

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU021619\
 Data File : VU029515.D
 Acq On : 15 Feb 2019 22:53
 Operator : JC/SP
 Sample : K1489-04
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 ESXG2

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\SOMUTR021219WMA.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	1.183	22	29	45	rBV	1118142	1028596	100.00%	13.191%
2	1.302	57	66	76	rBV	368253	391136	38.03%	5.016%
3	1.399	88	96	105	rBV2	257445	307397	29.89%	3.942%
4	1.682	177	184	196	rVB	78642	99016	9.63%	1.270%
5	1.756	201	207	222	rVB	76601	108942	10.59%	1.397%
6	2.093	307	312	321	rVB3	14495	19808	1.93%	0.254%
7	2.273	358	368	380	rBV	179254	275426	26.78%	3.532%
8	2.447	415	422	440	rVB	27427	52034	5.06%	0.667%
9	2.987	581	590	605	rVB7	18876	48979	4.76%	0.628%
10	4.219	950	973	1006	rBV3	181965	606119	58.93%	7.773%
11	4.646	1087	1106	1127	rBV2	98859	239565	23.29%	3.072%
12	5.338	1299	1321	1354	rBV2	213903	587461	57.11%	7.534%
13	5.884	1477	1491	1518	rBV	215098	419590	40.79%	5.381%
14	6.183	1574	1584	1599	rVB3	12661	22644	2.20%	0.290%
15	6.328	1614	1629	1648	rBV	134793	270926	26.34%	3.474%
16	7.228	1897	1909	1928	rBV	73902	134601	13.09%	1.726%
17	7.395	1951	1961	1974	rVB3	10586	19543	1.90%	0.251%
18	7.563	2000	2013	2031	rBV	313721	545765	53.06%	6.999%
19	7.849	2089	2102	2120	rBV	44640	80678	7.84%	1.035%
20	8.312	2235	2246	2276	rBV	373332	668451	64.99%	8.572%
21	9.087	2473	2487	2513	rBV	316958	535460	52.06%	6.867%
22	10.431	2892	2905	2918	rBV	117712	190558	18.53%	2.444%
23	10.604	2950	2959	2974	rVB2	10525	18640	1.81%	0.239%
24	11.482	3219	3232	3258	rBV	343264	561298	54.57%	7.198%
25	11.755	3306	3317	3336	rBV7	8645	20672	2.01%	0.265%
26	11.858	3336	3349	3372	rBV	328990	544518	52.94%	6.983%

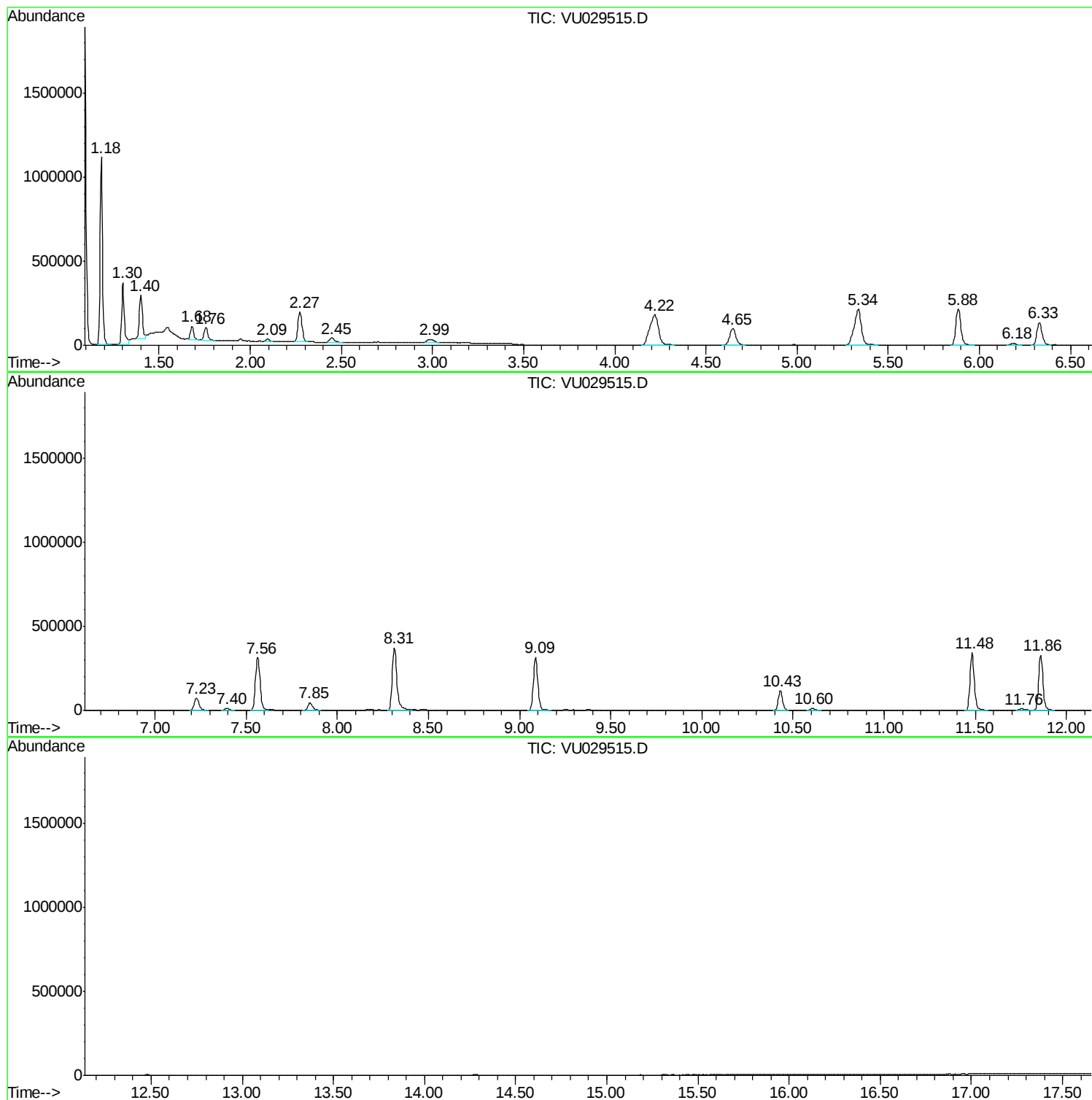
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR021219WMA.M
Quant Title : TRACE VOA SOM01.0

