LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV101723\
Data File : VV032440.D
Acq On : 17 Oct 2023 13:04
Operator : SY/MD
Sample : 04903-04
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : C0AE7

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\SFAMVTR092923WMA.M
Title : TRACE VOA SFAM1.0

Signal : TIC: VV032440.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.095	10	17	34	rVB2	211376	258807	71.10%	4.626%
2	1.169	35	40	46	rVB	7939	6894	1.89%	0.123%
3	1.204	46	51	56	rBV	15069	13787	3.79%	0.246%
4	1.298	71	80	94	rBV2	112439	169617	46.60%	3.032%
5	1.417	107	117	131	rBV4	33065	67629	18.58%	1.209%
6	1.587	162	170	186	rVB	46650	67307	18.49%	1.203%
7	1.661	187	193	200	rVB4	7430	9340	2.57%	0.167%
8	1.796	220	235	236	rBV6	27747	42549	11.69%	0.761%
9	1.822	237	243	256	rVB6	456748	359591	98.78%	6.427%
10	2.140	334	342	356	rBV	83602	127234	34.95%	2.274%
11	2.330	396	401	414	rVB	21796	33048	9.08%	0.591%
12	3.725	822	835	851	rVB5	14776	46559	12.79%	0.832%
13	4.346	1015	1028	1056	rVB	64403	167614	46.05%	2.996%
14	4.593	1092	1105	1119	rBV3	6971	17382	4.77%	0.311%
15	4.741	1143	1151	1165	rVB5	3200	7055	1.94%	0.126%
16	5.047	1228	1246	1268	rBV2	139123	343257	94.30%	6.135%
17	5.246	1292	1308	1326	rVB5	10465	29630	8.14%	0.530%
18	5.612	1410	1422	1440	rBV	129578	263443	72.37%	4.709%
19	5.767	1460	1470	1482	rVB4	6021	13120	3.60%	0.235%
20	5.895	1500	1510	1525	rVB4	8183	20845	5.73%	0.373%
21	6.059	1546	1561	1584	rVB2	78807	164494	45.19%	2.940%
22	6.185	1590	1600	1609	rBV5	7238	14228	3.91%	0.254%
23	6.297	1625	1635	1654	rBV4	12133	30973	8.51%	0.554%
24	6.963	1832	1842	1858	rBV	36310	68538	18.83%	1.225%
25	7.056	1863	1871	1881	rVB	2909	5888	1.62%	0.105%
26	7.236	1914	1927	1931	rBV6	4593	9416	2.59%	0.168%
27	7.284	1931	1942	1958	rVV	152040	274680	75.46%	4.910%
28	7.365	1958	1967	1977	rVB3	8433	17787	4.89%	0.318%
29	7.593	2029	2038	2053	rBV	20518	39820	10.94%	0.712%
30	7.902	2123	2134	2138	rBV2	20030	32584	8.95%	0.582%
31	7.934	2140	2144	2156	rVB3	25791	40953	11.25%	0.732%
32	8.063	2173	2184	2196	rBV	182414	364022	100.00%	6.507%
33	8.207	2219	2229	2254	rVB2	68192	132249	36.33%	2.364%
34	8.805	2404	2415	2441	rBV	199977	356397	97.91%	6.370%
35	9.095	2496	2505	2517	rBV2	12351	21721	5.97%	0.388%
36	9.188	2522	2534	2552	rBV4	11503	27342	7.51%	0.489%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV101723\
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Operator : SY/MD
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ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AE7

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

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Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR092923WMA.M
Title : TRACE VOA SFAM1.0

37	9.345	2576	2583	2604	rVB5	6822	15548	4.27%	0.278%
38	9.497	2620	2630	2637	rBV4	6925	11551	3.17%	0.206%
39	9.673	2676	2685	2697	rBV4	8507	15963	4.39%	0.285%
40	9.763	2703	2713	2736	rBV2	54368	104577	28.73%	1.869%
41	10.165	2828	2838	2853	rVB	58639	99598	27.36%	1.780%
42	10.333	2877	2890	2902	rVB2	72573	125330	34.43%	2.240%
43	10.458	2921	2929	2947	rVB4	26028	53033	14.57%	0.948%
44	10.554	2950	2959	2973	rVB4	6556	16753	4.60%	0.299%
45	10.670	2983	2995	3004	rBV4	4998	11899	3.27%	0.213%
46	10.860	3047	3054	3064	rBV2	12666	20414	5.61%	0.365%
47	10.995	3087	3096	3104	rBV3	7304	12470	3.43%	0.223%
48	11.101	3119	3129	3148	rBV2	53765	94641	26.00%	1.692%
49	11.194	3148	3158	3173	rVV	226823	356208	97.85%	6.367%
50	11.490	3233	3250	3265	rBV2	71115	173750	47.73%	3.106%
51	11.567	3265	3274	3303	rVB	182771	343288	94.30%	6.136%
52	11.776	3333	3339	3349	rVB3	16704	28847	7.92%	0.516%
53	11.860	3358	3365	3368	rBV6	5889	8408	2.31%	0.150%
54	11.963	3390	3397	3403	rBV	9511	12891	3.54%	0.230%
55	12.197	3459	3470	3482	rBV2	76371	130313	35.80%	2.329%
56	12.291	3490	3499	3514	rBV	64840	102973	28.29%	1.841%
57	12.358	3514	3520	3529	rVV	8779	14630	4.02%	0.262%
58	12.410	3529	3536	3547	rVB2	13463	22426	6.16%	0.401%
59	12.673	3611	3618	3622	rBV7	7948	11309	3.11%	0.202%
60	12.795	3646	3656	3659	rBV9	7224	10641	2.92%	0.190%
61	12.828	3659	3666	3677	rVB4	10121	18611	5.11%	0.333%
62	13.088	3739	3747	3762	rBV9	11698	24938	6.85%	0.446%
63	13.213	3778	3786	3794	rBV9	7375	11372	3.12%	0.203%
64	13.384	3832	3839	3846	rBV8	7659	12038	3.31%	0.215%
65	13.448	3853	3859	3868	rBV	6493	8654	2.38%	0.155%
66	13.988	4014	4027	4037	rBV	24613	41916	11.51%	0.749%
67	15.586	4515	4524	4529	rBV4	3844	6503	1.79%	0.116%
68	17.046	4972	4978	4987	rVB2	5721	7318	2.01%	0.131%

