LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011524\
Data File : VV033651.D
Acq On : 15 Jan 2024 18:33
Operator : SY/MD
Sample : P1131-19
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AY1

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\SFAMVTR011224WMA.M
Title : TRACE VOA SFAM1.0

Signal : TIC: VV033651.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.082	7	13	28	rVB	63587	63615	2.83%	0.724%
2	1.185	38	45	54	rBV	237808	205088	9.12%	2.334%
3	1.275	61	73	80	rBV3	141218	220730	9.82%	2.512%
4	1.320	83	87	97	rVV2	20825	25307	1.13%	0.288%
5	1.529	146	152	163	rVB	37069	46976	2.09%	0.535%
6	1.590	164	171	187	rVB2	64186	92509	4.12%	1.053%
7	1.757	218	223	237	rVB3	19387	28934	1.29%	0.329%
8	1.970	281	289	298	rBV4	17829	25927	1.15%	0.295%
9	2.053	307	315	337	rVB	71892	117657	5.23%	1.339%
10	2.462	437	442	449	rVV3	21714	36965	1.64%	0.421%
11	2.510	449	457	473	rVB5	30084	72301	3.22%	0.823%
12	2.683	501	511	527	rVB2	44552	83594	3.72%	0.952%
13	3.105	629	642	658	rBV	128632	262054	11.66%	2.983%
14	3.654	799	813	826	rBV2	19503	46859	2.08%	0.533%
15	3.806	845	860	914	rBV	929012	2248032	100.00%	25.588%
16	4.246	983	997	1029	rVB2	57116	147115	6.54%	1.675%
17	4.574	1084	1099	1111	rBV4	15717	39696	1.77%	0.452%
18	4.670	1111	1129	1147	rVB3	13078	38684	1.72%	0.440%
19	4.953	1203	1217	1225	rBV3	134697	305964	13.61%	3.483%
20	5.005	1226	1233	1263	rVV2	142414	384831	17.12%	4.380%
21	5.127	1264	1271	1290	rVV8	18041	51954	2.31%	0.591%
22	5.227	1290	1302	1319	rVB5	11063	25522	1.14%	0.291%
23	5.532	1386	1397	1429	rBV	132622	300440	13.36%	3.420%
24	5.841	1484	1493	1497	rBV6	13952	24026	1.07%	0.273%
25	5.989	1527	1539	1567	rBV2	74969	187888	8.36%	2.139%
26	6.696	1741	1759	1771	rBV6	9691	25514	1.13%	0.290%
27	6.918	1818	1828	1848	rVB3	54189	106252	4.73%	1.209%
28	7.088	1858	1881	1896	rBV5	23952	59390	2.64%	0.676%
29	7.243	1904	1929	1947	rBV	156040	299814	13.34%	3.413%
30	7.561	2008	2028	2044	rBV6	23803	70193	3.12%	0.799%
31	8.037	2158	2176	2203	rBV	133399	358664	15.95%	4.083%
32	8.786	2399	2409	2413	rBV	210911	334579	14.88%	3.808%
33	8.812	2414	2417	2440	rVB	232583	349764	15.56%	3.981%
34	9.194	2526	2536	2545	rBV5	20217	36475	1.62%	0.415%
35	9.403	2586	2601	2605	rBV10	10826	25136	1.12%	0.286%
36	9.641	2666	2675	2687	rBV7	11952	22810	1.01%	0.260%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011524\
Data File : VV033651.D
Acq On : 15 Jan 2024 18:33
Operator : SY/MD
Sample : P1131-19
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AY1

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\SFAMVTR011224WMA.M
Title : TRACE VOA SFAM1.0

37	9.735	2693	2704	2715	rBV7	13149	30016	1.34%	0.342%
38	10.156	2824	2835	2849	rVB	58074	104289	4.64%	1.187%
39	10.288	2863	2876	2888	rBV	9895	29048	1.29%	0.331%
40	10.799	3015	3035	3045	rBV2	19901	49752	2.21%	0.566%
41	10.870	3050	3057	3070	rVB4	15255	25420	1.13%	0.289%
42	10.992	3087	3095	3101	rBV	46909	68749	3.06%	0.783%
43	11.120	3127	3135	3143	rBV2	26940	46550	2.07%	0.530%
44	11.185	3146	3155	3180	rVB2	233257	538978	23.98%	6.135%
45	11.577	3261	3277	3291	rVB3	323145	707028	31.45%	8.048%
46	11.921	3376	3384	3393	rBV9	13791	26108	1.16%	0.297%
47	12.233	3474	3481	3490	rVB2	23352	38490	1.71%	0.438%
48	12.622	3594	3602	3612	rBV2	40572	58138	2.59%	0.662%
49	13.201	3773	3782	3790	rBV	42034	69491	3.09%	0.791%
50	13.246	3790	3796	3813	rVV2	31614	57032	2.54%	0.649%
51	13.545	3882	3889	3903	rVB10	11254	22481	1.00%	0.256%
52	13.628	3905	3915	3926	rBV7	19773	41809	1.86%	0.476%
53	14.236	4097	4104	4113	rVB3	16931	24382	1.08%	0.278%
54	14.384	4140	4150	4160	rVB4	13013	26899	1.20%	0.306%
55	14.615	4215	4222	4237	rVB4	10537	25936	1.15%	0.295%
56	15.750	4568	4575	4582	rBV3	15501	23512	1.05%	0.268%