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



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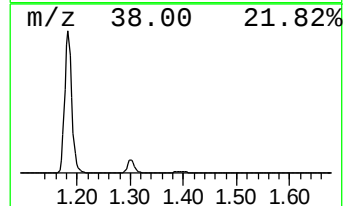
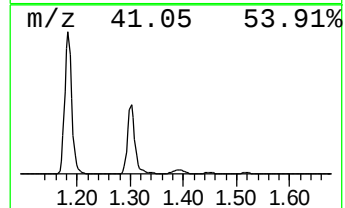
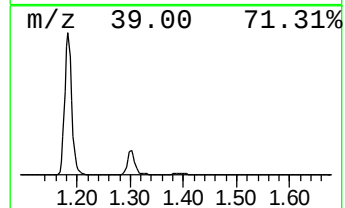
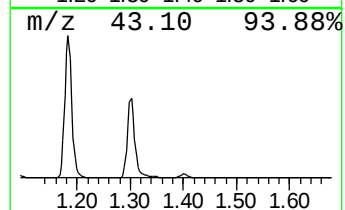
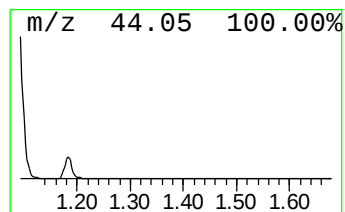
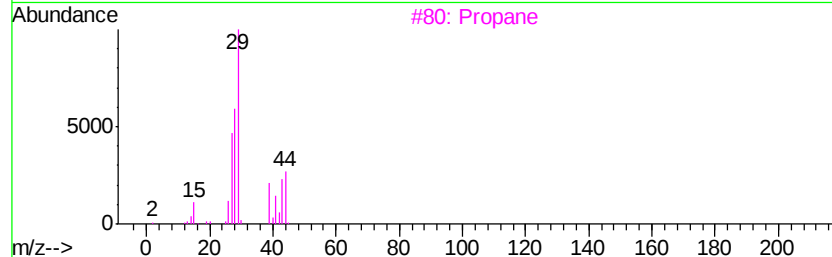
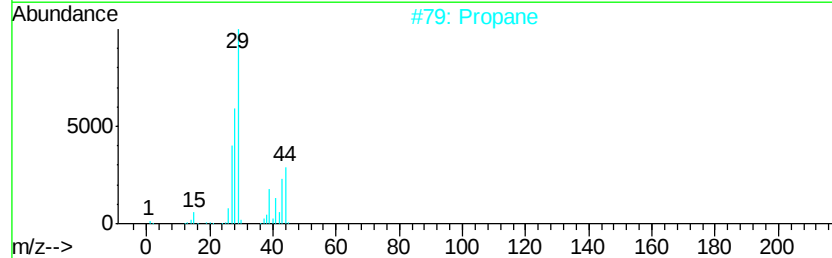
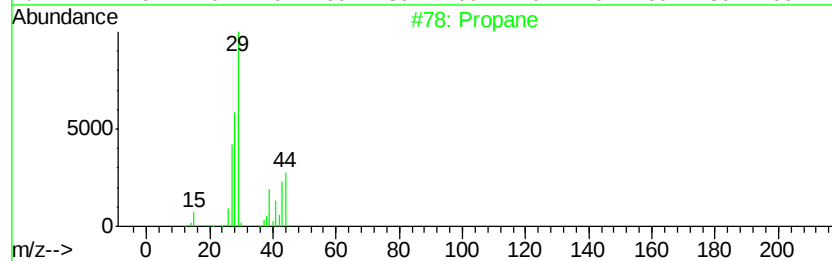
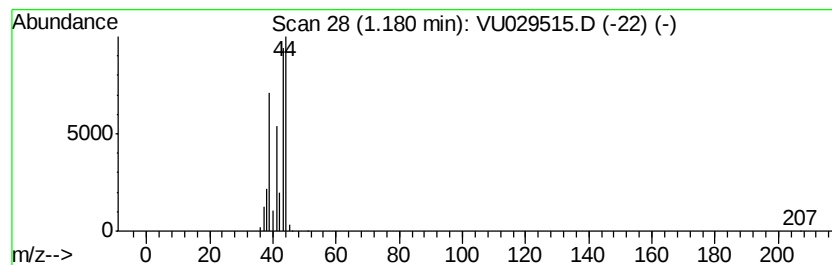
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.18	12.26 ug/L	1028600	1,4-Difluorobenzene	5.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane	44	C3H8	000074-98-6	90
2		Propane	44	C3H8	000074-98-6	83
3		Propane	44	C3H8	000074-98-6	9
4		Propene	42	C3H6	000115-07-1	9
5		Acetaldehyde	44	C2H4O	000075-07-0	5



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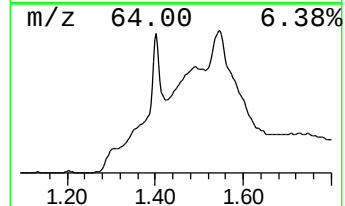
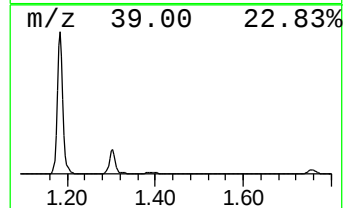