Sum of corrected areas: 5594611

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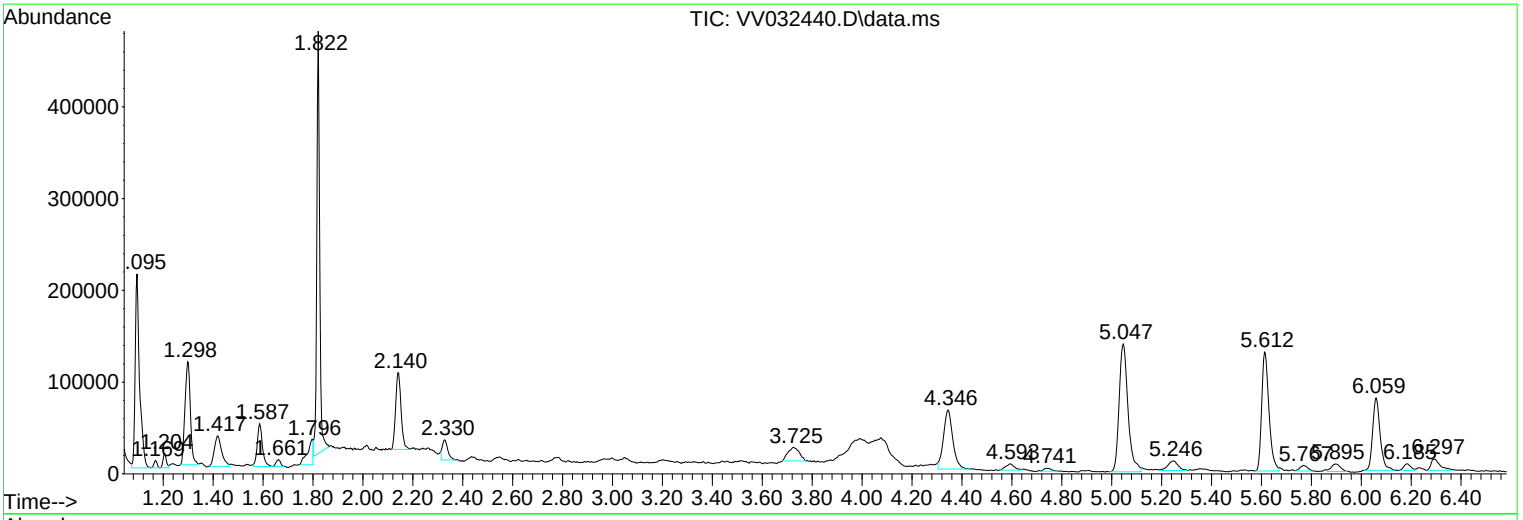
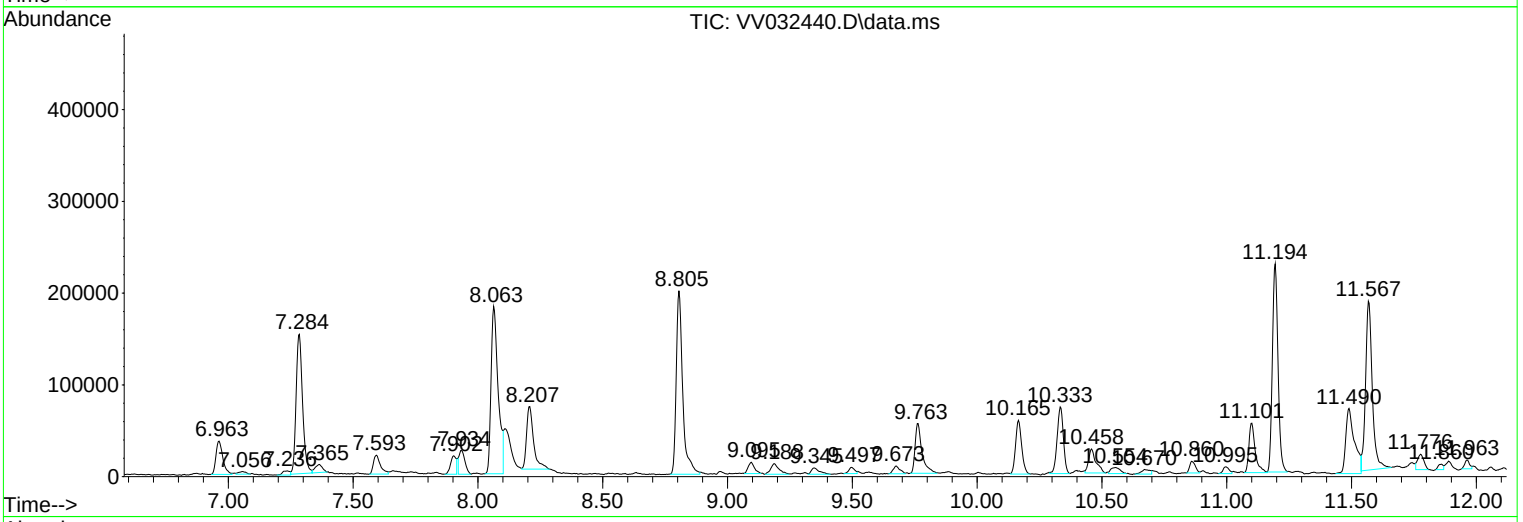
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 6 Sample Multiplier: 1

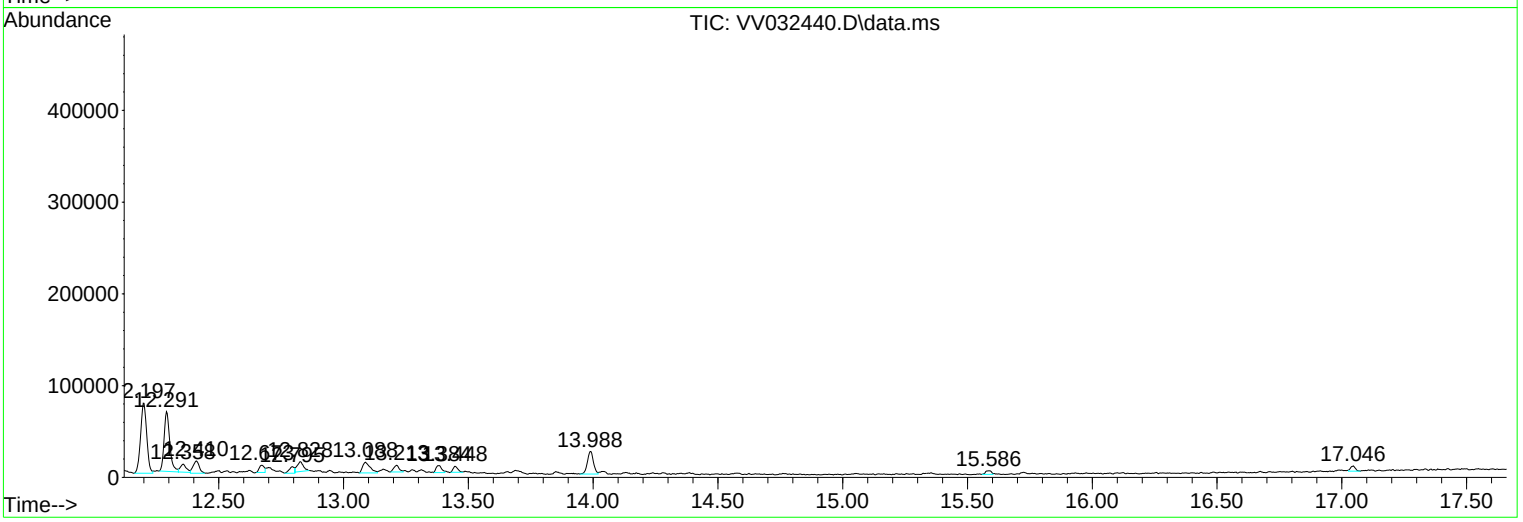
Instrument :
MSVOA_V
ClientSampleId :
C0AE7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR092923WMA.M
Quant Title : TRACE VOA SFAM1.0

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







Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV101723\
Data File : VV032440.D
Acq On : 17 Oct 2023 13:04
Operator : SY/MD
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ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AE7

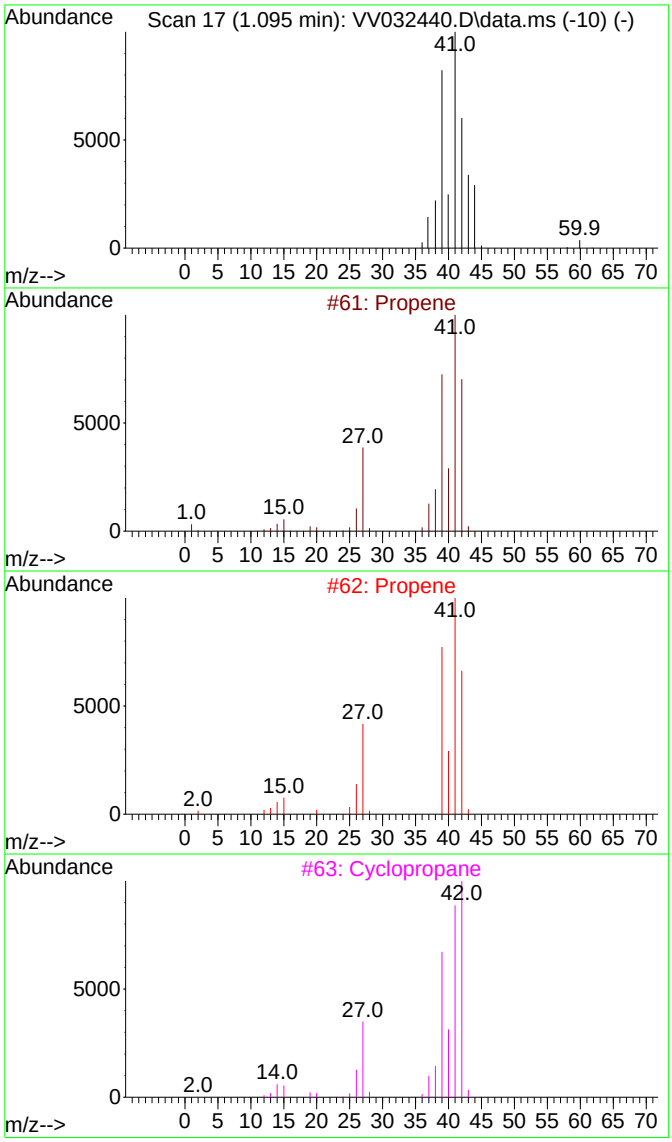
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR092923WMA.M
Quant Title : TRACE VOA SFAM1.0

