LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU063021\
Data File : VU044266.D
Acq On : 30 Jun 2021 17:39
Operator : SY/MD
Sample : M2905-10
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DBK33

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_U\Method\SFAMUTR062121WMA.M
Title : TRACE VOA SFAM1.0

Signal : TIC: VU044266.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.594	70	79	104	rBV3	268984	544747	2.15%	0.450%
2	2.571	366	383	403	rBV2	132977	307992	1.21%	0.254%
3	3.591	683	700	732	rBV	139750	406417	1.60%	0.335%
4	4.459	951	970	1015	rBV	958614	2778737	10.95%	2.294%
5	5.735	1346	1367	1374	rBV2	196675	509112	2.01%	0.420%
6	5.784	1375	1382	1402	rVB	247348	490315	1.93%	0.405%
7	6.256	1516	1529	1541	rBV	171266	324785	1.28%	0.268%
8	7.906	2031	2042	2057	rVB	248652	429463	1.69%	0.355%
9	8.198	2120	2133	2151	rBV3	184292	518103	2.04%	0.428%
10	8.639	2254	2270	2294	rBV	963280	1714118	6.76%	1.415%
11	9.427	2506	2515	2519	rBV	228657	368032	1.45%	0.304%
12	9.603	2557	2570	2579	rBV	201917	315420	1.24%	0.260%
13	9.690	2579	2597	2610	rBV7	577217	1512721	5.96%	1.249%
14	10.709	2902	2914	2926	rBV	1704112	2827403	11.15%	2.334%
15	10.967	2985	2994	3014	rVB3	144386	258404	1.02%	0.213%
16	11.098	3026	3035	3045	rVB	871470	1354436	5.34%	1.118%
17	11.729	3221	3231	3240	rBV	568133	937047	3.69%	0.774%
18	11.851	3251	3269	3287	rVB	16168388	25366092	100.00%	20.940%
19	11.947	3289	3299	3306	rBV	631090	1013779	4.00%	0.837%
20	12.050	3320	3331	3340	rBV	2163368	3279743	12.93%	2.707%
21	12.105	3340	3348	3358	rVB3	801552	1265700	4.99%	1.045%
22	12.198	3367	3377	3384	rBV2	377339	694197	2.74%	0.573%
23	12.262	3391	3397	3405	rVB	490741	739814	2.92%	0.611%
24	12.385	3428	3435	3448	rVB6	219709	379204	1.49%	0.313%
25	12.487	3449	3467	3474	rBV2	534372	1190634	4.69%	0.983%
26	12.581	3487	3496	3506	rVB2	860943	1299750	5.12%	1.073%
27	12.729	3533	3542	3561	rVB2	874806	2022054	7.97%	1.669%
28	12.835	3561	3575	3584	rBV3	1206625	2571644	10.14%	2.123%
29	12.947	3598	3610	3617	rBV5	534256	946295	3.73%	0.781%
30	13.018	3617	3632	3644	rVB	5078235	8327846	32.83%	6.875%
31	13.111	3646	3661	3674	rBV	4851136	7615356	30.02%	6.286%
32	13.182	3675	3683	3690	rVB2	229593	319177	1.26%	0.263%
33	13.262	3692	3708	3725	rVB3	416299	1214587	4.79%	1.003%
34	13.352	3726	3736	3741	rBV3	339384	607329	2.39%	0.501%
35	13.462	3761	3770	3773	rBV4	334293	468374	1.85%	0.387%
36	13.494	3773	3780	3788	rVV	1274091	1910685	7.53%	1.577%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU063021\
Data File : VU044266.D
Acq On : 30 Jun 2021 17:39
Operator : SY/MD
Sample : M2905-10
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DBK33

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_U\Method\SFAMUTR062121WMA.M
Title : TRACE VOA SFAM1.0

37	13.542	3788	3795	3803	rVB	258340	373457	1.47%	0.308%
38	13.671	3826	3835	3846	rBV3	613184	1070343	4.22%	0.884%
39	13.738	3847	3856	3874	rVB2	1924773	3615809	14.25%	2.985%
40	13.828	3874	3884	3894	rBV2	1000221	1559949	6.15%	1.288%
41	13.976	3915	3930	3952	rBV	8761606	15488079	61.06%	12.785%
42	14.195	3983	3998	4009	rBV4	488878	1361105	5.37%	1.124%
43	14.327	4030	4039	4047	rBV	582585	917331	3.62%	0.757%
44	14.381	4047	4056	4062	rBV4	635670	1086679	4.28%	0.897%
45	14.442	4070	4075	4085	rVB	610923	860248	3.39%	0.710%
46	14.536	4095	4104	4111	rBV3	381956	615970	2.43%	0.508%
47	14.584	4111	4119	4134	rVB	1784057	2534701	9.99%	2.092%
48	14.655	4135	4141	4162	rVB5	300680	668543	2.64%	0.552%
49	14.761	4162	4174	4181	rBV	2128503	3285157	12.95%	2.712%
50	14.899	4209	4217	4228	rVB	238681	377926	1.49%	0.312%
51	15.188	4297	4307	4318	rVB	2892950	4353086	17.16%	3.593%
52	15.272	4326	4333	4342	rVB	237959	375898	1.48%	0.310%
53	15.368	4349	4363	4369	rBV6	727933	1532118	6.04%	1.265%
54	15.410	4370	4376	4384	rVB	702893	1046194	4.12%	0.864%
55	15.655	4441	4452	4472	rVB	650907	1345708	5.31%	1.111%
56	15.809	4492	4500	4513	rBV2	171257	286651	1.13%	0.237%
57	16.204	4614	4623	4638	rVB5	165533	359645	1.42%	0.297%
58	16.394	4672	4682	4699	rBV5	137255	299638	1.18%	0.247%
59	16.835	4807	4819	4839	rVB2	183496	345479	1.36%	0.285%
60	17.021	4866	4877	4891	rBV2	356980	550244	2.17%	0.454%