Sum of corrected areas: 8785367

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011524\
Data File : VV033651.D
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Operator : SY/MD
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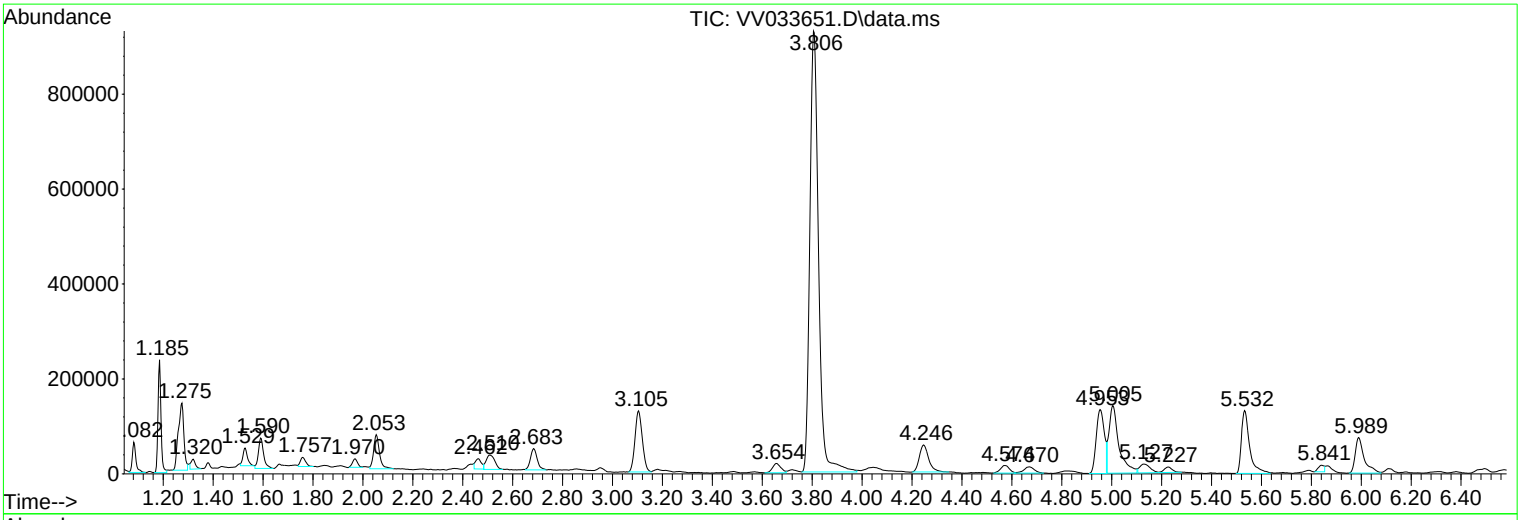
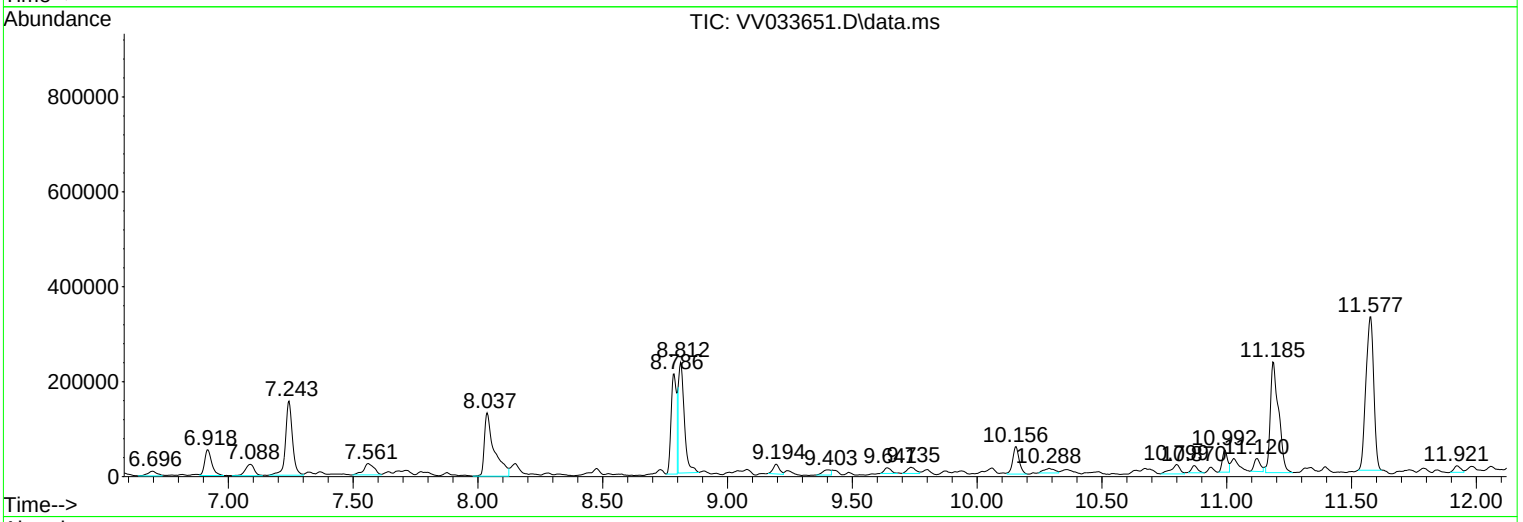
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 26 Sample Multiplier: 1