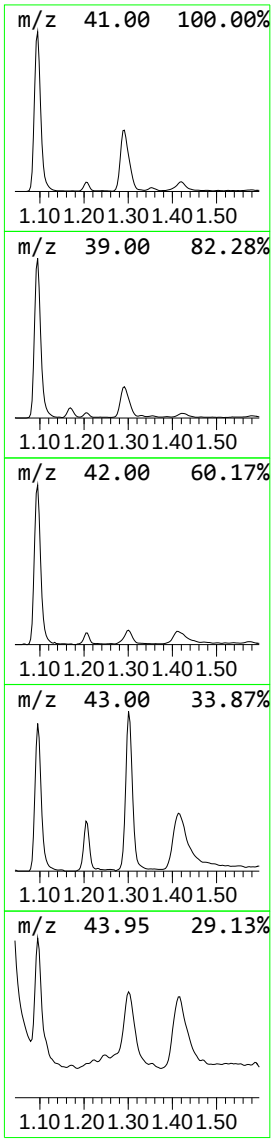
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.095	4.91 ug/L	258807	1,4-Difluorobenzene	5.612

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propene	42	C3H6	000115-07-1	90
2	Propene	42	C3H6	000115-07-1	42
3	Cyclopropane	42	C3H6	000075-19-4	40
4	Cyclopropane	42	C3H6	000075-19-4	9
5	Cyclopropene	40	C3H4	002781-85-3	9





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV101723\
Data File : VV032440.D
Acq On : 17 Oct 2023 13:04
Operator : SY/MD
Sample : 04903-04
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AE7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR092923WMA.M
Quant Title : TRACE VOA SFAM1.0

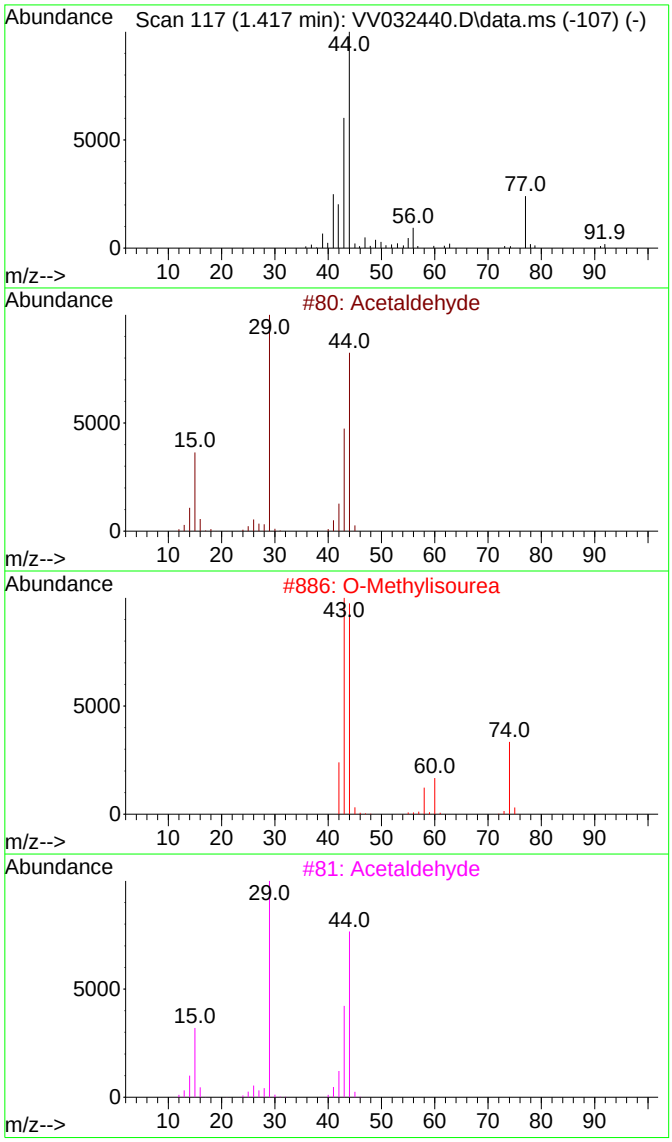
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

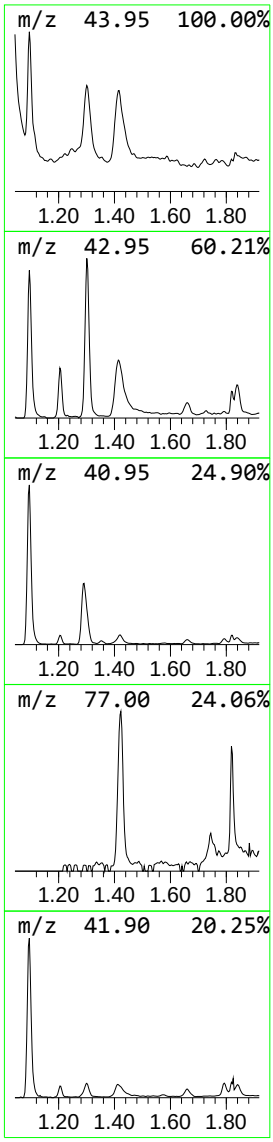
****

Peak Number 2 unknown-01 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.417	1.28 ug/L	67629	1,4-Difluorobenzene	5.612

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetaldehyde	44	C2H4O	000075-07-0	9
2	O-Methylisourea	74	C2H6N2O	002440-60-0	9
3	Acetaldehyde	44	C2H4O	000075-07-0	7
4	Ethylene oxide	44	C2H4O	000075-21-8	5
5	1,2-Propanediamine	74	C3H10N2	000078-90-0	5





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV101723\
Data File : VV032440.D
Acq On : 17 Oct 2023 13:04
Operator : SY/MD
Sample : 04903-04
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AE7

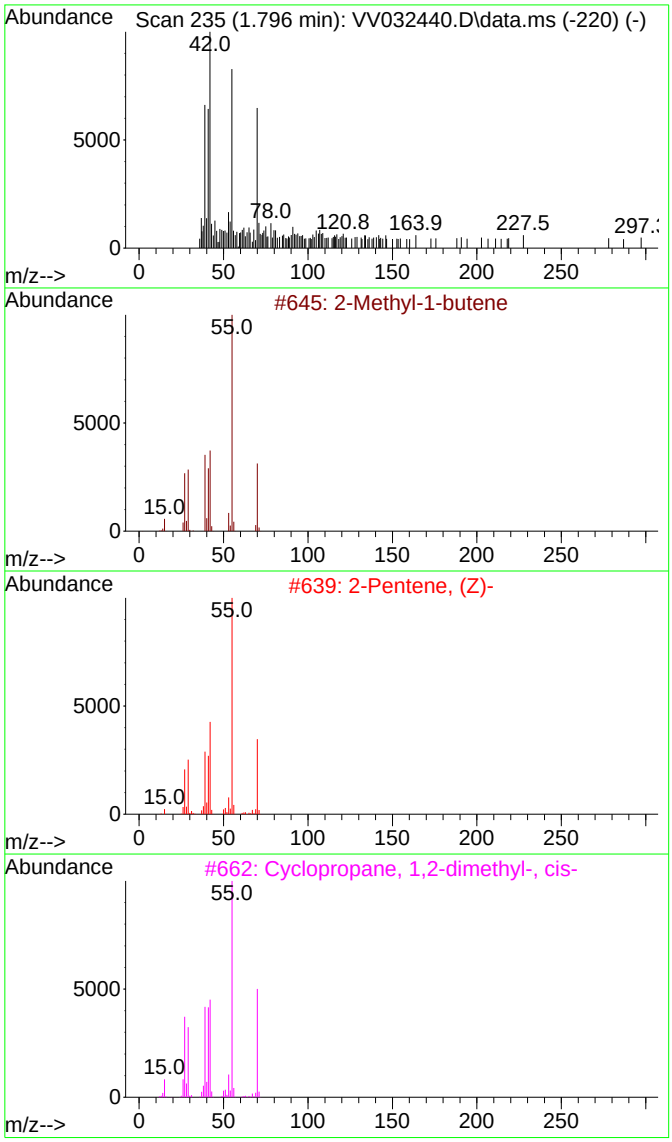
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR092923WMA.M
Quant Title : TRACE VOA SFAM1.0