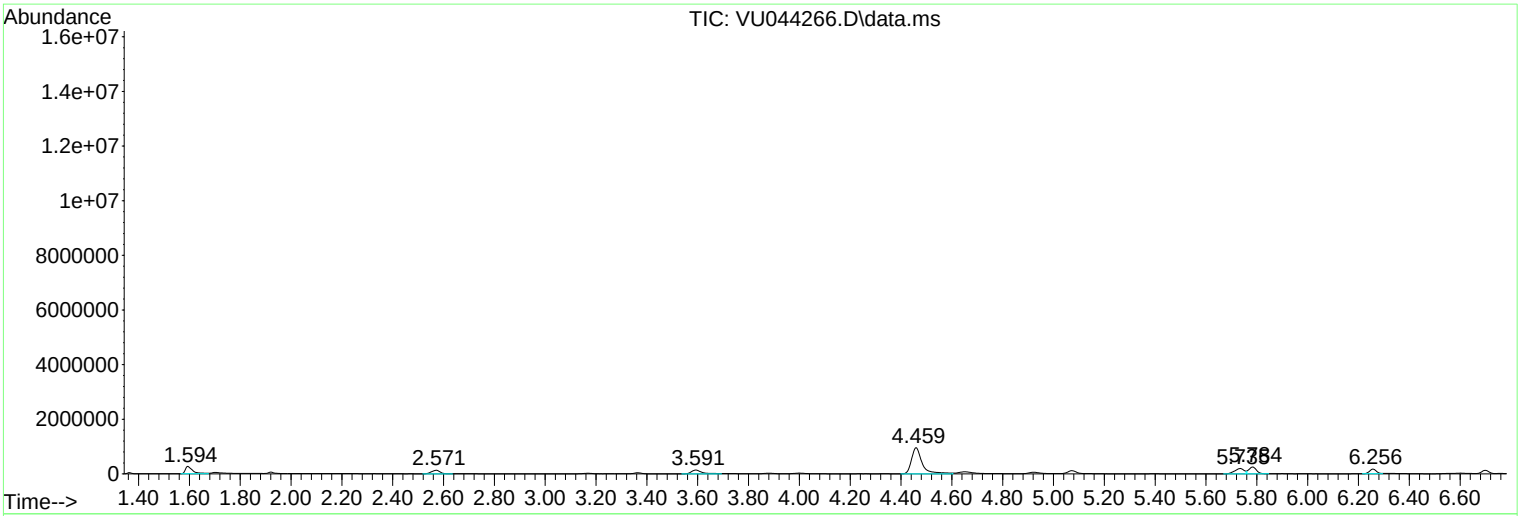
Sum of corrected areas: 121139470

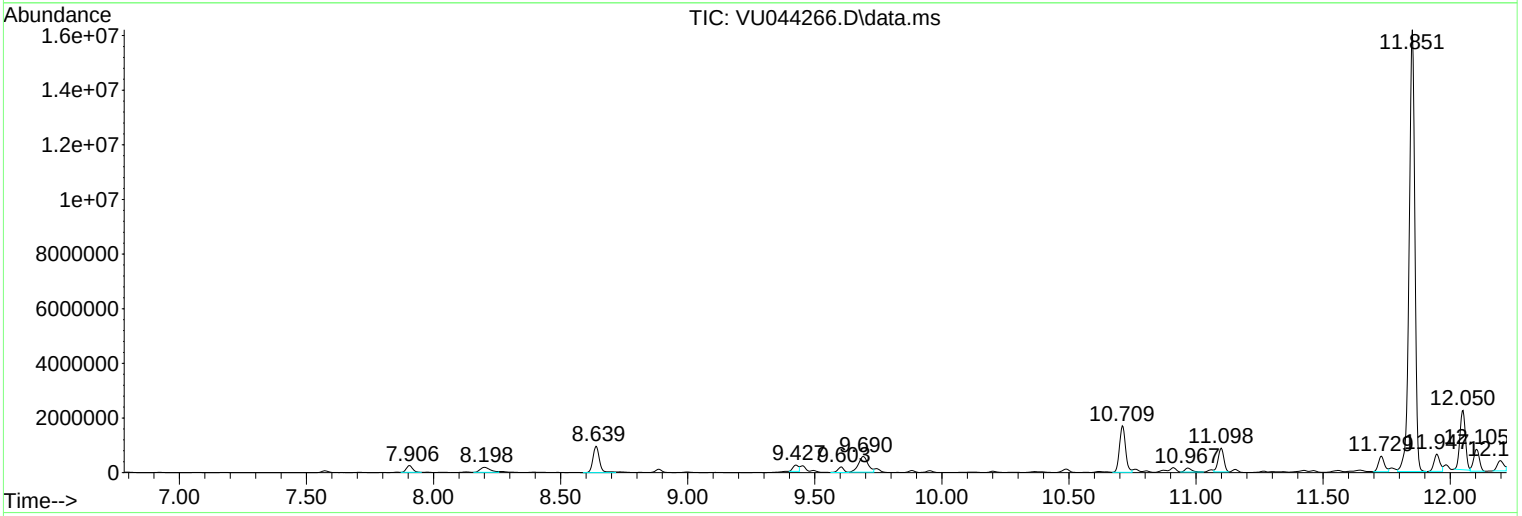
Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU063021\
Data File : VU044266.D
Acq On : 30 Jun 2021 17:39
Operator : SY/MD
Sample : M2905-10
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

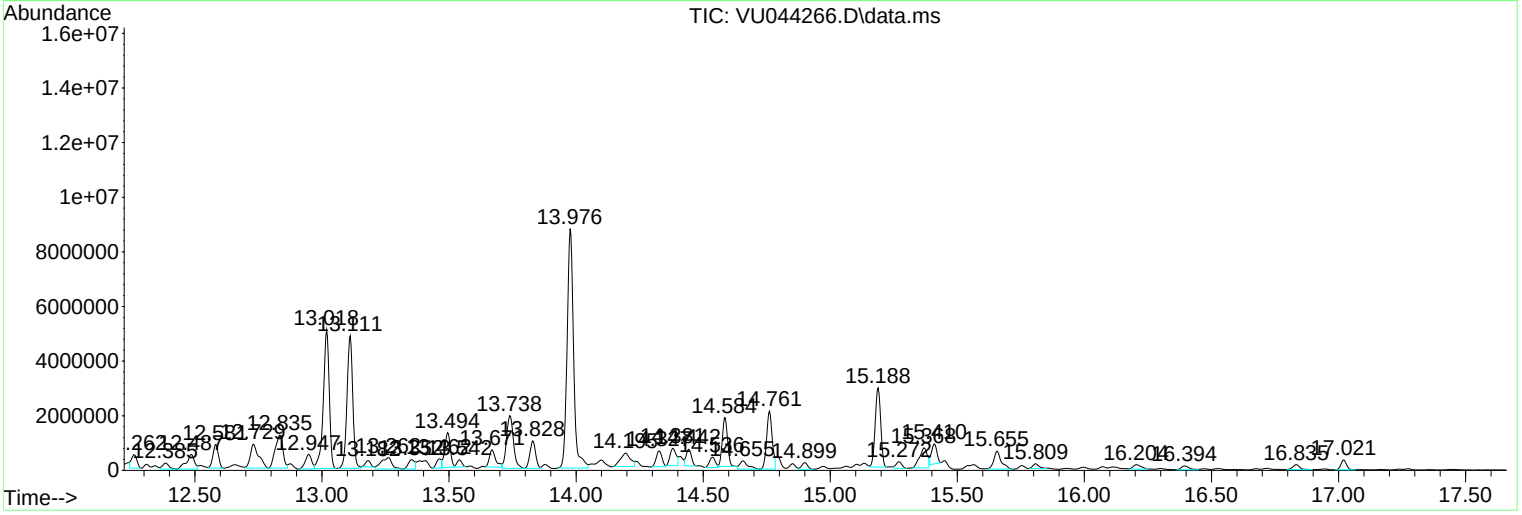
Instrument :
MSVOA_U
ClientSampleId :
DBK33

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR062121WMA.M
Quant Title : TRACE VOA SFAM1.0

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







Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU063021\
Data File : VU044266.D
Acq On : 30 Jun 2021 17:39
Operator : SY/MD
Sample : M2905-10
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DBK33

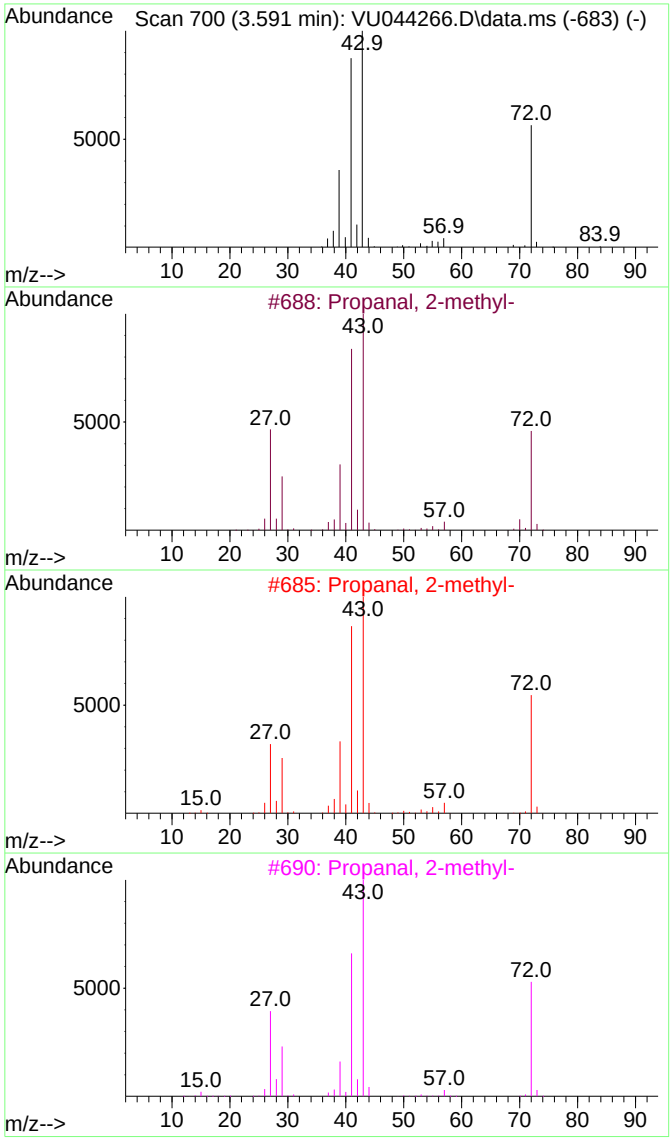
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR062121WMA.M
Quant Title : TRACE VOA SFAM1.0