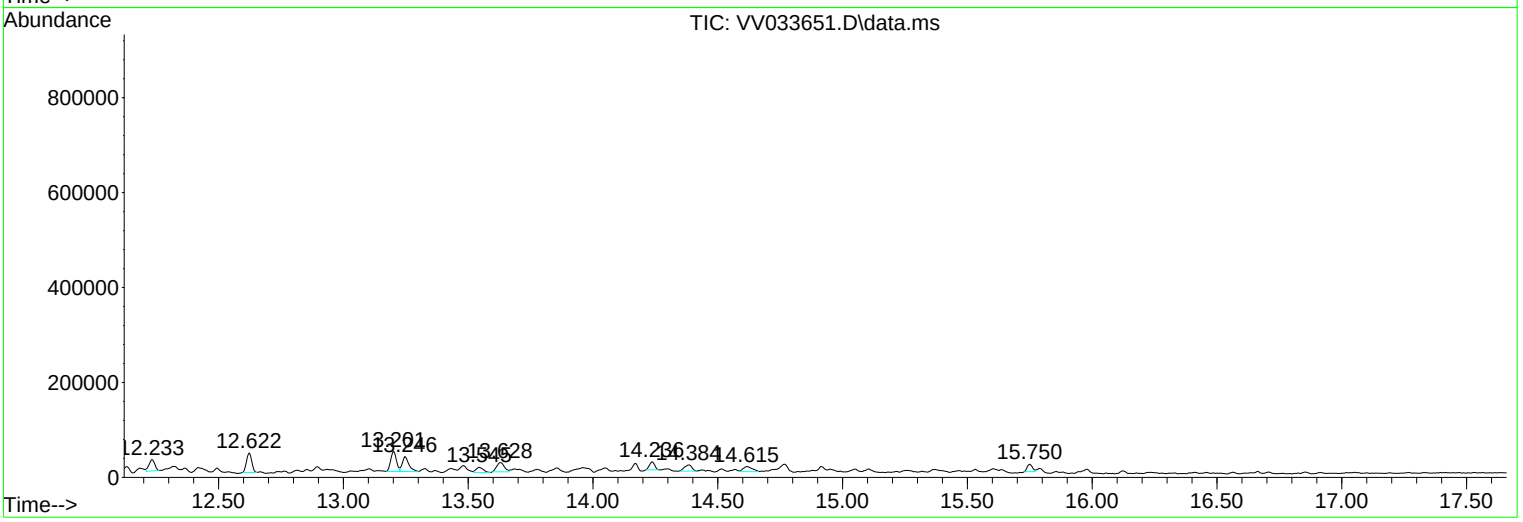
Instrument :
MSVOA_V
ClientSampleId :
C0AY1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR011224WMA.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\VV011524\
Data File : VV033651.D
Acq On : 15 Jan 2024 18:33
Operator : SY/MD
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Instrument :
MSVOA_V
ClientSampleId :
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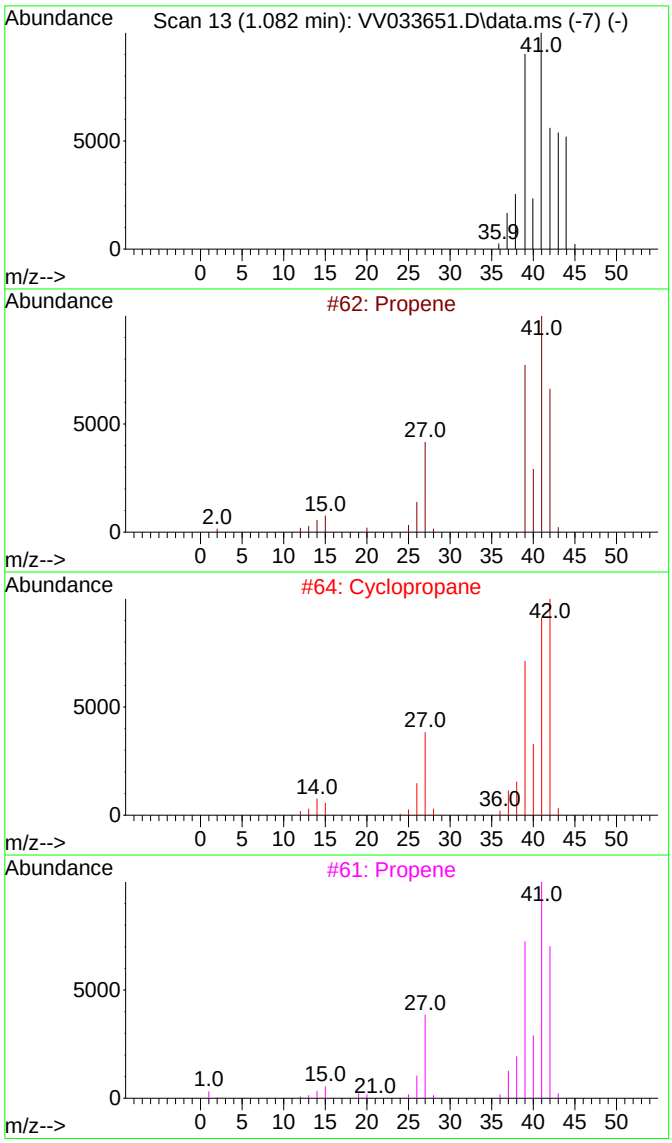
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR011224WMA.M
Quant Title : TRACE VOA SFAM1.0

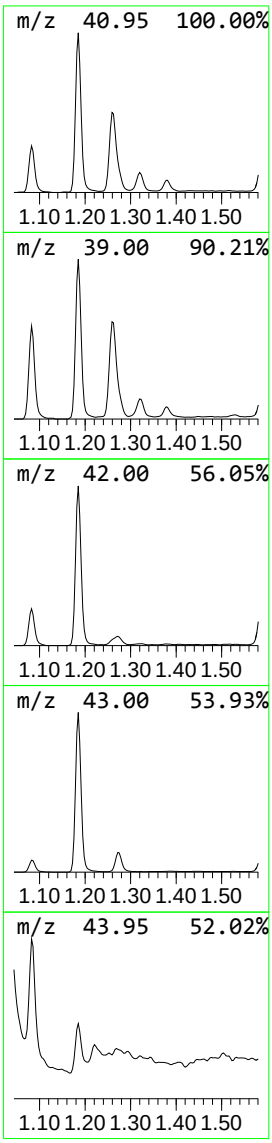
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.082	1.06 ug/L	63615	1,4-Difluorobenzene	5.532

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propene	42	C3H6	000115-07-1	42
2	Cyclopropane	42	C3H6	000075-19-4	7
3	Propene	42	C3H6	000115-07-1	7
4	Cyclopropane	42	C3H6	000075-19-4	4
5	Cyclopropene	40	C3H4	002781-85-3	4





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011524\
Data File : VV033651.D
Acq On : 15 Jan 2024 18:33
Operator : SY/MD
Sample : P1131-19
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AY1

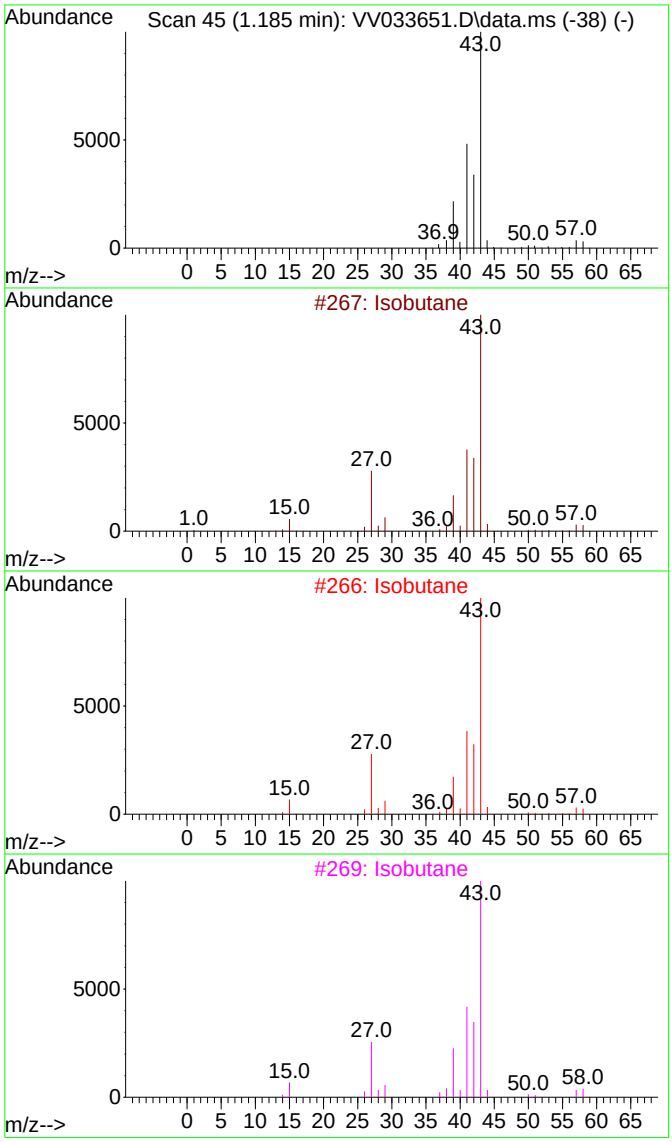
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR011224WMA.M
Quant Title : TRACE VOA SFAM1.0