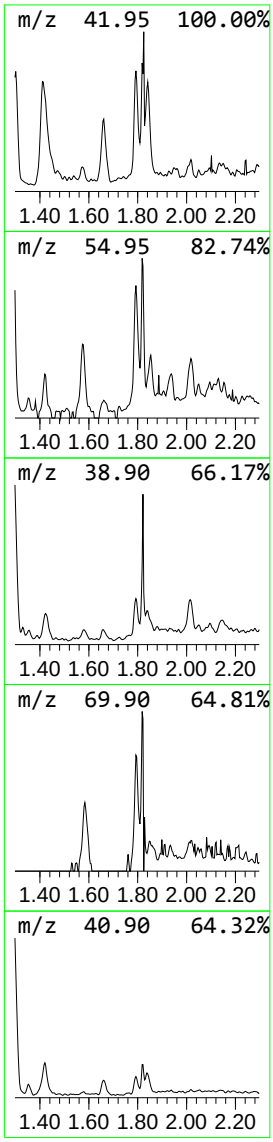
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.796	0.81 ug/L	42549	1,4-Difluorobenzene	5.612

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methyl-1-butene	70	C5H10	000563-46-2	38
2	2-Pentene, (Z)-	70	C5H10	000627-20-3	38
3	Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	38
4	Pentane, 1-chloro-	106	C5H11Cl	000543-59-9	35
5	3-Ethyltetrahydrofuran	100	C6H12O	093716-28-0	35





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV101723\
Data File : VV032440.D
Acq On : 17 Oct 2023 13:04
Operator : SY/MD
Sample : 04903-04
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AE7

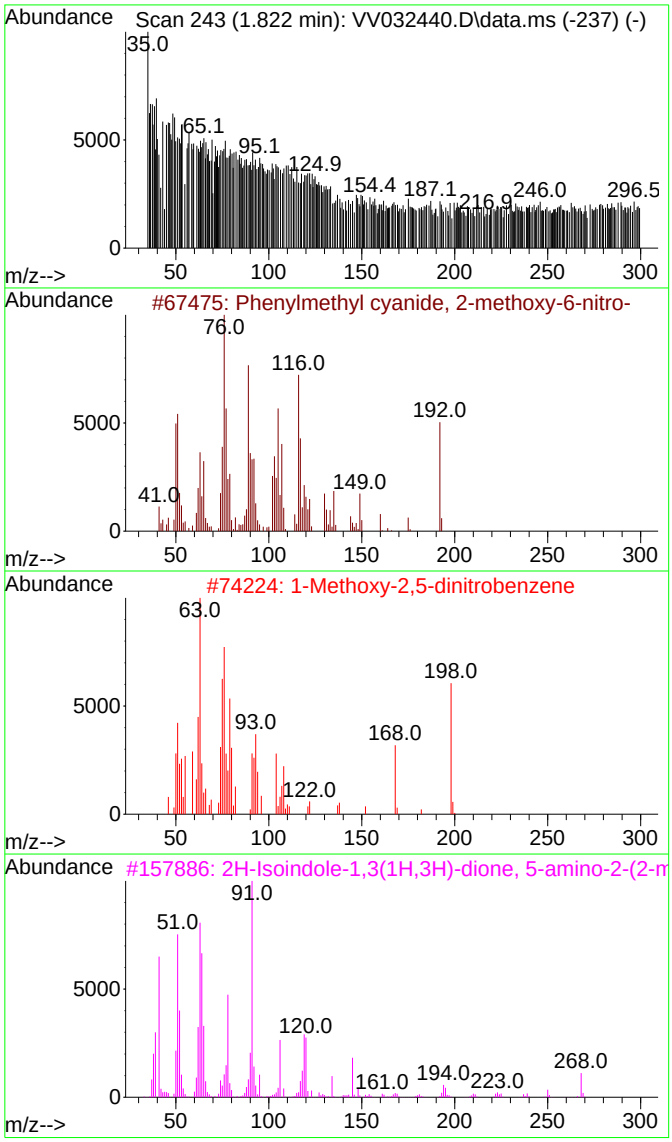
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR092923WMA.M
Quant Title : TRACE VOA SFAM1.0

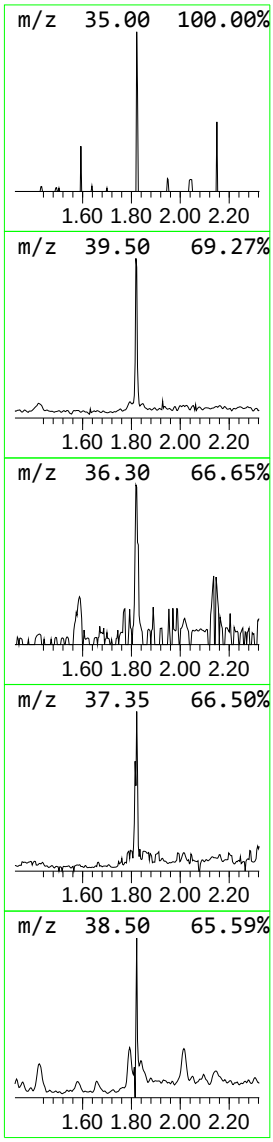
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 unknown-02 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.822	6.82 ug/L	359591	1,4-Difluorobenzene	5.612

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenylmethyl cyanide, 2-methoxy-...	192	C9H8N2O3	020876-27-1	48
2	1-Methoxy-2,5-dinitrobenzene	198	C7H6N2O5	1000362-87-5	38
3	2H-Isoindole-1,3(1H,3H)-dione, 5...	268	C15H12N2O3	313515-42-3	35
4	1,2-Benzisoxazole	119	C7H5NO	000271-95-4	35
5	4-(2-Hydroxyethylamino)-3-nitroc...	250	C11H10N2O5	088353-20-2	30





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV101723\
Data File : VV032440.D
Acq On : 17 Oct 2023 13:04
Operator : SY/MD
Sample : 04903-04
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AE7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR092923WMA.M
Quant Title : TRACE VOA SFAM1.0

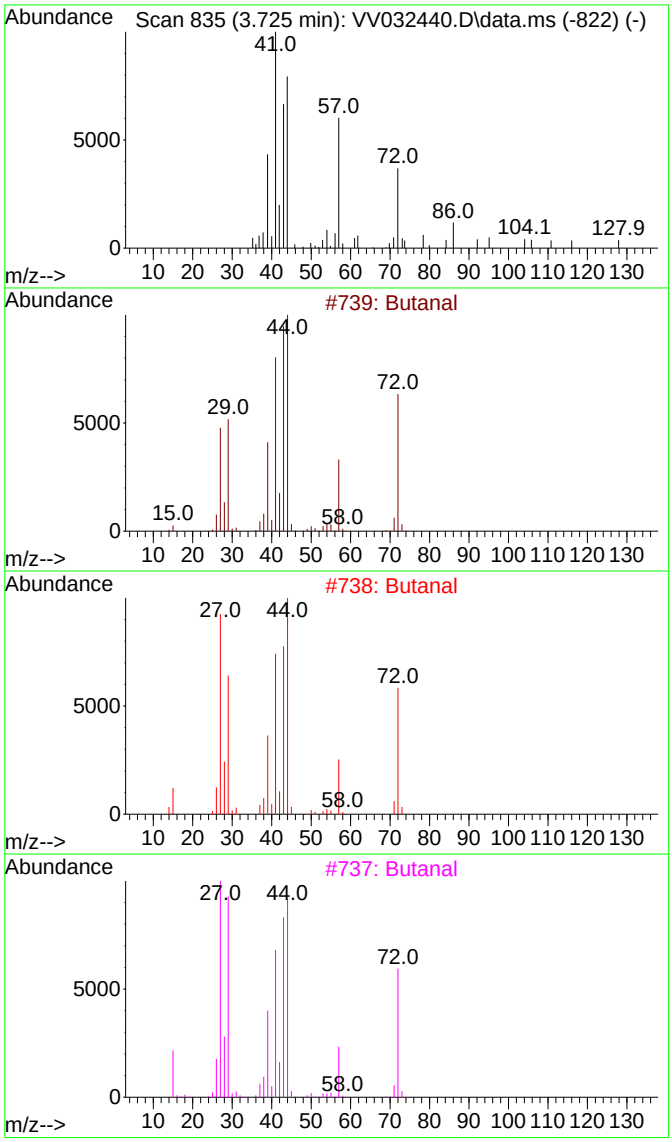
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