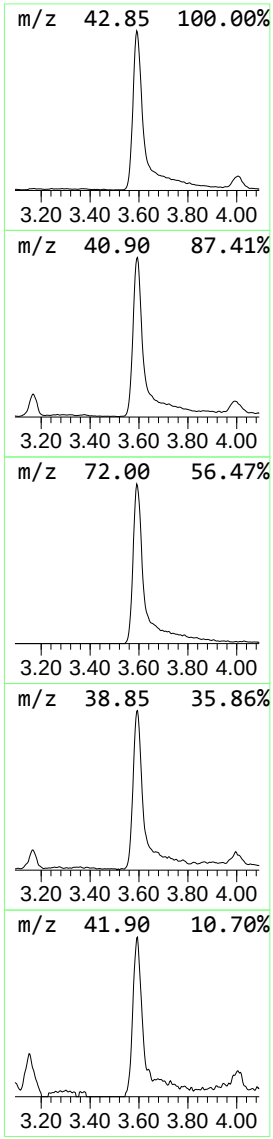
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Propanal, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.591	6.26 ug/L	406417	1,4-Difluorobenzene	6.256

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanal, 2-methyl-	72	C4H8O	000078-84-2	91
2	Propanal, 2-methyl-	72	C4H8O	000078-84-2	91
3	Propanal, 2-methyl-	72	C4H8O	000078-84-2	78
4	1-Propene, 3-methoxy-	72	C4H8O	000627-40-7	53
5	Propanal, 2-methyl-	72	C4H8O	000078-84-2	43





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU063021\
Data File : VU044266.D
Acq On : 30 Jun 2021 17:39
Operator : SY/MD
Sample : M2905-10
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DBK33

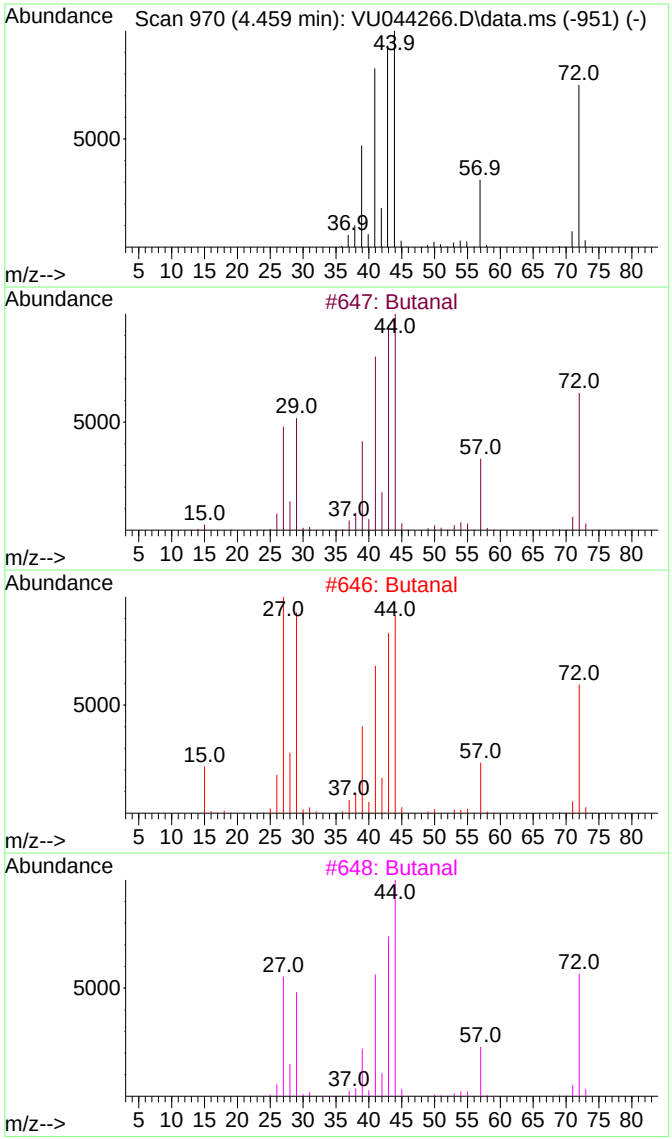
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR062121WMA.M
Quant Title : TRACE VOA SFAM1.0

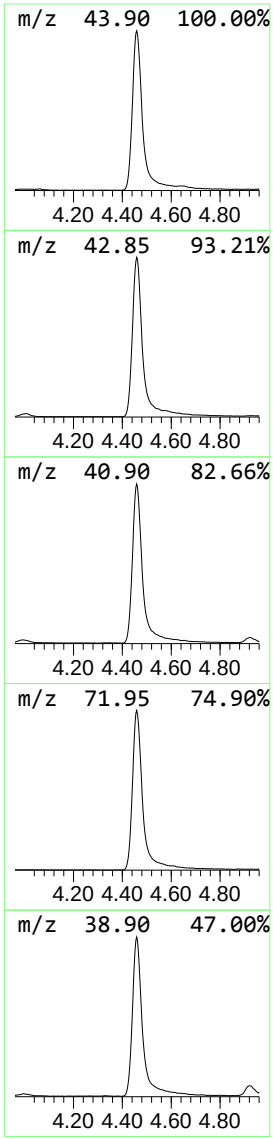
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Butanal Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.459	42.78 ug/L	2778740	1,4-Difluorobenzene	6.256

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanal	72	C4H8O	000123-72-8	94
2	Butanal	72	C4H8O	000123-72-8	91
3	Butanal	72	C4H8O	000123-72-8	72
4	Ethene, ethoxy-	72	C4H8O	000109-92-2	52
5	Butanal	72	C4H8O	000123-72-8	46





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU063021\
Data File : VU044266.D
Acq On : 30 Jun 2021 17:39
Operator : SY/MD
Sample : M2905-10
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DBK33

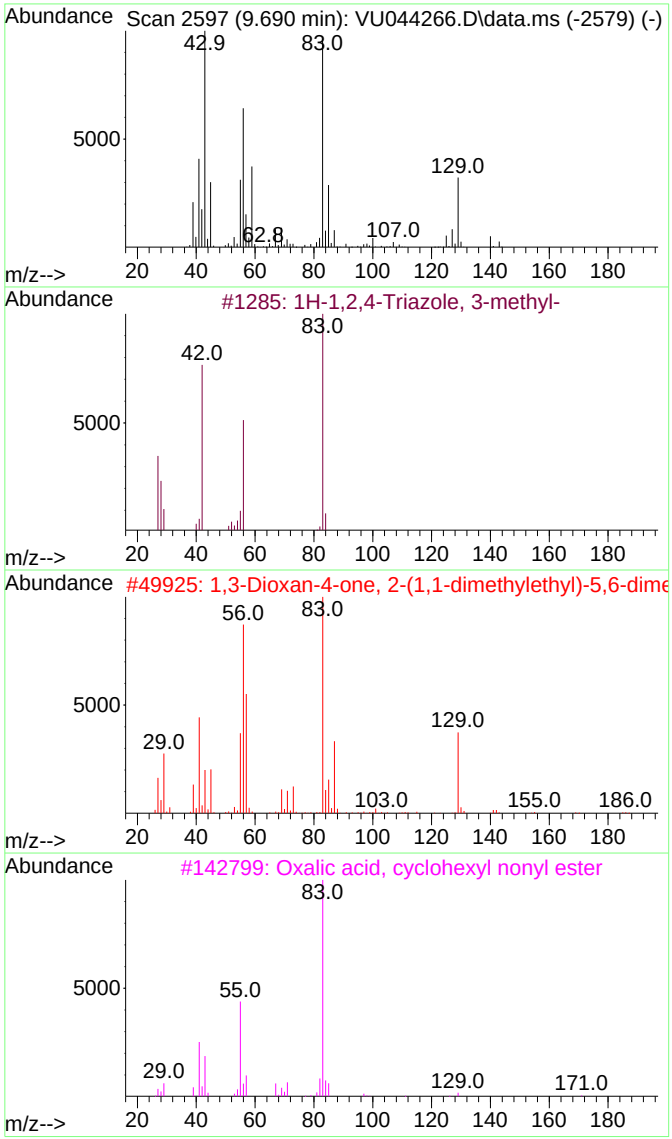
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR062121WMA.M
Quant Title : TRACE VOA SFAM1.0