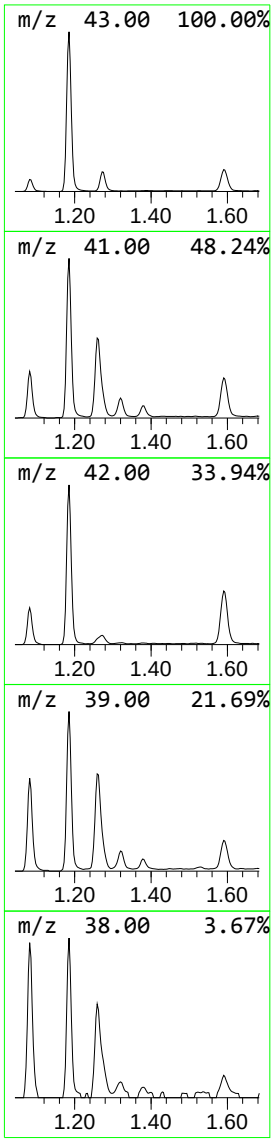
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.185	3.41 ug/L	205088	1,4-Difluorobenzene	5.532

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isobutane	58	C4H10	000075-28-5	80
2	Isobutane	58	C4H10	000075-28-5	72
3	Isobutane	58	C4H10	000075-28-5	64
4	Isobutane	58	C4H10	000075-28-5	64
5	Isobutane	58	C4H10	000075-28-5	59





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011524\
Data File : VV033651.D
Acq On : 15 Jan 2024 18:33
Operator : SY/MD
Sample : P1131-19
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AY1

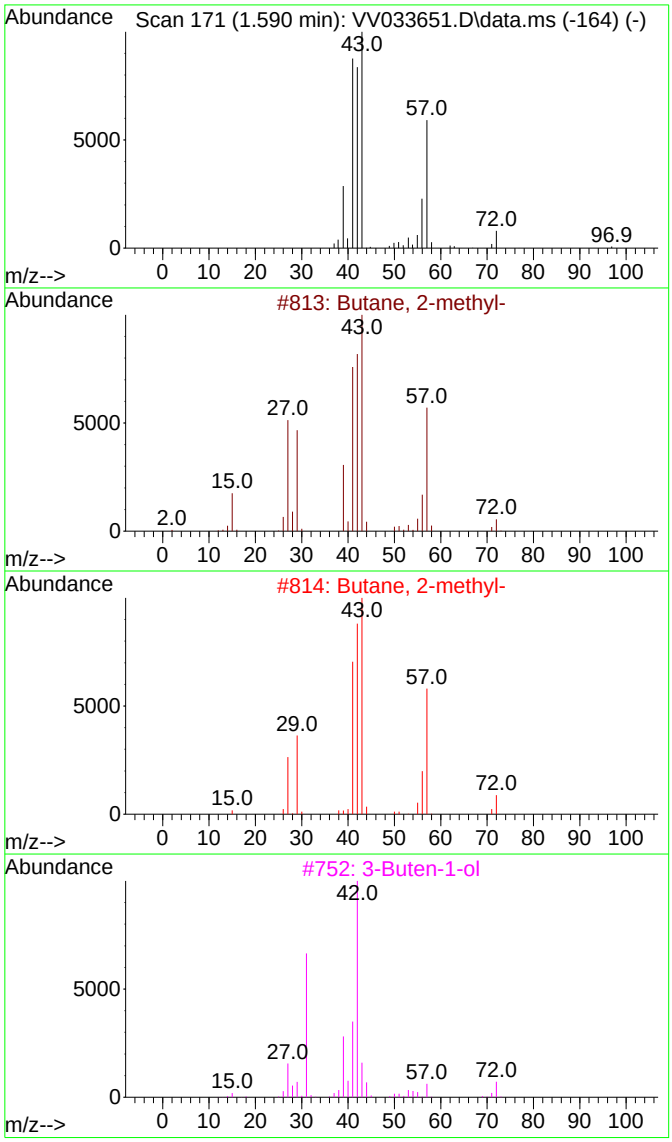
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR011224WMA.M
Quant Title : TRACE VOA SFAM1.0

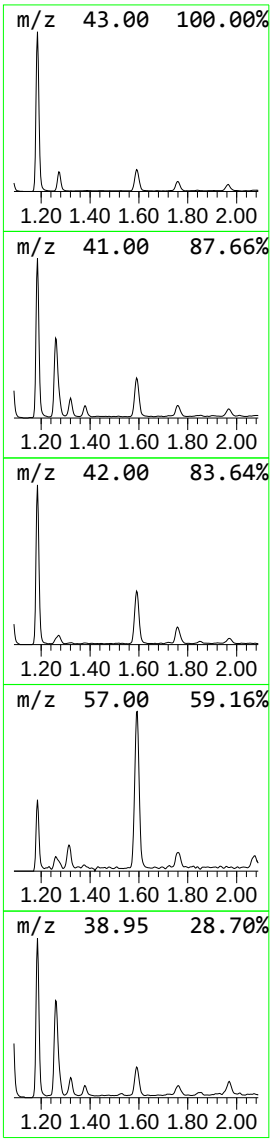
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.590	1.54 ug/L	92509	1,4-Difluorobenzene	5.532