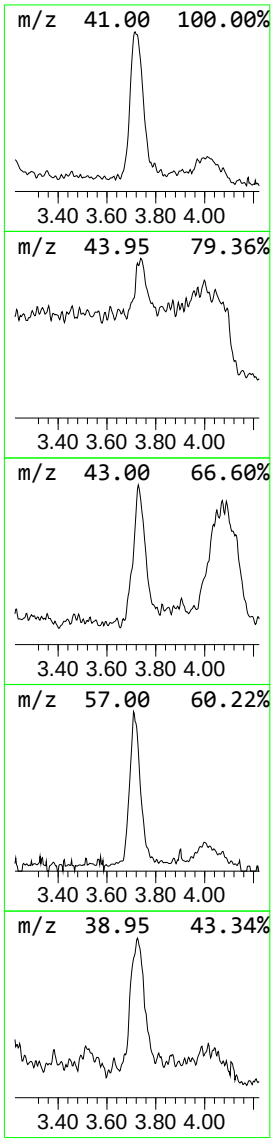
****

Peak Number 5 unknown-03 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.725	0.88 ug/L	46559	1,4-Difluorobenzene	5.612

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanal	72	C4H8O	000123-72-8	46
2	Butanal	72	C4H8O	000123-72-8	43
3	Butanal	72	C4H8O	000123-72-8	43
4	Ethene, ethoxy-	72	C4H8O	000109-92-2	32
5	Propanamide, 2-methyl-	87	C4H9NO	000563-83-7	28





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Data File : VV032440.D
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ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AE7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR092923WMA.M
Quant Title : TRACE VOA SFAM1.0

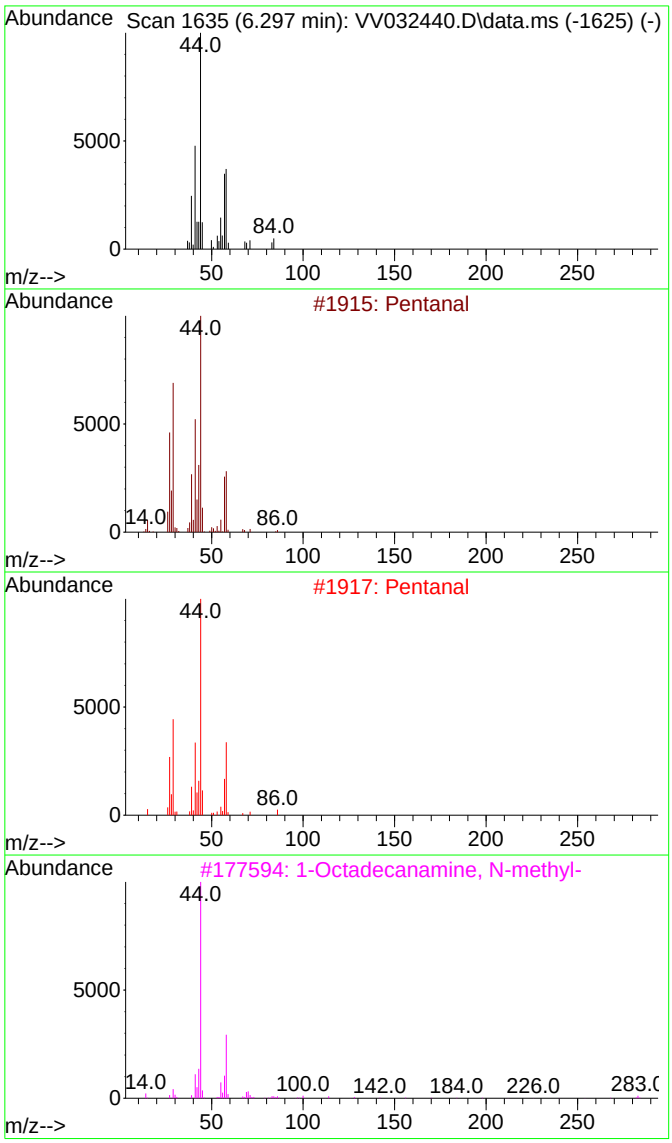
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

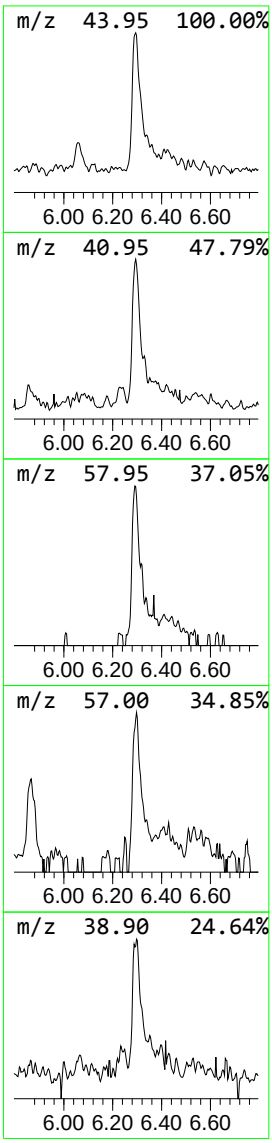
****

Peak Number 7 Pentanal Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.297	0.59 ug/L	30973	1,4-Difluorobenzene	5.612

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentanal	86	C5H10O	000110-62-3	72
2	Pentanal	86	C5H10O	000110-62-3	64
3	1-Octadecanamine, N-methyl-	283	C19H41N	002439-55-6	9
4	Cyclopropyl carbinol	72	C4H8O	002516-33-8	9
5	Cyclopropyl carbinol	72	C4H8O	002516-33-8	9





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV101723\
Data File : VV032440.D
Acq On : 17 Oct 2023 13:04
Operator : SY/MD
Sample : 04903-04
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AE7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR092923WMA.M
Quant Title : TRACE VOA SFAM1.0

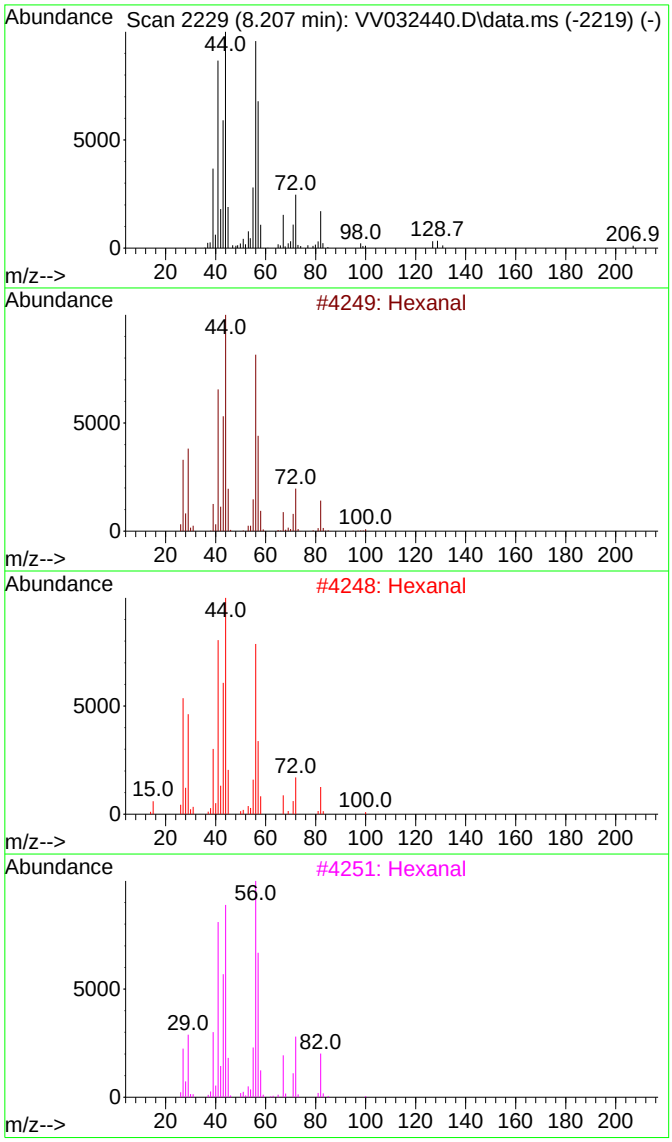
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

