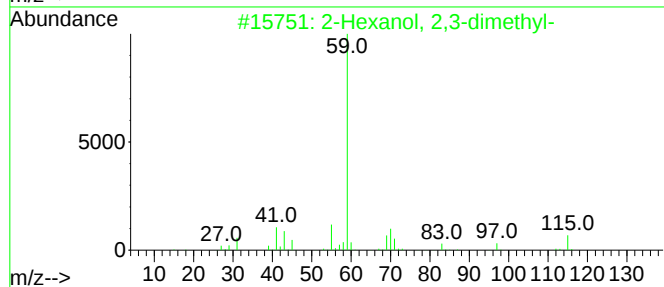
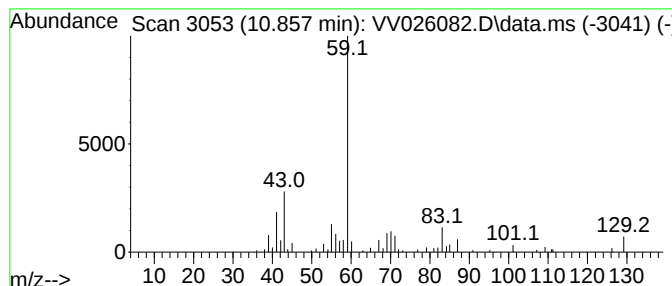
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Quant Title : TRACE VOA SFAM1.0

TIC Library : C:\Database\NIST20.L
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Peak Number 5 2-Hexanol, 2,3-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.857	0.74 ug/L	68678	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Hexanol, 2,3-dimethyl-	130	C8H18O	019550-03-9	59
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	59
3			Butanamide	87	C4H9NO	000541-35-5	53
4			2-Pentanol, 3-chloro-2-methyl-	136	C6H13ClO	074685-49-7	53
5			2-Octanol, 2,6-dimethyl-	158	C10H22O	018479-57-7	50



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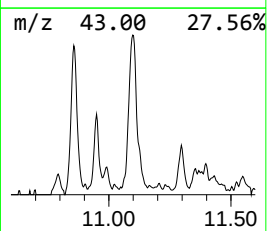
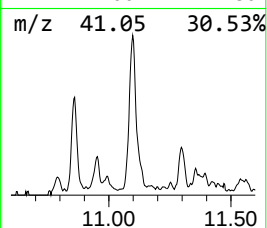
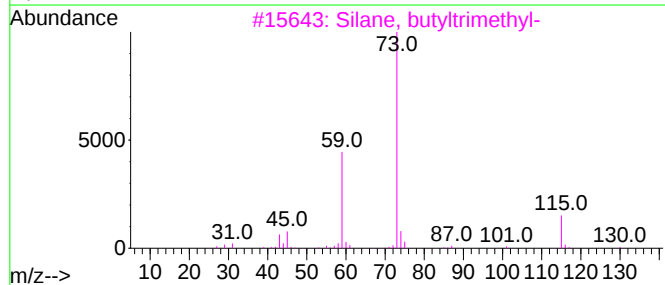
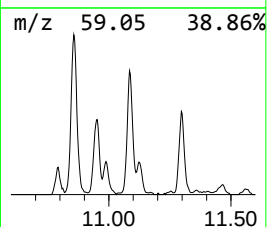
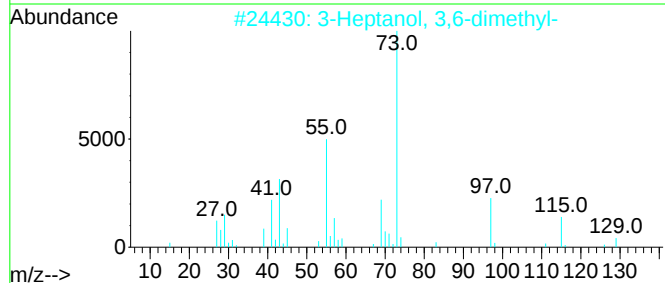
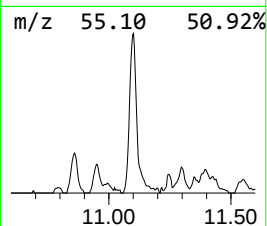
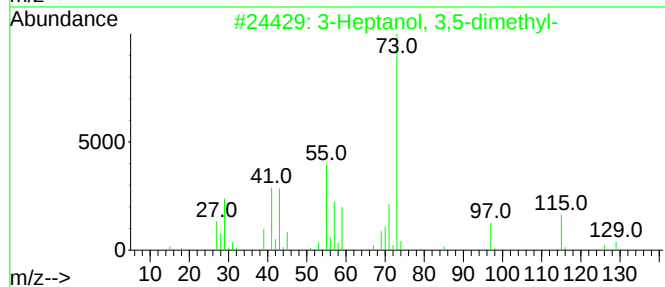
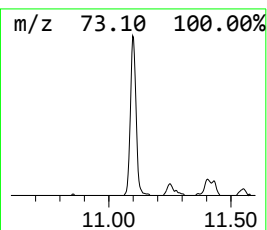
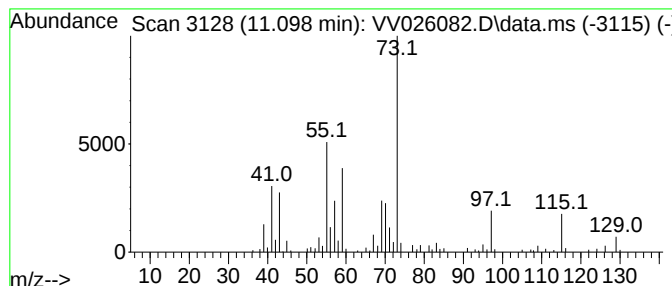
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Peak Number 6 unknown-02 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.098	1.62 ug/L	150175	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Heptanol, 3,5-dimethyl-		144	C9H20O		019549-74-7	38
2	3-Heptanol, 3,6-dimethyl-		144	C9H20O		001573-28-0	32
3	Silane, butyltrimethyl-		130	C7H18Si		001000-49-3	32
4	3-Octanol, 3-methyl-		144	C9H20O		005340-36-3	16
5	Cyclopentane, (1-methylbutyl)-		140	C10H20		004737-43-3	16



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	8.513	0.7	ug/L	76877	2	8.854	516100	5.0
unknown-01	8.799	0.6	ug/L	62416	2	8.854	516100	5.0
2-Hexanol, 2,3-...	10.857	0.7	ug/L	68678	3	11.249	464040	5.0
unknown-02	11.098	1.6	ug/L	150175	3	11.249	464040	5.0