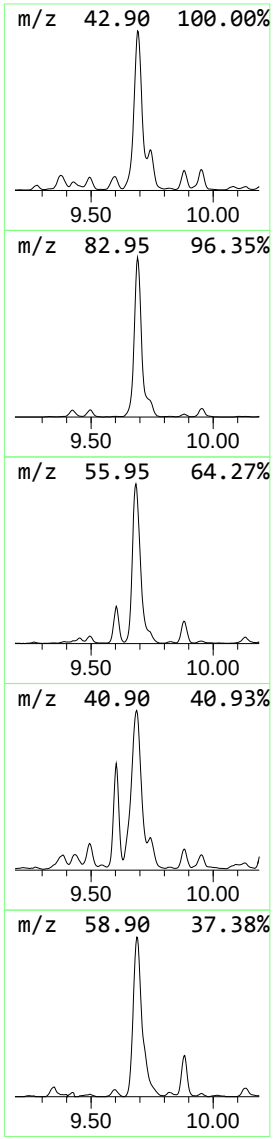
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.690	20.55 ug/L	1512720	Chlorobenzene-d5	9.423

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-1,2,4-Triazole, 3-methyl-	83	C3H5N3	007170-01-6	43
2	1,3-Dioxan-4-one, 2-(1,1-dimethy...	186	C10H18O3	107289-14-5	42
3	Oxalic acid, cyclohexyl nonyl ester	298	C17H30O4	1000309-31-1	35
4	Oxalic acid, cyclohexyl decyl ester	312	C18H32O4	1000309-31-2	27
5	2-Butyl-1,2-azaborolidine	125	C7H16BN	005357-10-8	27





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU063021\
Data File : VU044266.D
Acq On : 30 Jun 2021 17:39
Operator : SY/MD
Sample : M2905-10
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DBK33

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR062121WMA.M
Quant Title : TRACE VOA SFAM1.0

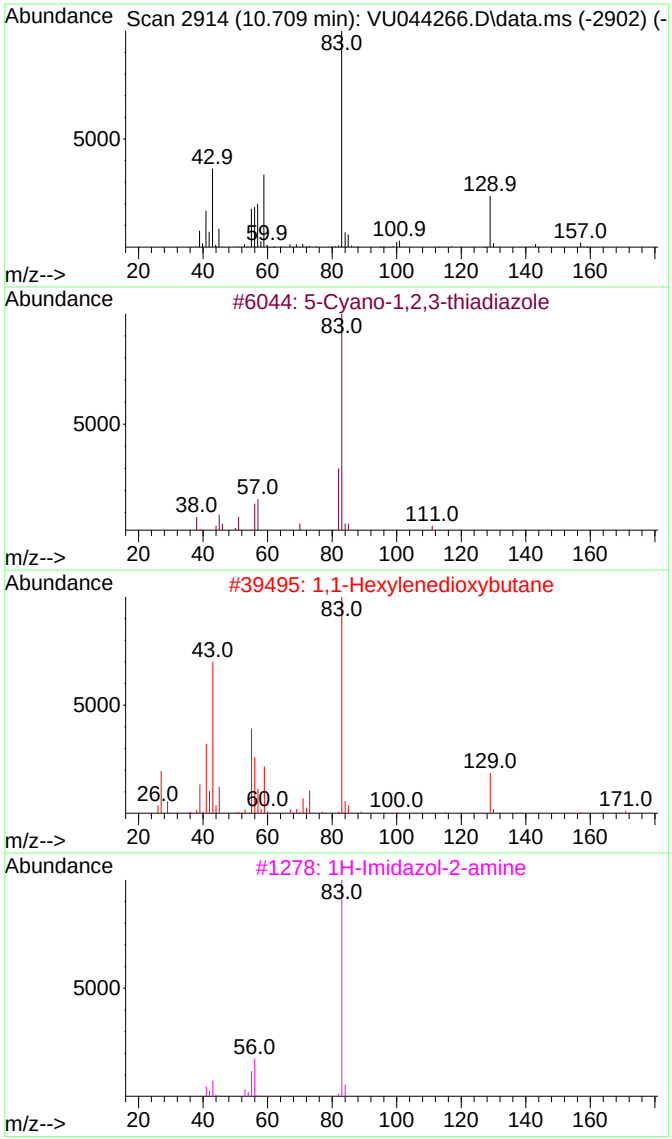
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

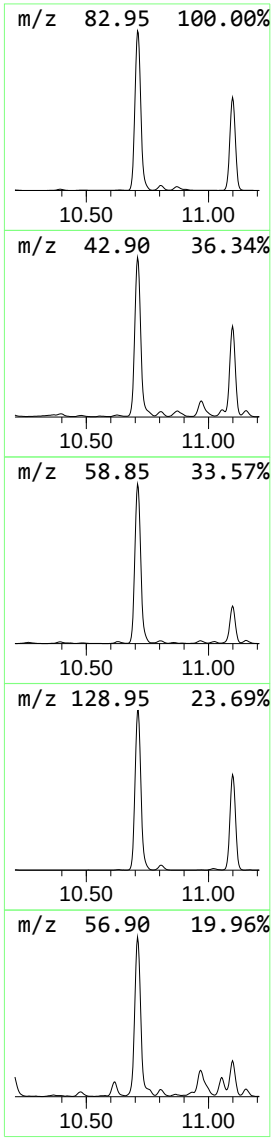
****

Peak Number 4 unknown-02 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.709	0.56 ug/L	2827400	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Cyano-1,2,3-thiadiazole	111	C3HN3S	057352-02-0	40
2	1,1-Hexylenedioxybutane	172	C10H20O2	1000249-77-4	39
3	1H-Imidazol-2-amine	83	C3H5N3	007720-39-0	9
4	1,4-Pentadien-3-ol	84	C5H8O	000922-65-6	9
5	.alpha.-Amino-2,5-dihydro-5-meth...	157	C7H11NO3	018455-25-9	9





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU063021\
Data File : VU044266.D
Acq On : 30 Jun 2021 17:39
Operator : SY/MD
Sample : M2905-10
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DBK33

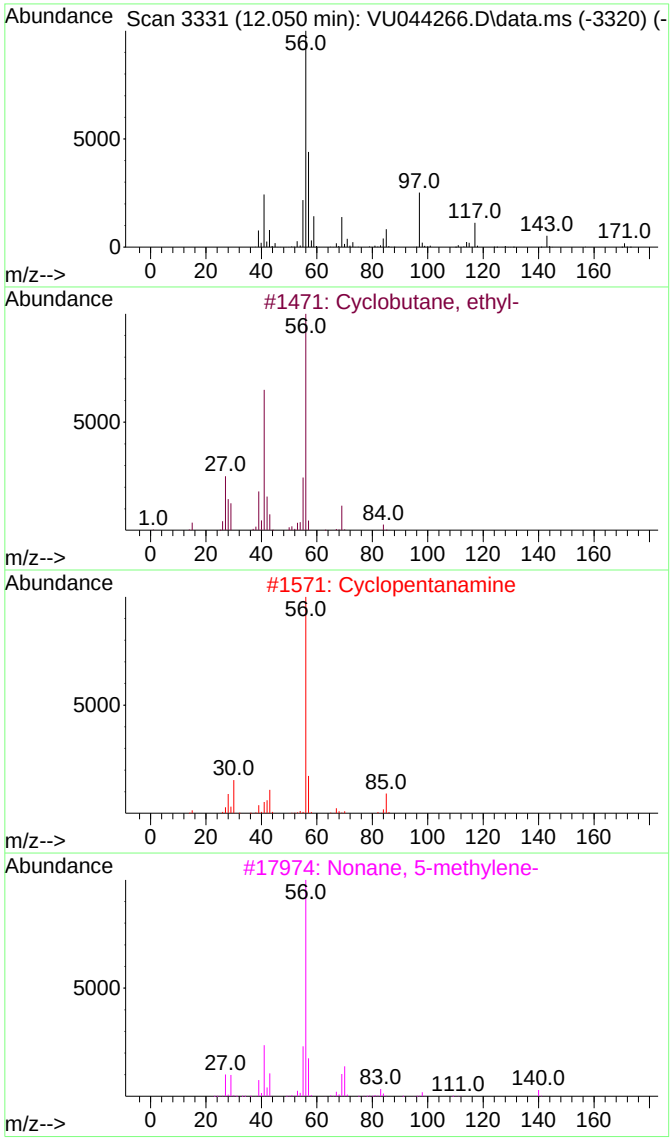
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR062121WMA.M
Quant Title : TRACE VOA SFAM1.0