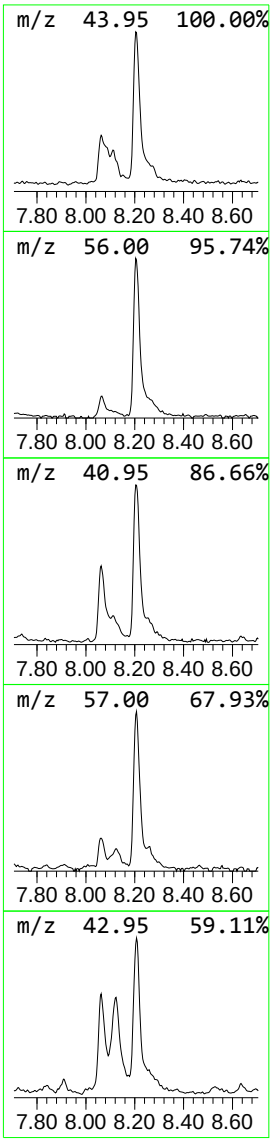
****

Peak Number 9 Hexanal Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.207	1.86 ug/L	132249	Chlorobenzene-d5	8.805

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanal	100	C6H12O	000066-25-1	91
2	Hexanal	100	C6H12O	000066-25-1	90
3	Hexanal	100	C6H12O	000066-25-1	90
4	Hexanal	100	C6H12O	000066-25-1	83
5	Hexanal	100	C6H12O	000066-25-1	78





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV101723\
Data File : VV032440.D
Acq On : 17 Oct 2023 13:04
Operator : SY/MD
Sample : 04903-04
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AE7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR092923WMA.M
Quant Title : TRACE VOA SFAM1.0

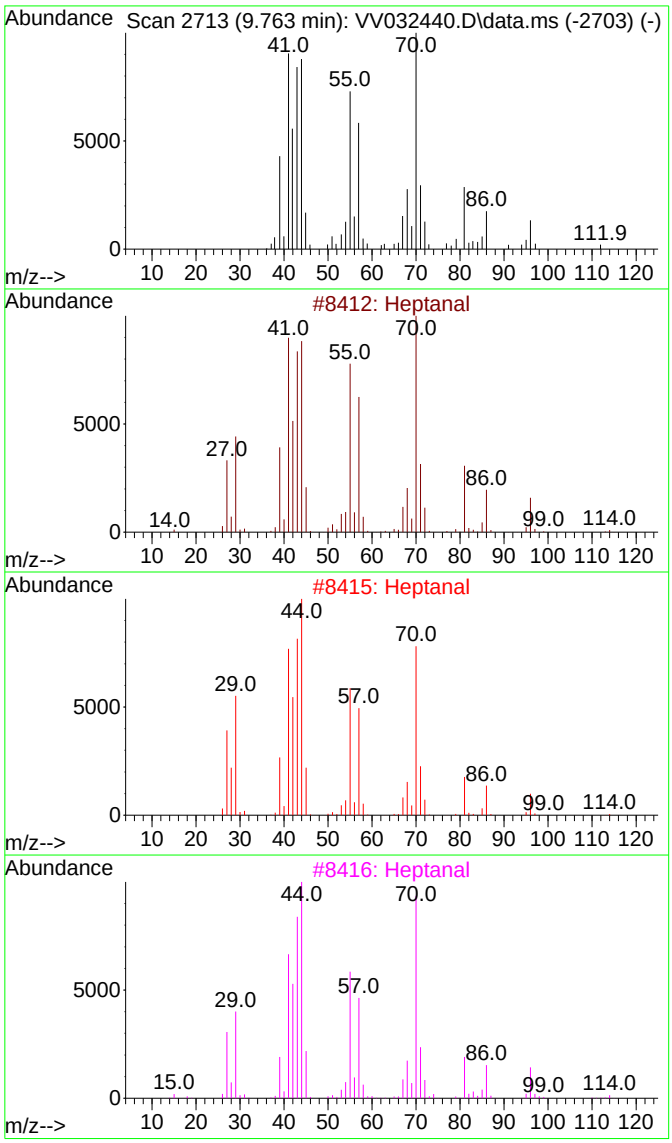
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

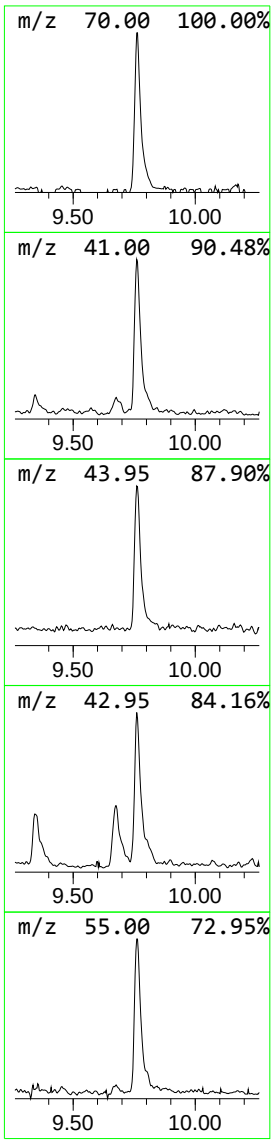
****

Peak Number 10 Heptanal Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.763	1.47 ug/L	104577	Chlorobenzene-d5	8.805

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptanal	114	C7H14O	000111-71-7	90
2	Heptanal	114	C7H14O	000111-71-7	72
3	Heptanal	114	C7H14O	000111-71-7	64
4	Heptanal	114	C7H14O	000111-71-7	64
5	Heptanal	114	C7H14O	000111-71-7	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV101723\
Data File : VV032440.D
Acq On : 17 Oct 2023 13:04
Operator : SY/MD
Sample : 04903-04
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AE7

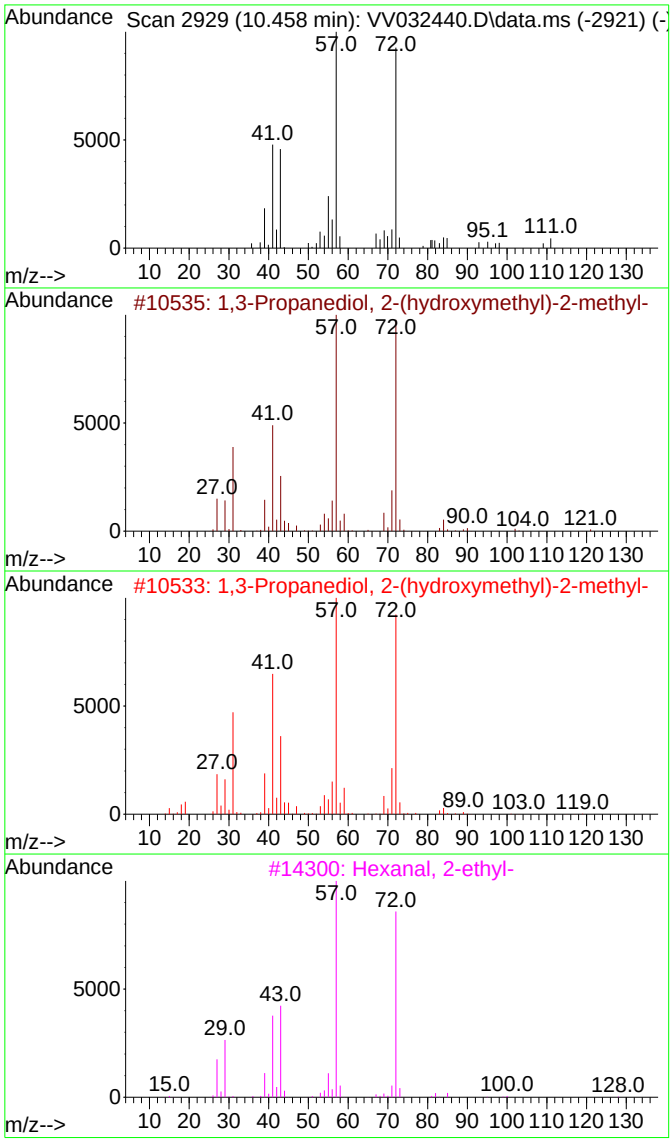
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR092923WMA.M
Quant Title : TRACE VOA SFAM1.0