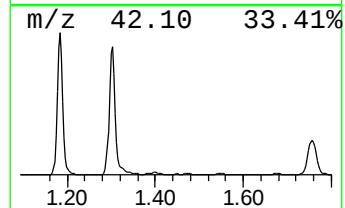
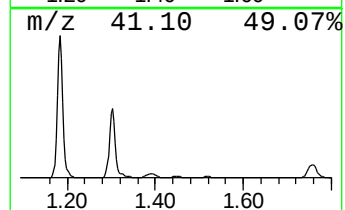
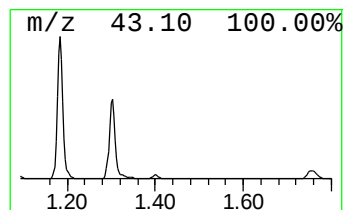
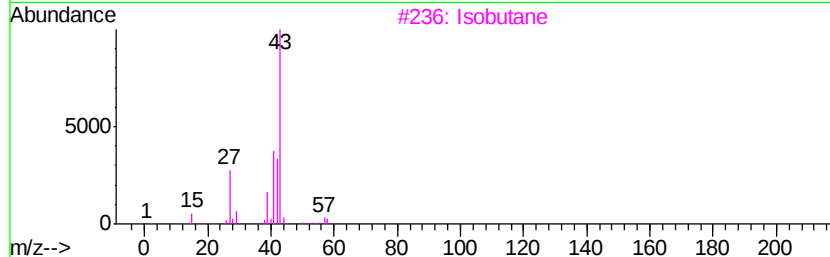
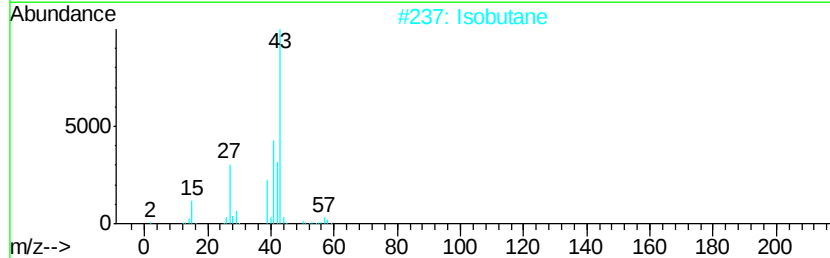
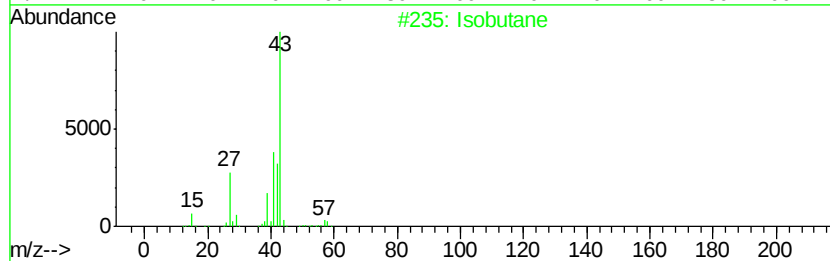
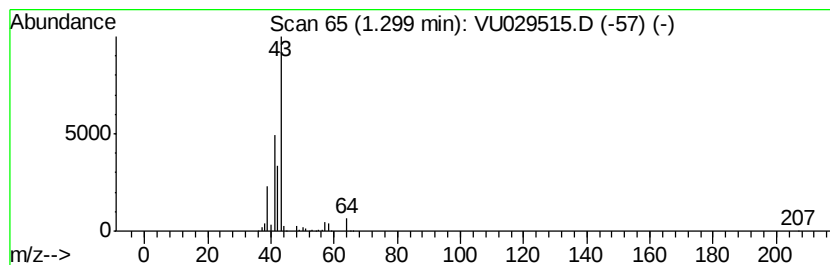
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Quant Title : TRACE VOA SOM01.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.30	4.66 ug/L	391136	1,4-Difluorobenzene	5.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Isobutane	58	C4H10	000075-28-5	59
2		Isobutane	58	C4H10	000075-28-5	59
3		Isobutane	58	C4H10	000075-28-5	53
4		Isobutane	58	C4H10	000075-28-5	53
5		Isopropylsulfonyl chloride	142	C3H7ClO2S	010147-37-2	9



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ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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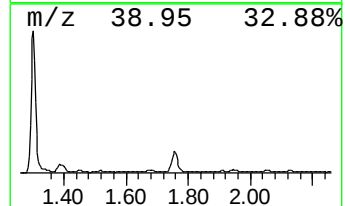
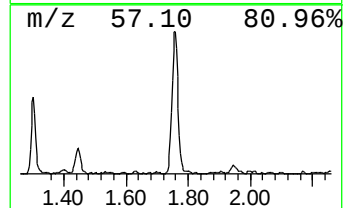