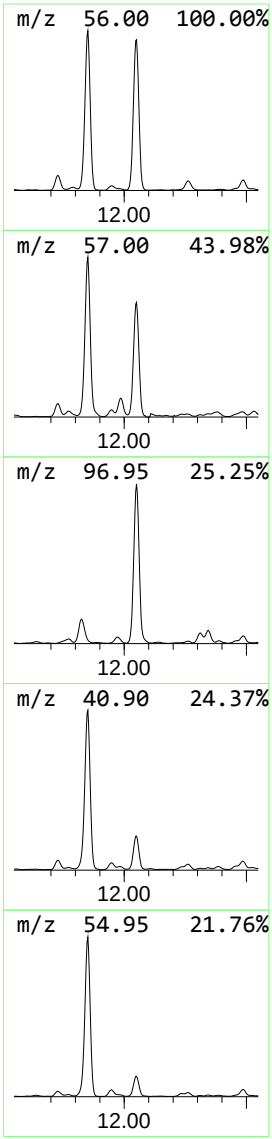
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 (DEL) Alkane: Cyclic12.050 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.050	0.65 ug/L	3279740	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclobutane, ethyl-	84	C6H12	004806-61-5	47
2	Cyclopentanamine	85	C5H11N	001003-03-8	38
3	Nonane, 5-methylene-	140	C10H20	006795-79-5	33
4	1H-Pyrrole-2,5-dione	97	C4H3NO2	000541-59-3	9
5	Hexane, 3,4-dimethyl-	114	C8H18	000583-48-2	9





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU063021\
Data File : VU044266.D
Acq On : 30 Jun 2021 17:39
Operator : SY/MD
Sample : M2905-10
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DBK33

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR062121WMA.M
Quant Title : TRACE VOA SFAM1.0

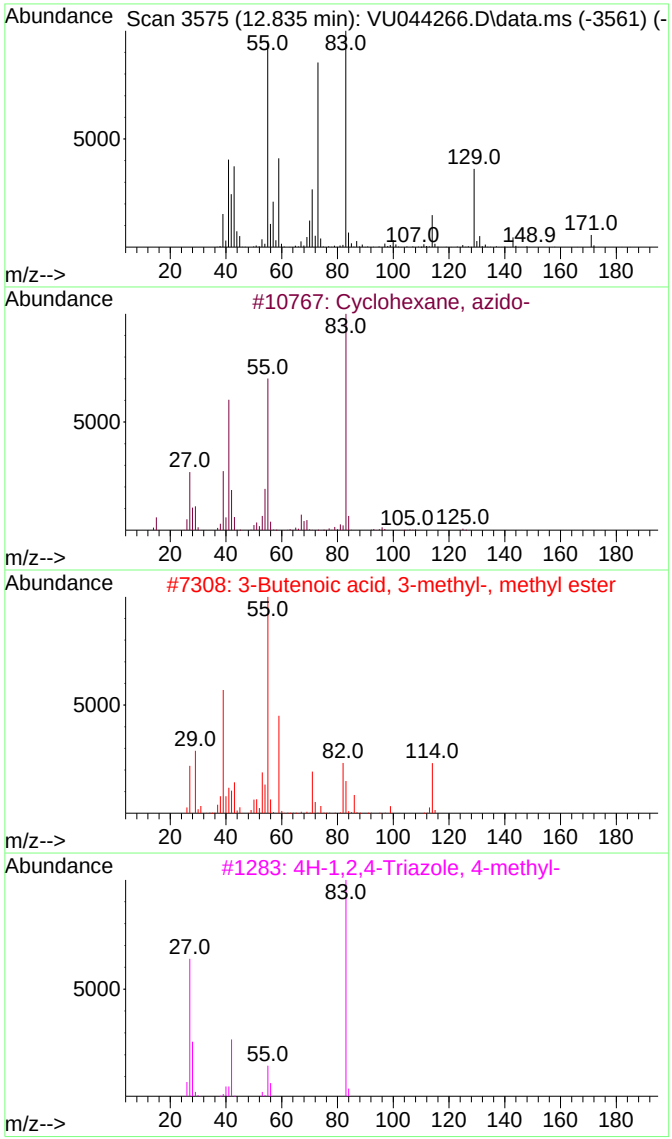
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

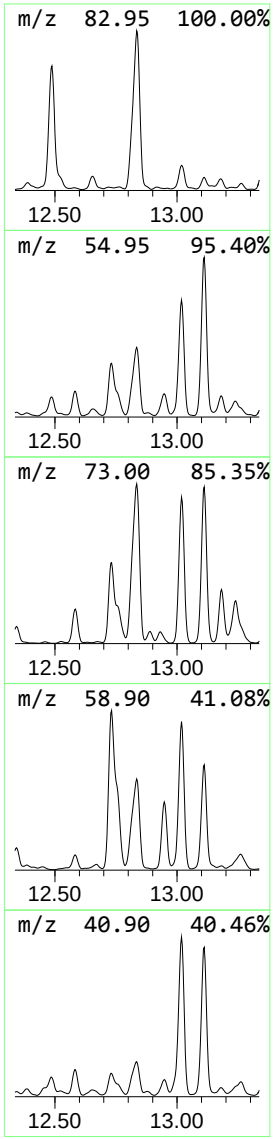
****

Peak Number 6 unknown-03 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.835	0.51 ug/L	2571640	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, azido-	125	C6H11N3	019573-22-9	38
2	3-Butenoic acid, 3-methyl-, meth...	114	C6H10O2	025859-52-3	30
3	4H-1,2,4-Triazole, 4-methyl-	83	C3H5N3	010570-40-8	27
4	2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	27
5	2-Hydroxy-2-methyl-but-3-enyl 2-...	184	C10H16O3	080758-67-4	17





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU063021\
Data File : VU044266.D
Acq On : 30 Jun 2021 17:39
Operator : SY/MD
Sample : M2905-10
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DBK33

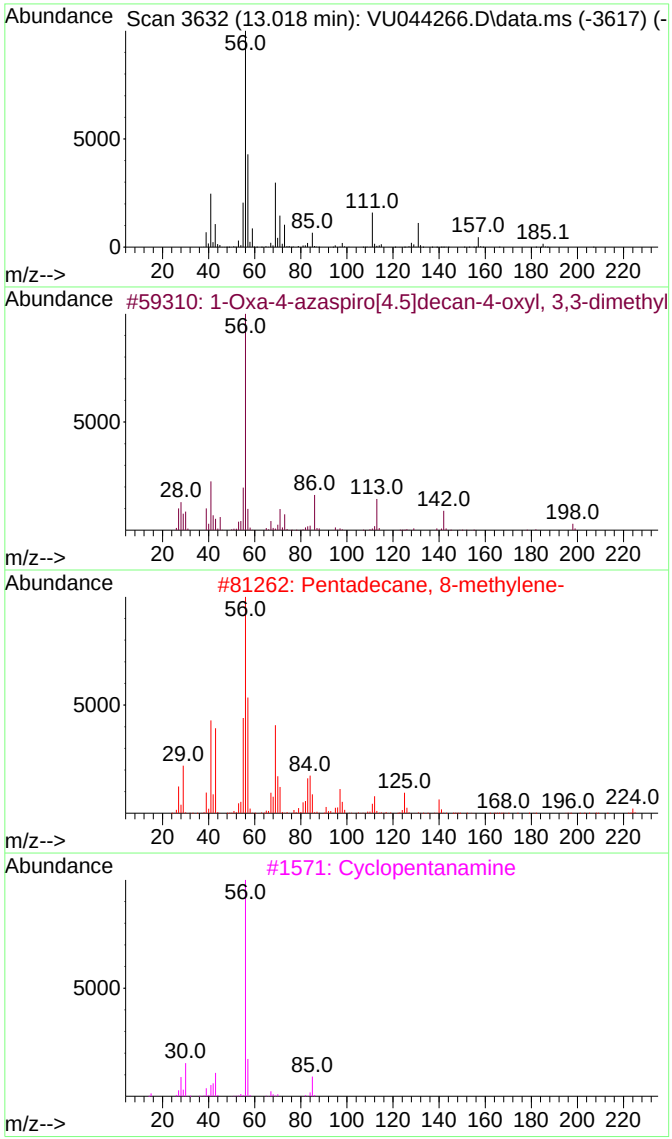
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR062121WMA.M
Quant Title : TRACE VOA SFAM1.0