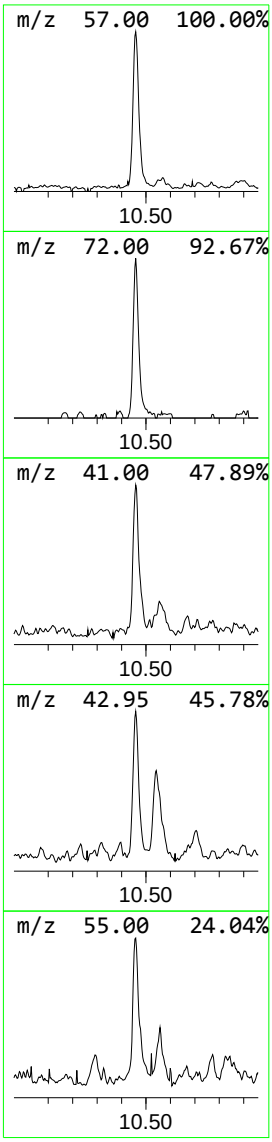
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 1,3-Propanediol, 2-(hydroxy... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.458	0.74 ug/L	53033	1,4-Dichlorobenzene-d4	11.194

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Propanediol, 2-(hydroxymethy...	120	C5H12O3	000077-85-0	72
2	1,3-Propanediol, 2-(hydroxymethy...	120	C5H12O3	000077-85-0	72
3	Hexanal, 2-ethyl-	128	C8H16O	000123-05-7	64
4	Hexanal, 2-ethyl-	128	C8H16O	000123-05-7	64
5	2-Propanone, hydrazone	72	C3H8N2	005281-20-9	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV101723\
Data File : VV032440.D
Acq On : 17 Oct 2023 13:04
Operator : SY/MD
Sample : 04903-04
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AE7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR092923WMA.M
Quant Title : TRACE VOA SFAM1.0

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

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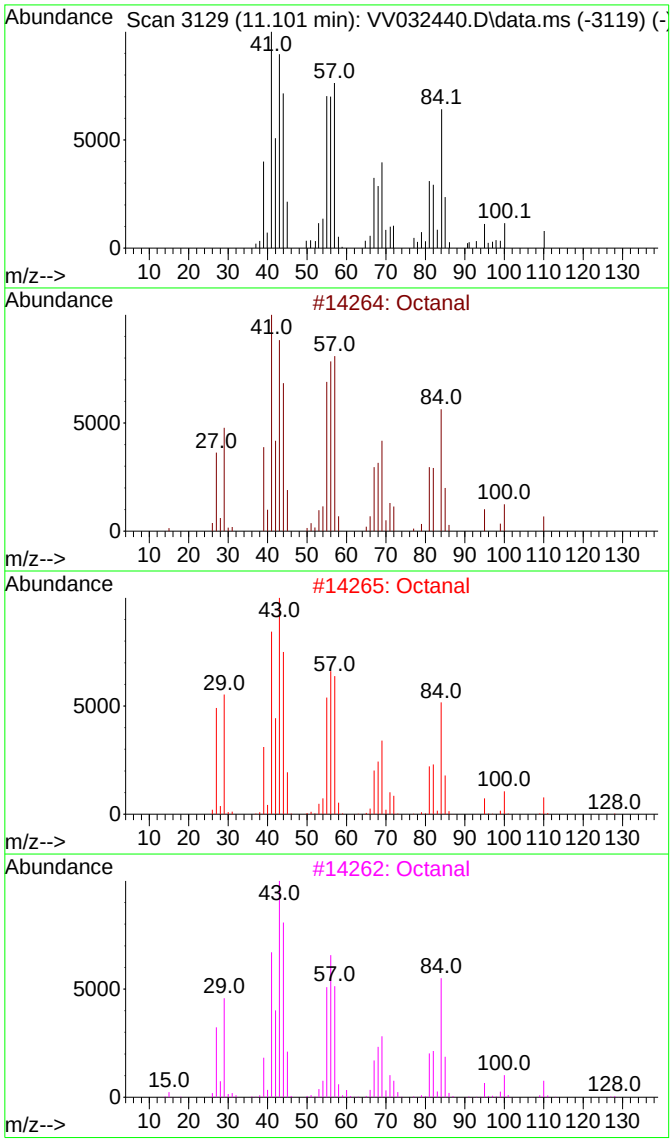
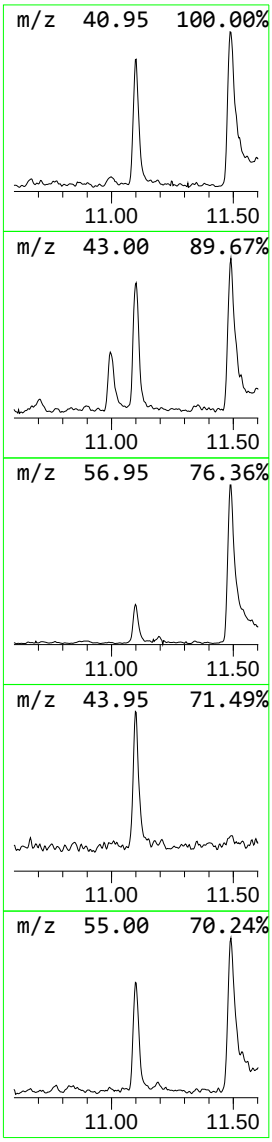
Peak Number 13 Octanal Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.101	1.33 ug/L	94641	1,4-Dichlorobenzene-d4	11.194

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octanal	128	C8H16O	000124-13-0	91
2	Octanal	128	C8H16O	000124-13-0	91
3	Octanal	128	C8H16O	000124-13-0	90
4	Octanal	128	C8H16O	000124-13-0	90
5	Octanal	128	C8H16O	000124-13-0	53





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV101723\
Data File : VV032440.D
Acq On : 17 Oct 2023 13:04
Operator : SY/MD
Sample : 04903-04
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AE7