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	90
2	Butane, 2-methyl-	72	C5H12	000078-78-4	80
3	3-Buten-1-ol	72	C4H8O	000627-27-0	47
4	1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	39
5	2(3H)-Furanone, dihydro-4-methyl-	100	C5H8O2	001679-49-8	33





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011524\
Data File : VV033651.D
Acq On : 15 Jan 2024 18:33
Operator : SY/MD
Sample : P1131-19
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AY1

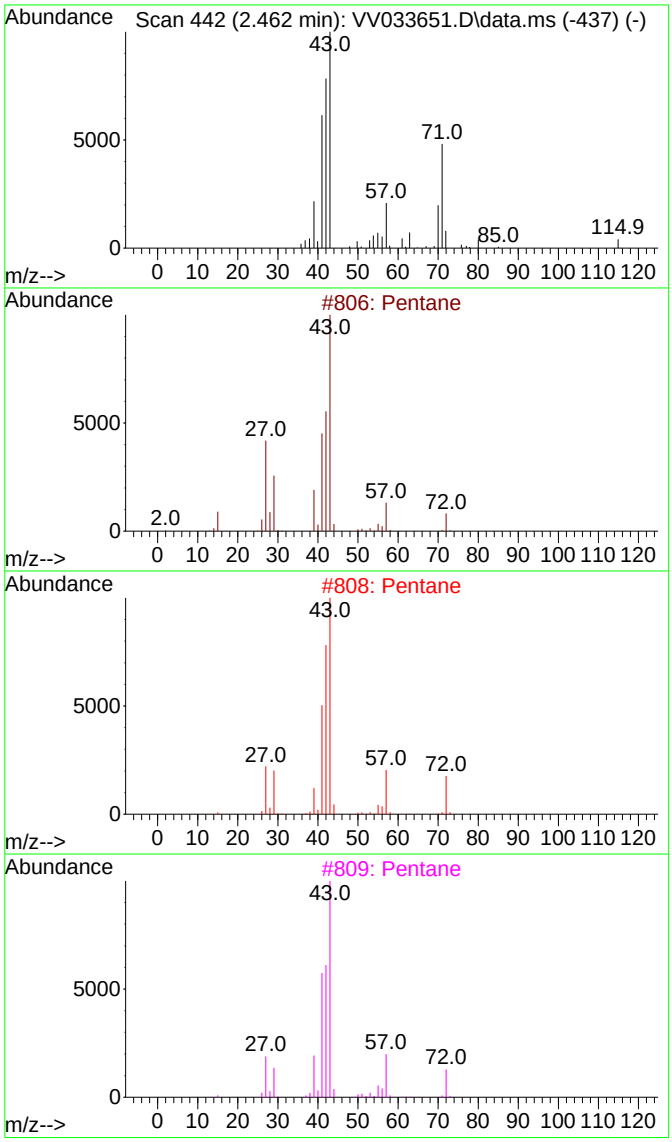
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR011224WMA.M
Quant Title : TRACE VOA SFAM1.0

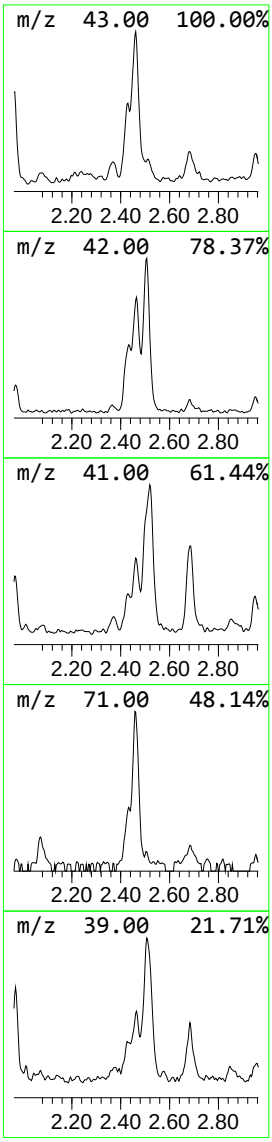
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.462	0.62 ug/L	36965	1,4-Difluorobenzene	5.532

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane	72	C5H12	000109-66-0	43
2	Pentane	72	C5H12	000109-66-0	43
3	Pentane	72	C5H12	000109-66-0	43
4	1-Pentene	70	C5H10	000109-67-1	43
5	Pentane, 2-methyl-	86	C6H14	000107-83-5	39





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011524\
Data File : VV033651.D
Acq On : 15 Jan 2024 18:33
Operator : SY/MD
Sample : P1131-19
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AY1

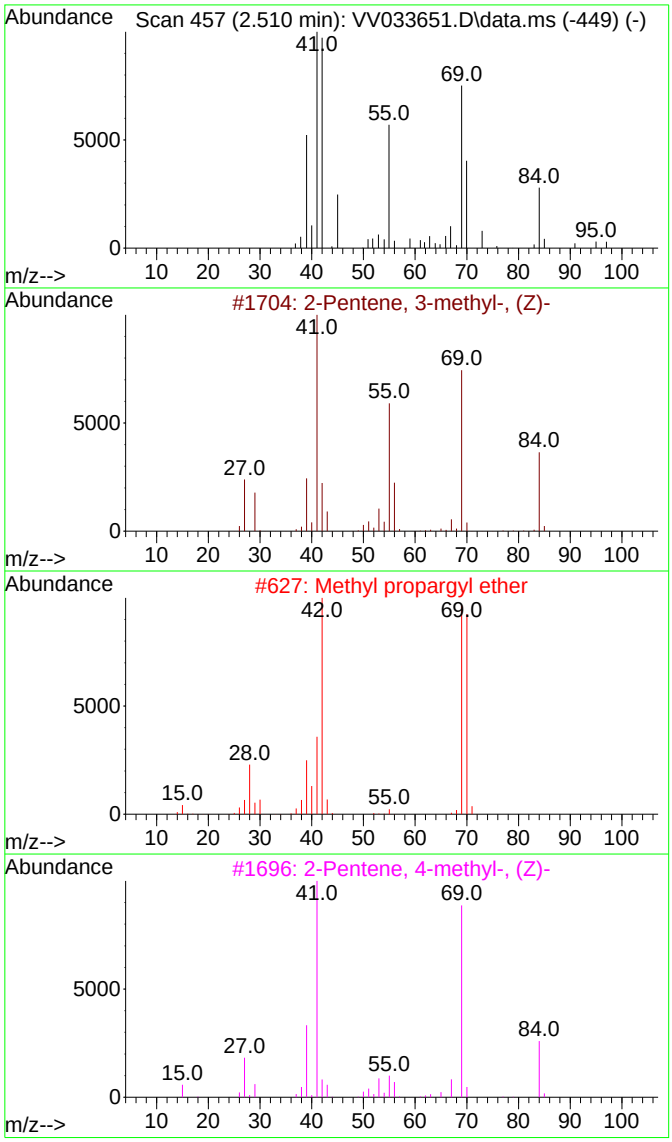
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR011224WMA.M
Quant Title : TRACE VOA SFAM1.0