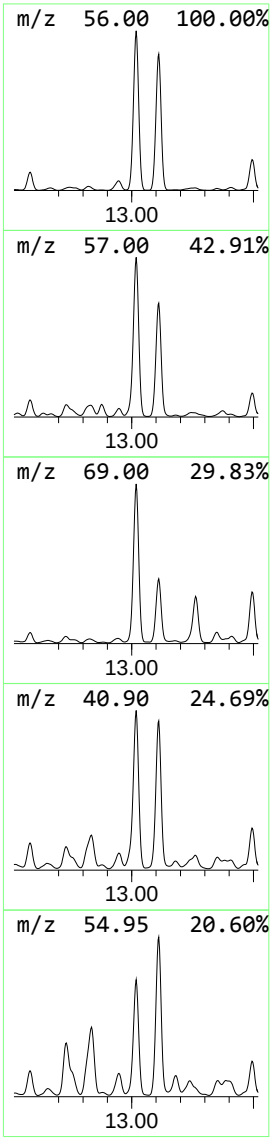
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 unknown-04 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.018	1.64 ug/L	8327850	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Oxa-4-azaspiro[4.5]decan-4-oxy...	198	C10H16NO3	1000115-84-0	42
2	Pentadecane, 8-methylene-	224	C16H32	055668-09-2	42
3	Cyclopentanamine	85	C5H11N	001003-03-8	40
4	Cyclopentane, (1,1-dimethylethyl)-	126	C9H18	003875-52-3	38
5	2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	36





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU063021\
Data File : VU044266.D
Acq On : 30 Jun 2021 17:39
Operator : SY/MD
Sample : M2905-10
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DBK33

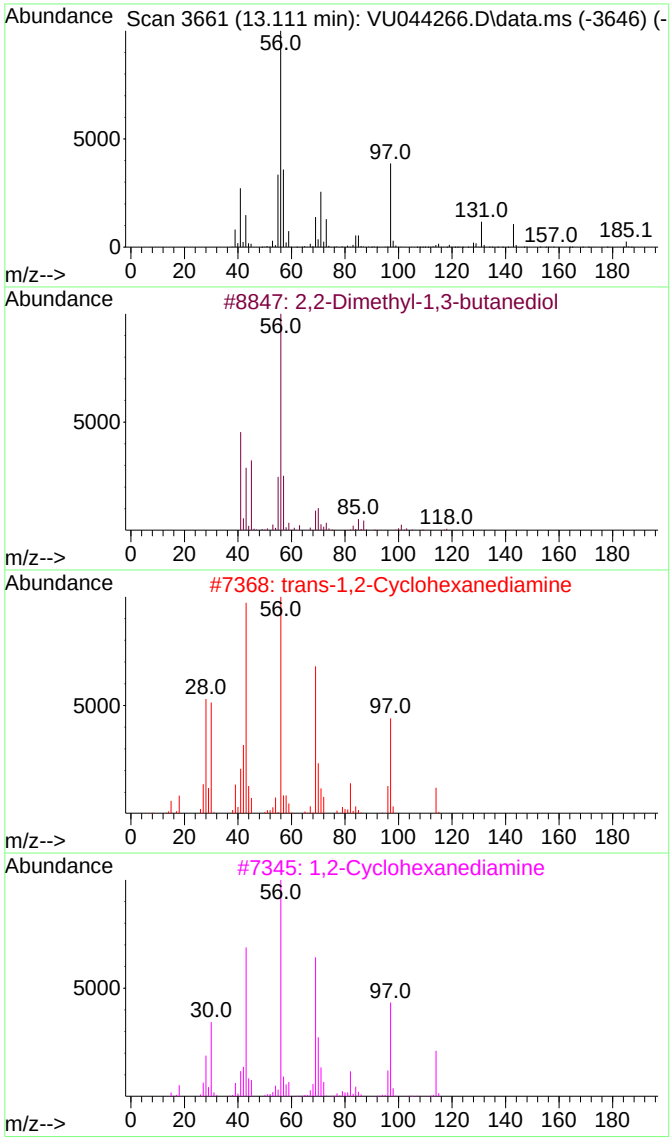
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR062121WMA.M
Quant Title : TRACE VOA SFAM1.0

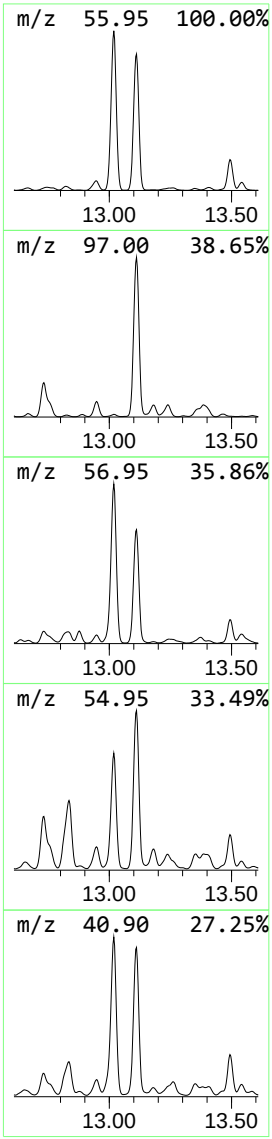
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 unknown-05 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.111	1.50 ug/L	7615360	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2-Dimethyl-1,3-butanediol	118	C6H14O2	000076-35-7	38
2	trans-1,2-Cyclohexanediamine	114	C6H14N2	001121-22-8	36
3	1,2-Cyclohexanediamine	114	C6H14N2	000694-83-7	36
4	1-Hexene, 2,5-dimethyl-	112	C8H16	006975-92-4	25
5	1-Hexene, 2,5-dimethyl-	112	C8H16	006975-92-4	25





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU063021\
Data File : VU044266.D
Acq On : 30 Jun 2021 17:39
Operator : SY/MD
Sample : M2905-10
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DBK33

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR062121WMA.M
Quant Title : TRACE VOA SFAM1.0