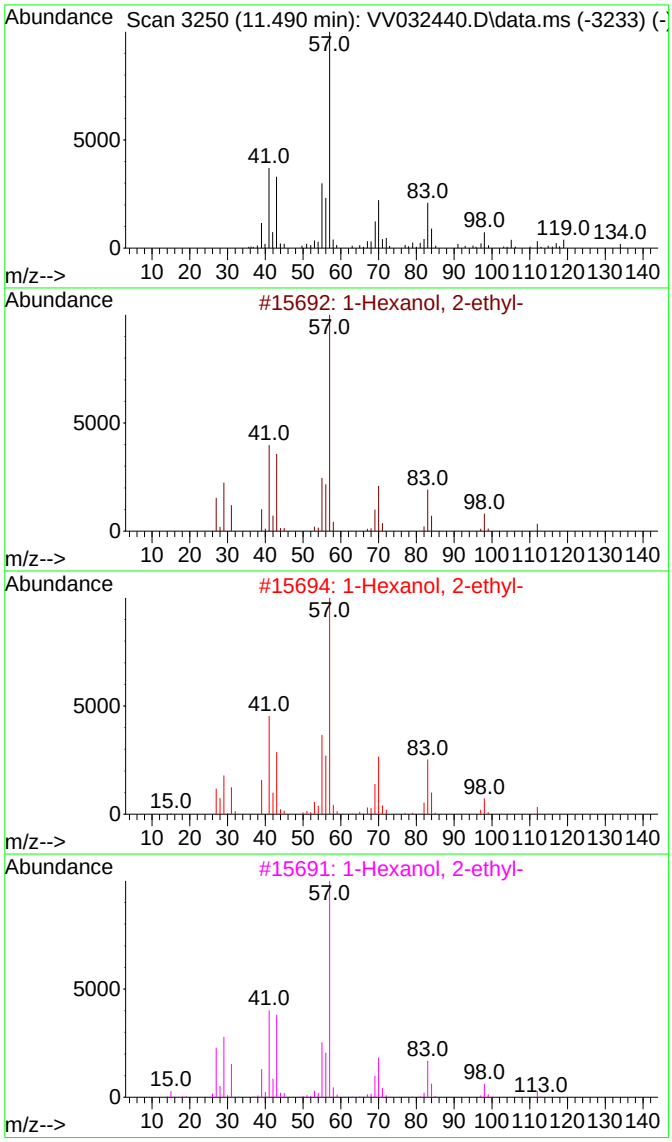
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR092923WMA.M
Quant Title : TRACE VOA SFAM1.0

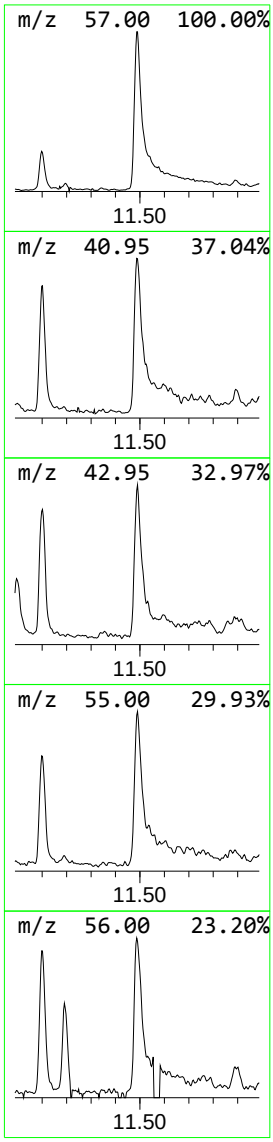
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 1-Hexanol, 2-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.490	2.44 ug/L	173750	1,4-Dichlorobenzene-d4	11.194

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	86
2	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	86
3	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	86
4	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
5	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV101723\
Data File : VV032440.D
Acq On : 17 Oct 2023 13:04
Operator : SY/MD
Sample : 04903-04
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AE7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR092923WMA.M
Quant Title : TRACE VOA SFAM1.0

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

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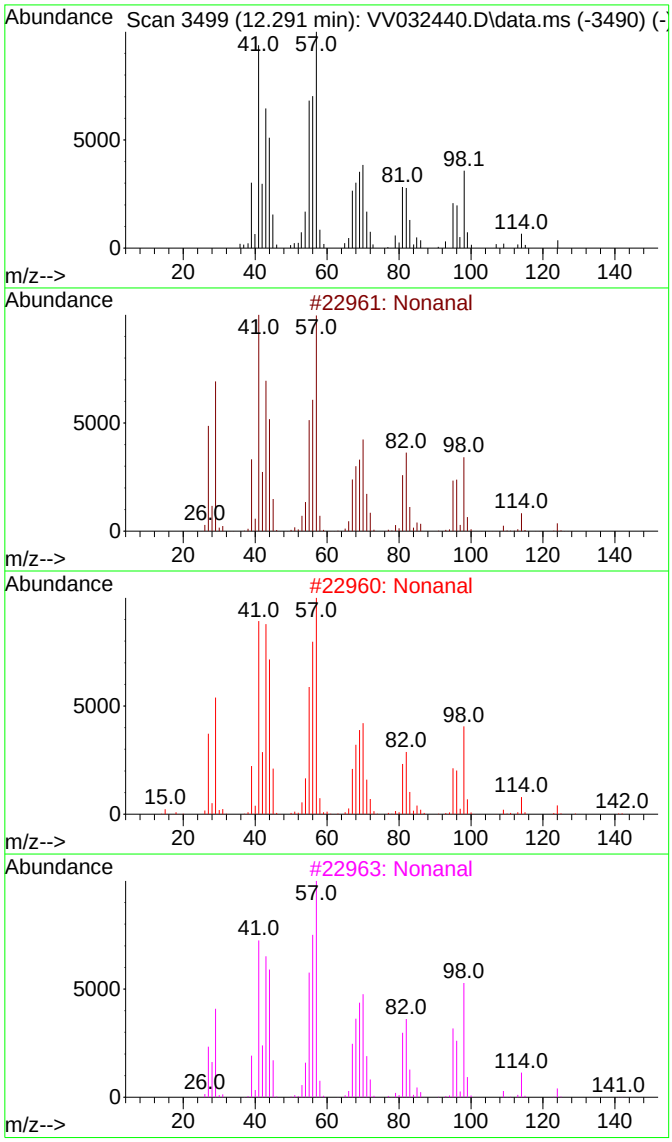
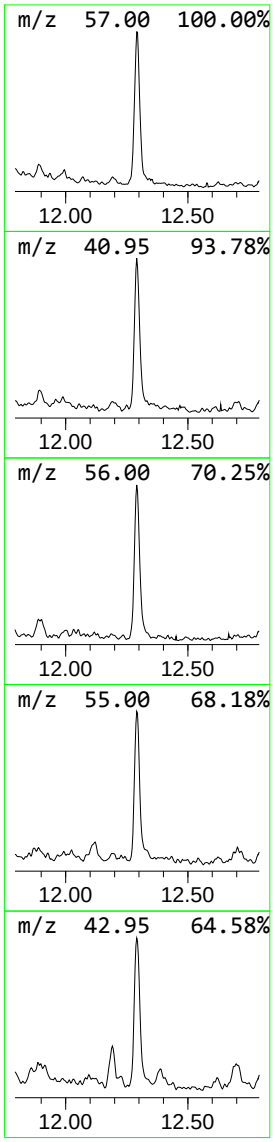
Peak Number 16 Nonanal Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.291	1.45 ug/L	102973	1,4-Dichlorobenzene-d4	11.194

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonanal	142	C9H18O	000124-19-6	91
2	Nonanal	142	C9H18O	000124-19-6	91
3	Nonanal	142	C9H18O	000124-19-6	90
4	Nonanal	142	C9H18O	000124-19-6	80
5	Bromoacetic acid, 10-undecenyl e...	290	C13H23BrO2	195373-65-0	25





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV101723\
Data File : VV032440.D
Acq On : 17 Oct 2023 13:04
Operator : SY/MD
Sample : 04903-04
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AE7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR092923WMA.M
Quant Title : TRACE VOA SFAM1.0

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
(DEL) Alkane: S...	1.095	4.9	ug/L	258807	1	5.612	263443	5.0
unknown-01	1.417	1.3	ug/L	67629	1	5.612	263443	5.0
(DEL) Alkane: S...	1.796	0.8	ug/L	42549	1	5.612	263443	5.0
unknown-02	1.822	6.8	ug/L	359591	1	5.612	263443	5.0
unknown-03	3.725	0.9	ug/L	46559	1	5.612	263443	5.0
Pentanal	6.297	0.6	ug/L	30973	1	5.612	263443	5.0
Hexanal	8.207	1.9	ug/L	132249	2	8.805	356397	5.0
Heptanal	9.763	1.5	ug/L	104577	2	8.805	356397	5.0
1,3-Propanediol...	10.458	0.7	ug/L	53033	3	11.194	356208	5.0
Octanal	11.101	1.3	ug/L	94641	3	11.194	356208	5.0
1-Hexanol, 2-et...	11.490	2.4	ug/L	173750	3	11.194	356208	5.0
Nonanal	12.291	1.5	ug/L	102973	3	11.194	356208	5.0