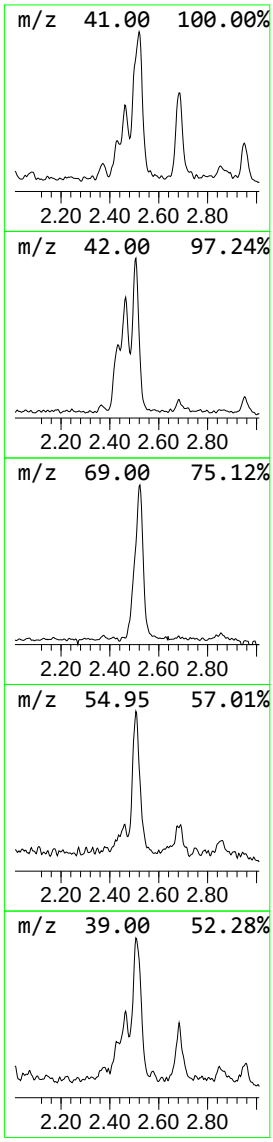
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.510	1.20 ug/L	72301	1,4-Difluorobenzene	5.532

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentene, 3-methyl-, (Z)-	84	C6H12	000922-62-3	35
2	Methyl propargyl ether	70	C4H6O	000627-41-8	35
3	2-Pentene, 4-methyl-, (Z)-	84	C6H12	000691-38-3	30
4	2-Butene, 2,3-dimethyl-	84	C6H12	000563-79-1	30
5	2-Butene, 2,3-dimethyl-	84	C6H12	000563-79-1	30





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Instrument :
MSVOA_V
ClientSampleId :
C0AY1

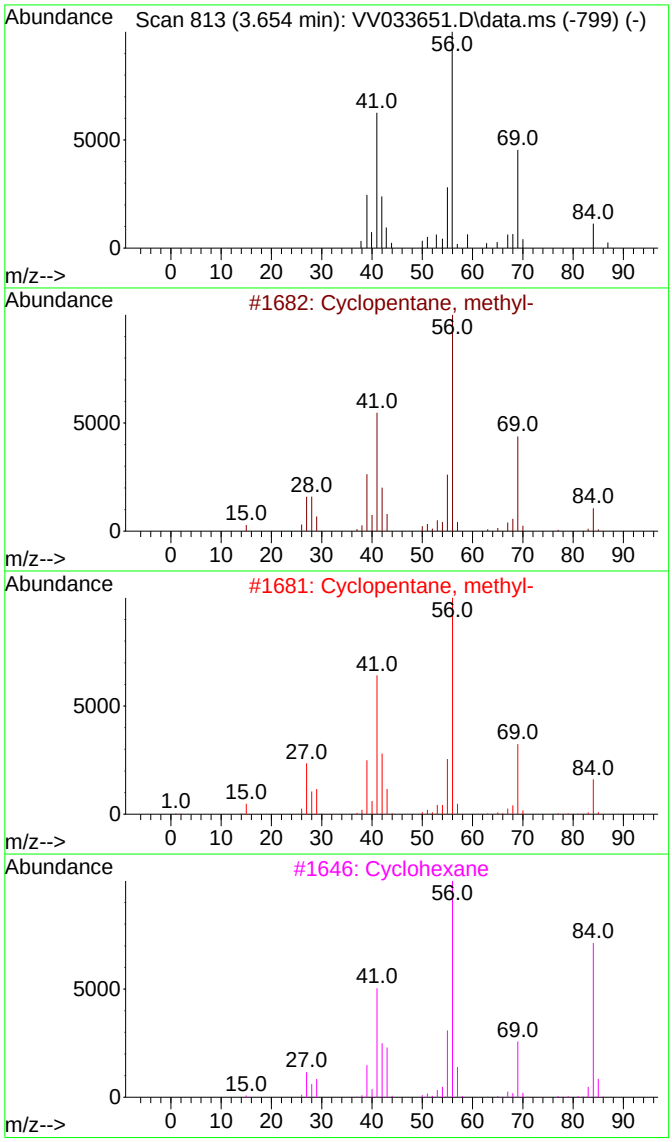
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR011224WMA.M
Quant Title : TRACE VOA SFAM1.0

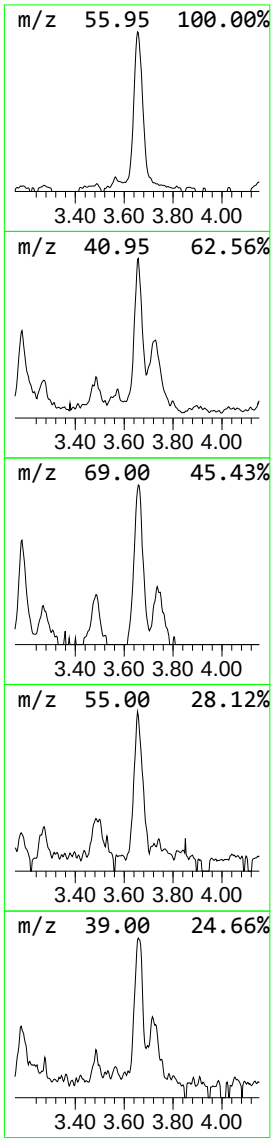
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 (DEL) Alkane: Cyclic3.654 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.654	0.78 ug/L	46859	1,4-Difluorobenzene	5.532

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2	Cyclopentane, methyl-	84	C6H12	000096-37-7	80
3	Cyclohexane	84	C6H12	000110-82-7	78
4	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	72
5	Cyclopentane, methyl-	84	C6H12	000096-37-7	72