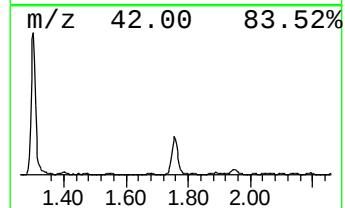
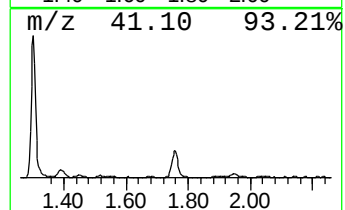
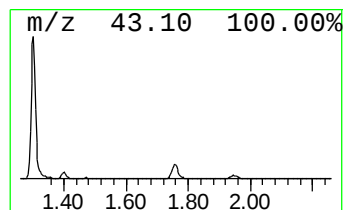
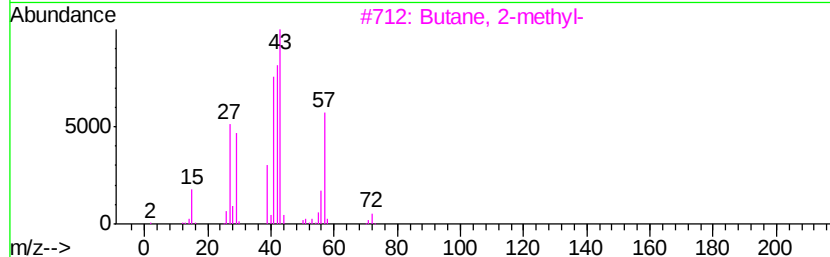
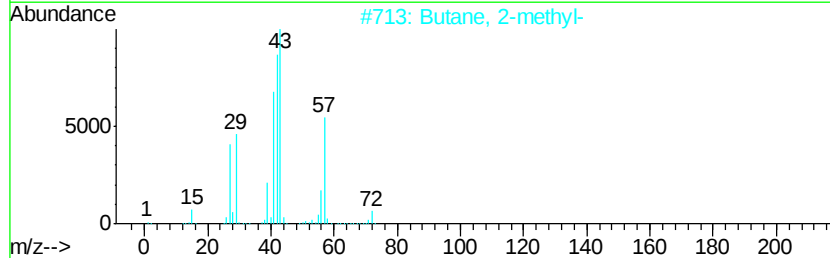
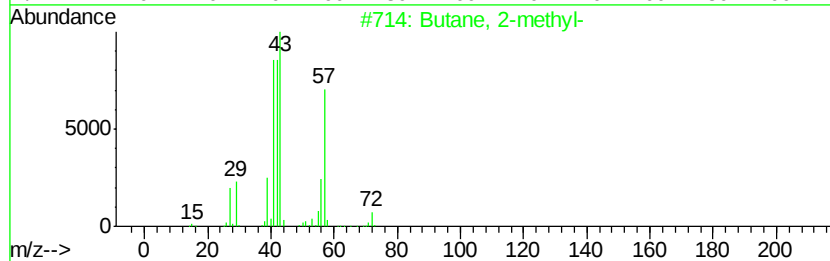
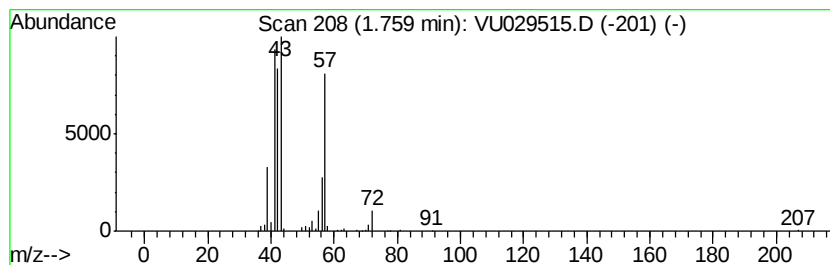
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Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.76	1.30 ug/L	108942	1,4-Difluorobenzene	5.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methyl-	72	C5H12	000078-78-4	91
2		Butane, 2-methyl-	72	C5H12	000078-78-4	86
3		Butane, 2-methyl-	72	C5H12	000078-78-4	80
4		Pentane	72	C5H12	000109-66-0	33
5		Butane, 2-chloro-3-methyl-	106	C5H11Cl	000631-65-2	25



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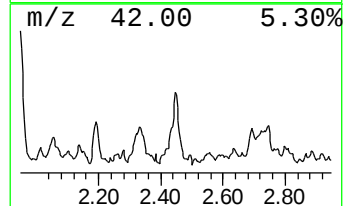
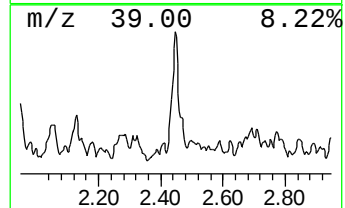
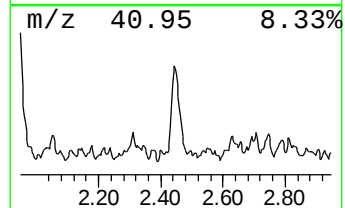
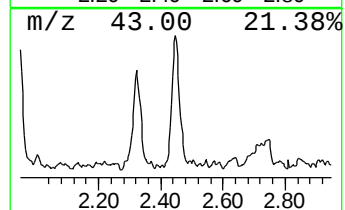