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

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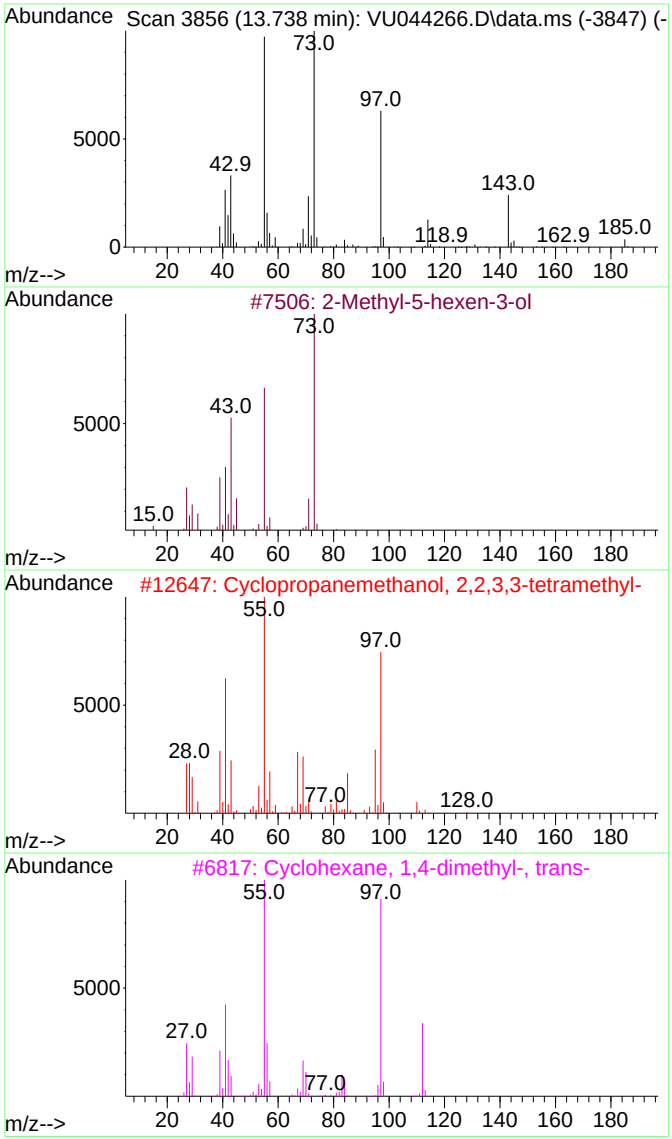
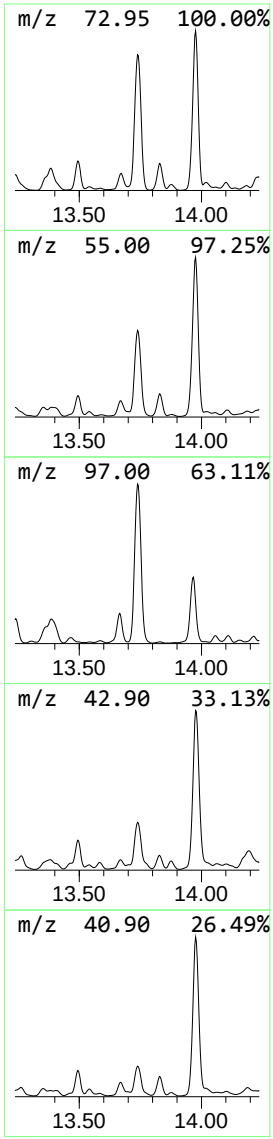
Peak Number 9 unknown-06 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.738	0.71 ug/L	3615810	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	27
2	Cyclopropanemethanol, 2,2,3,3-te...	128	C8H16O	002415-96-5	25
3	Cyclohexane, 1,4-dimethyl-, trans-	112	C8H16	002207-04-7	22
4	5-Octen-4-one, 7-methyl-	140	C9H16O	032064-78-1	17
5	Cyclohexane, 1,4-dimethyl-	112	C8H16	000589-90-2	16





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU063021\
Data File : VU044266.D
Acq On : 30 Jun 2021 17:39
Operator : SY/MD
Sample : M2905-10
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DBK33

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR062121WMA.M
Quant Title : TRACE VOA SFAM1.0

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

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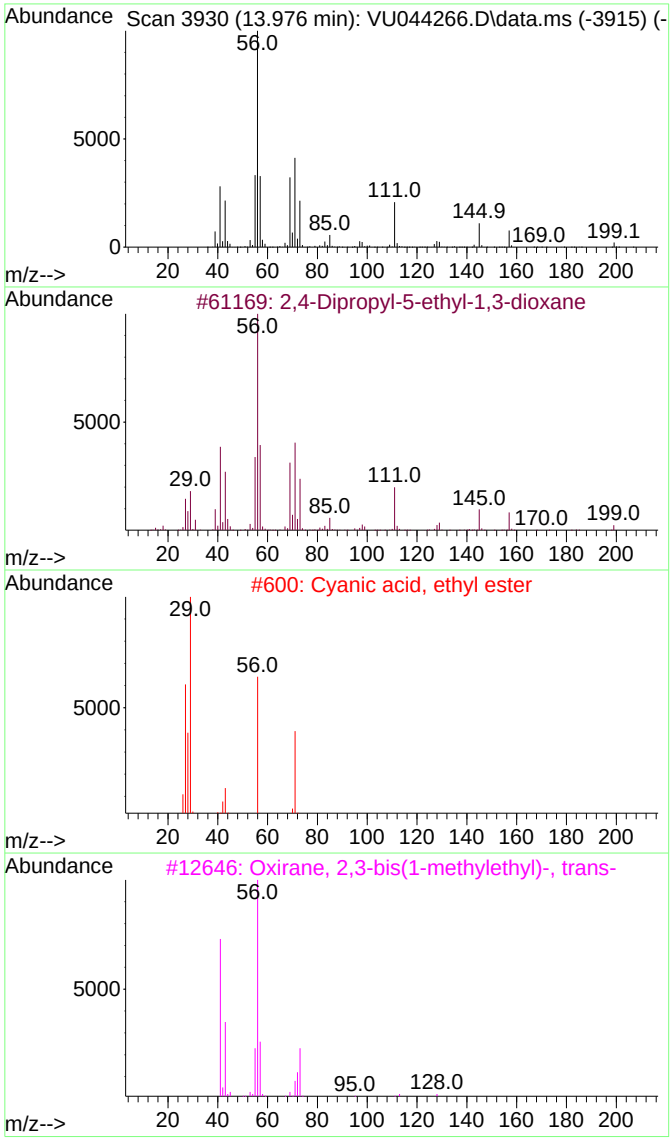
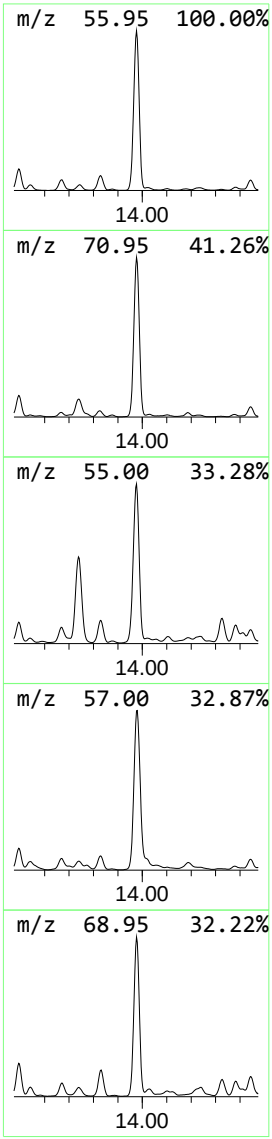
Peak Number 10 2,4-Dipropyl-5-ethyl-1,3-di... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.976	3.05 ug/L	15488100	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	72
2	Cyanic acid, ethyl ester	71	C3H5NO	000627-48-5	52
3	Oxirane, 2,3-bis(1-methylethyl)-...	128	C8H16O	054644-32-5	37
4	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	33
5	1,3-Pentanediol, 2,2,4-trimethyl-	146	C8H18O2	000144-19-4	28





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU063021\
Data File : VU044266.D
Acq On : 30 Jun 2021 17:39
Operator : SY/MD
Sample : M2905-10
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DBK33

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR062121WMA.M
Quant Title : TRACE VOA SFAM1.0