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011524\
Data File : VV033651.D
Acq On : 15 Jan 2024 18:33
Operator : SY/MD
Sample : P1131-19
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AY1

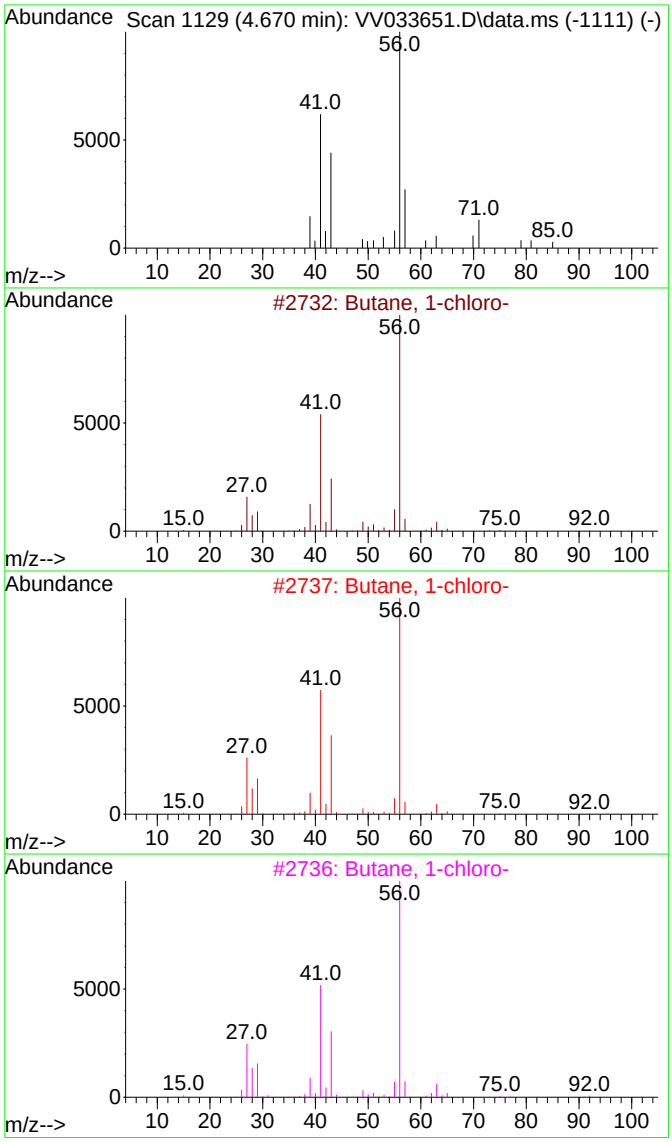
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR011224WMA.M
Quant Title : TRACE VOA SFAM1.0

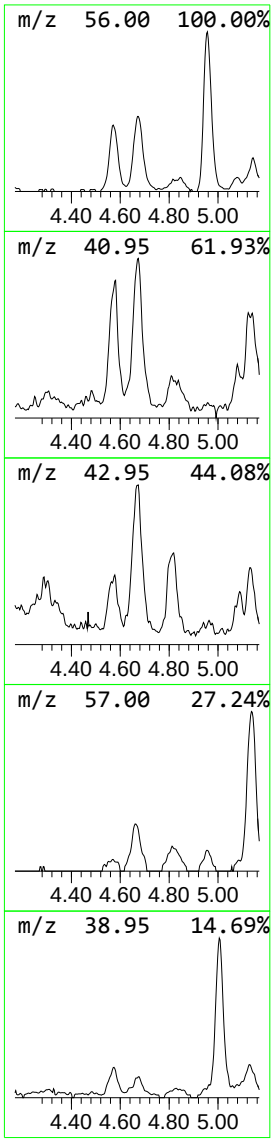
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Butane, 1-chloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.670	0.64 ug/L	38684	1,4-Difluorobenzene	5.532

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 1-chloro-	92	C4H9Cl	000109-69-3	53
2	Butane, 1-chloro-	92	C4H9Cl	000109-69-3	45
3	Butane, 1-chloro-	92	C4H9Cl	000109-69-3	45
4	4-Penten-2-ol, 4-methyl-	100	C6H12O	002004-67-3	40
5	Butane, 1-chloro-	92	C4H9Cl	000109-69-3	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011524\
Data File : VV033651.D
Acq On : 15 Jan 2024 18:33
Operator : SY/MD
Sample : P1131-19
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AY1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR011224WMA.M
Quant Title : TRACE VOA SFAM1.0

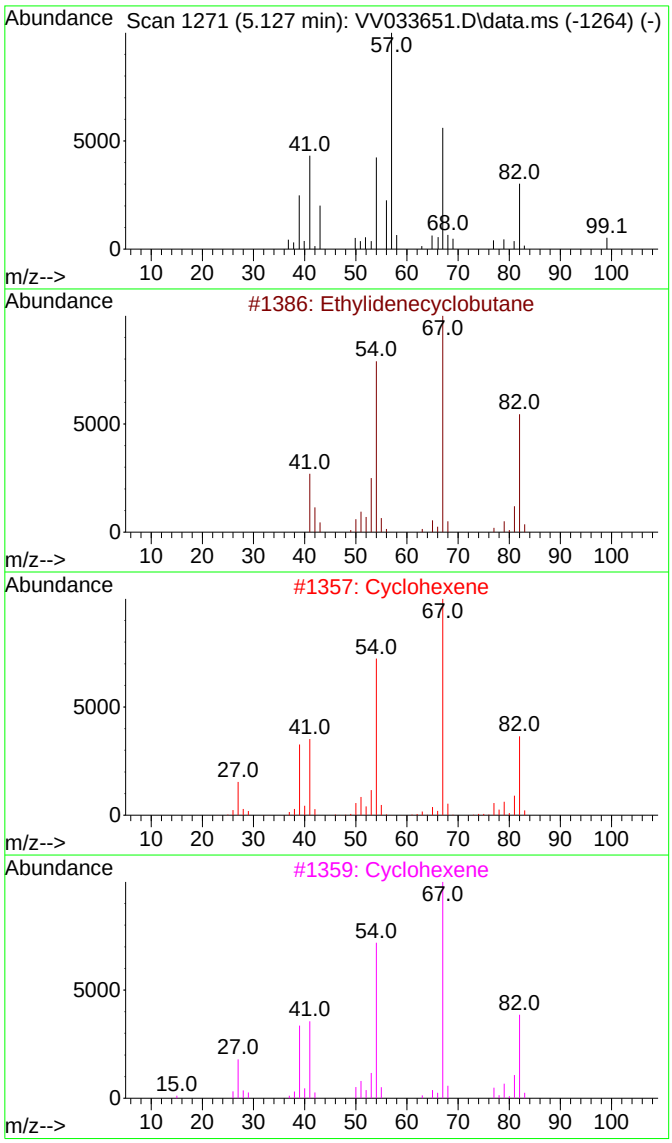
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