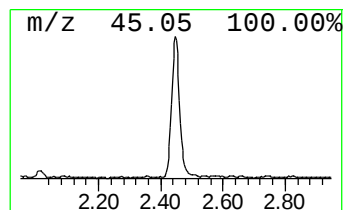
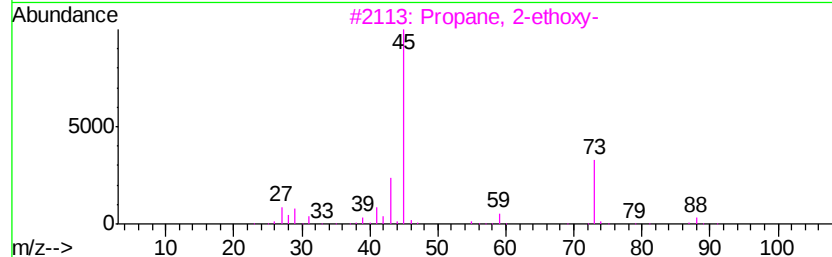
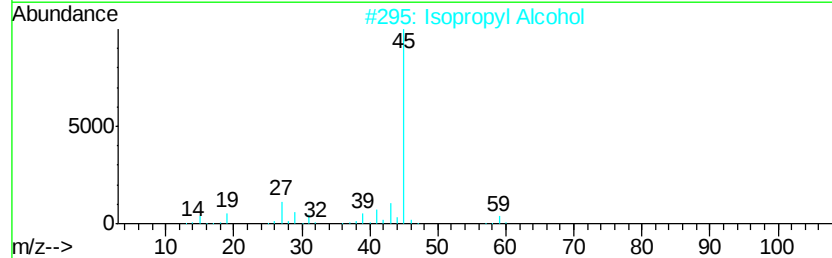
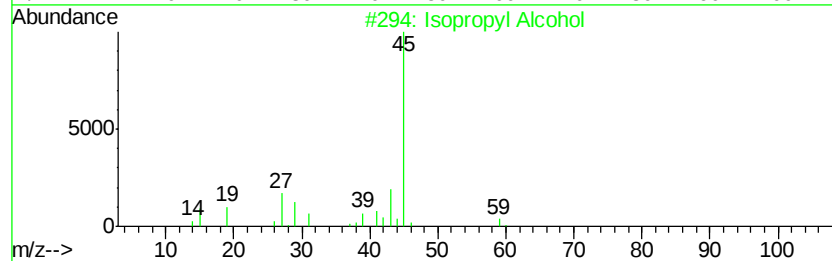
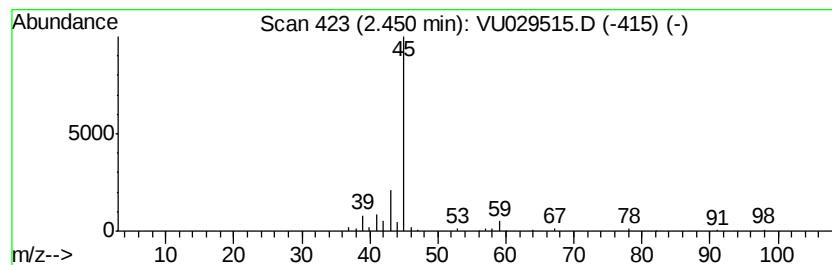
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Peak Number 4 Isopropyl Alcohol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.45	0.62 ug/L	52034	1,4-Difluorobenzene	5.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Isopropyl Alcohol	60	C3H8O	000067-63-0	72
2		Isopropyl Alcohol	60	C3H8O	000067-63-0	72
3		Propane, 2-ethoxy-	88	C5H12O	000625-54-7	56
4		Propane, 2-ethoxy-	88	C5H12O	000625-54-7	56
5		Isopropyl Alcohol	60	C3H8O	000067-63-0	56



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: Str...	1.18	12.3	ug/L	1028600	1	5.88	419590	5.0
(DEL) Alkane: Str...	1.30	4.7	ug/L	391136	1	5.88	419590	5.0
(DEL) Alkane: Str...	1.76	1.3	ug/L	108942	1	5.88	419590	5.0
Isopropyl Alcohol	2.45	0.6	ug/L	52034	1	5.88	419590	5.0