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

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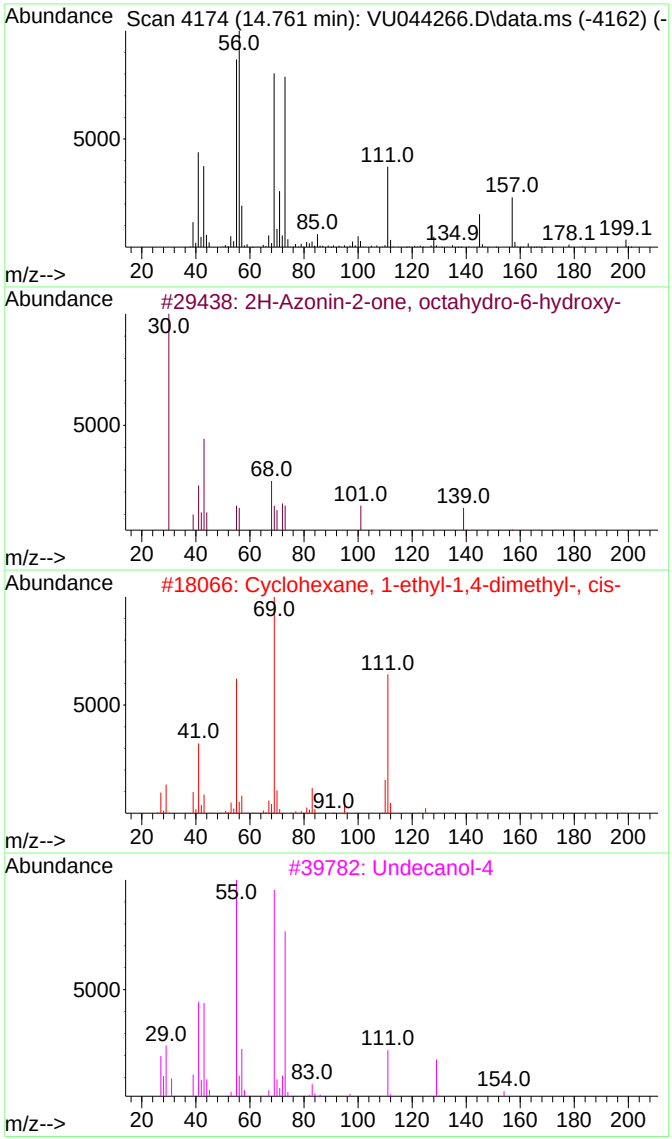
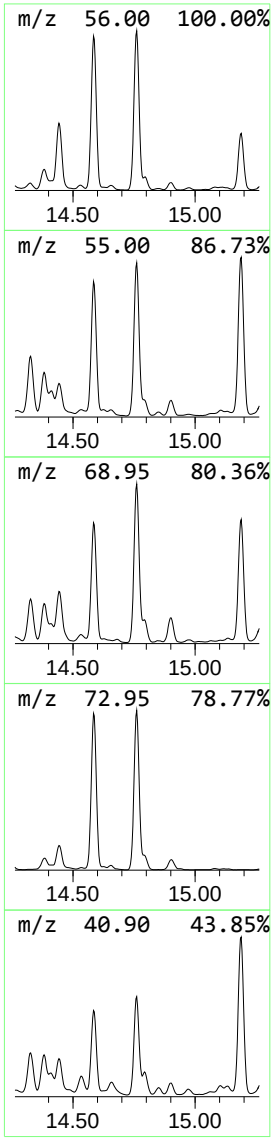
Peak Number 11 unknown-07 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.761	0.65 ug/L	3285160	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Azonin-2-one, octahydro-6-hyd...	157	C8H15NO2	023435-07-6	38
2	Cyclohexane, 1-ethyl-1,4-dimethy...	140	C10H20	062238-30-6	35
3	Undecanol-4	172	C11H24O	004272-06-4	35
4	2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	32
5	Neopentyl glycol	104	C5H12O2	000126-30-7	25





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU063021\
Data File : VU044266.D
Acq On : 30 Jun 2021 17:39
Operator : SY/MD
Sample : M2905-10
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DBK33

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR062121WMA.M
Quant Title : TRACE VOA SFAM1.0

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

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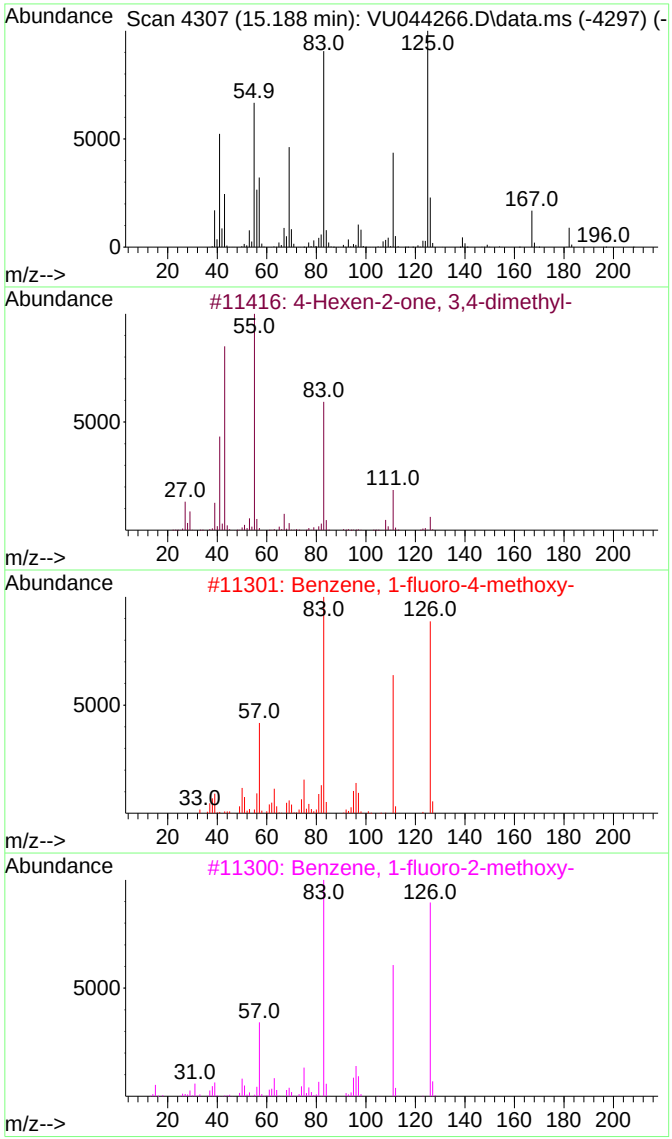
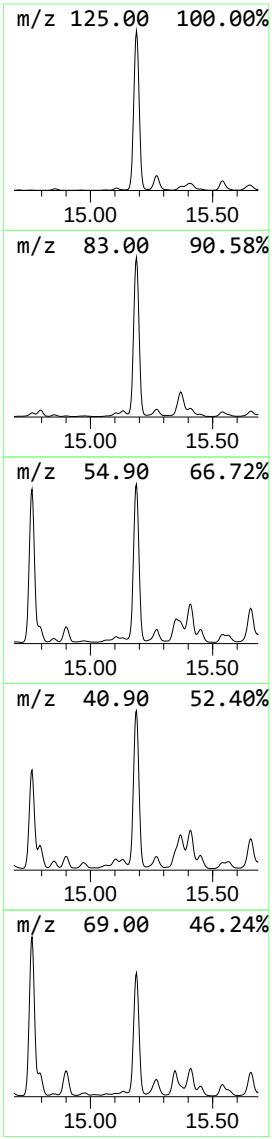
Peak Number 12 unknown-08 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.188	0.86 ug/L	4353090	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Hexen-2-one, 3,4-dimethyl-	126	C8H14O	053252-21-4	46
2	Benzene, 1-fluoro-4-methoxy-	126	C7H7FO	000459-60-9	43
3	Benzene, 1-fluoro-2-methoxy-	126	C7H7FO	000321-28-8	43
4	5-Amino-1-ethylpyrazole	111	C5H9N3	003528-58-3	35
5	Cyclohexanecarboxylic acid, (1H-...	195	C8H13N5O	1000310-78-1	35





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU063021\
Data File : VU044266.D
Acq On : 30 Jun 2021 17:39
Operator : SY/MD
Sample : M2905-10
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DBK33

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR062121WMA.M
Quant Title : TRACE VOA SFAM1.0

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Propanal, 2-met...	3.591	6.3	ug/L	406417	1	6.256	324785	5.0
Butanal	4.459	42.8	ug/L	2778740	1	6.256	324785	5.0
unknown-01	9.690	20.6	ug/L	1512720	2	9.423	368032	5.0
unknown-02	10.709	0.6	ug/L	2827400	3	11.819	25366100	5.0
(DEL) Alkane: C...	12.050	0.7	ug/L	3279740	3	11.819	25366100	5.0
unknown-03	12.835	0.5	ug/L	2571640	3	11.819	25366100	5.0
unknown-04	13.018	1.6	ug/L	8327850	3	11.819	25366100	5.0
unknown-05	13.111	1.5	ug/L	7615360	3	11.819	25366100	5.0
unknown-06	13.738	0.7	ug/L	3615810	3	11.819	25366100	5.0
2,4-Dipropyl-5-...	13.976	3.0	ug/L	15488100	3	11.819	25366100	5.0
unknown-07	14.761	0.7	ug/L	3285160	3	11.819	25366100	5.0
unknown-08	15.188	0.9	ug/L	4353090	3	11.819	25366100	5.0