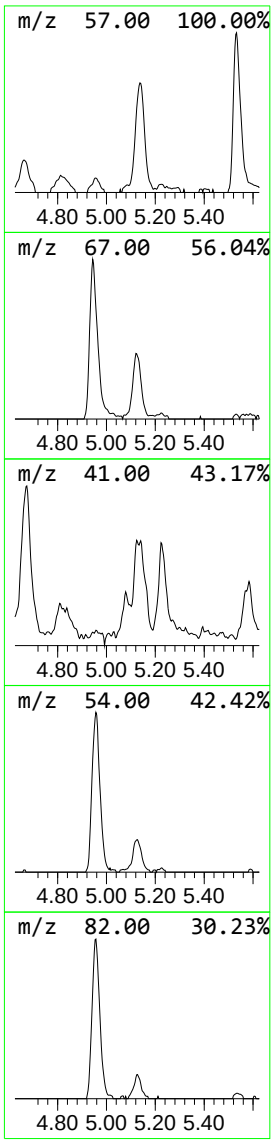
****

Peak Number 8 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.127	0.86 ug/L	51954	1,4-Difluorobenzene	5.532

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethylidenecyclobutane	82	C6H10	001528-21-8	40
2	Cyclohexene	82	C6H10	000110-83-8	38
3	Cyclohexene	82	C6H10	000110-83-8	38
4	Cyclohexene	82	C6H10	000110-83-8	38
5	Cyclohexene	82	C6H10	000110-83-8	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011524\
Data File : VV033651.D
Acq On : 15 Jan 2024 18:33
Operator : SY/MD
Sample : P1131-19
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AY1

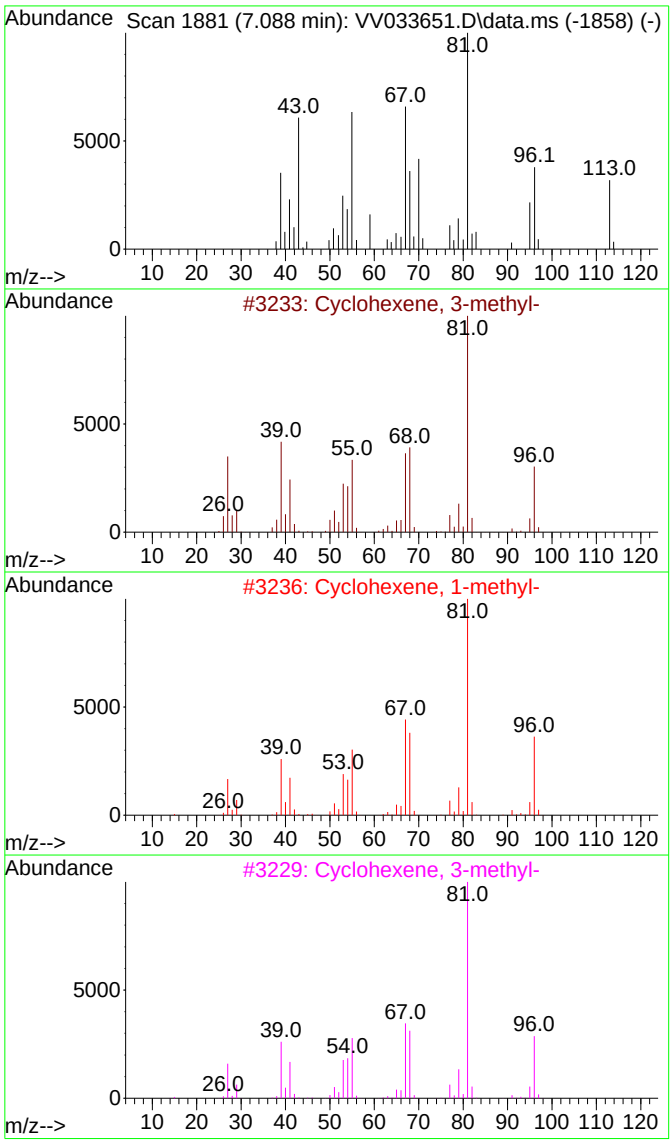
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR011224WMA.M
Quant Title : TRACE VOA SFAM1.0

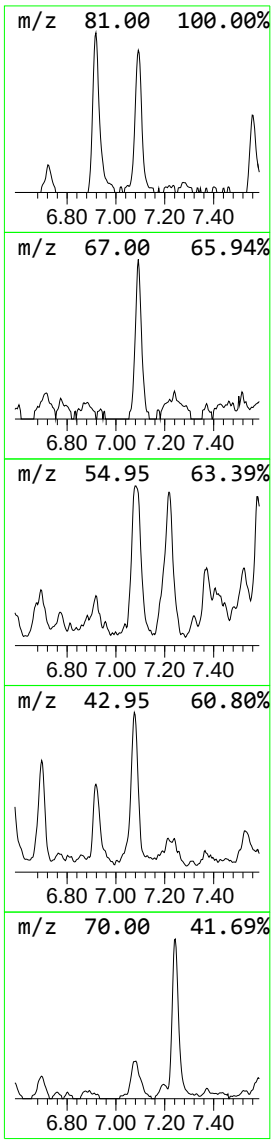
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Cyclohexene, 3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.088	0.99 ug/L	59390	1,4-Difluorobenzene	5.532

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexene, 3-methyl-	96	C7H12	000591-48-0	90
2	Cyclohexene, 1-methyl-	96	C7H12	000591-49-1	64
3	Cyclohexene, 3-methyl-	96	C7H12	000591-48-0	60
4	Cyclohexene, 1-methyl-	96	C7H12	000591-49-1	60
5	Cyclohexane, methylene-	96	C7H12	001192-37-6	58





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011524\
Data File : VV033651.D
Acq On : 15 Jan 2024 18:33
Operator : SY/MD
Sample : P1131-19
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AY1

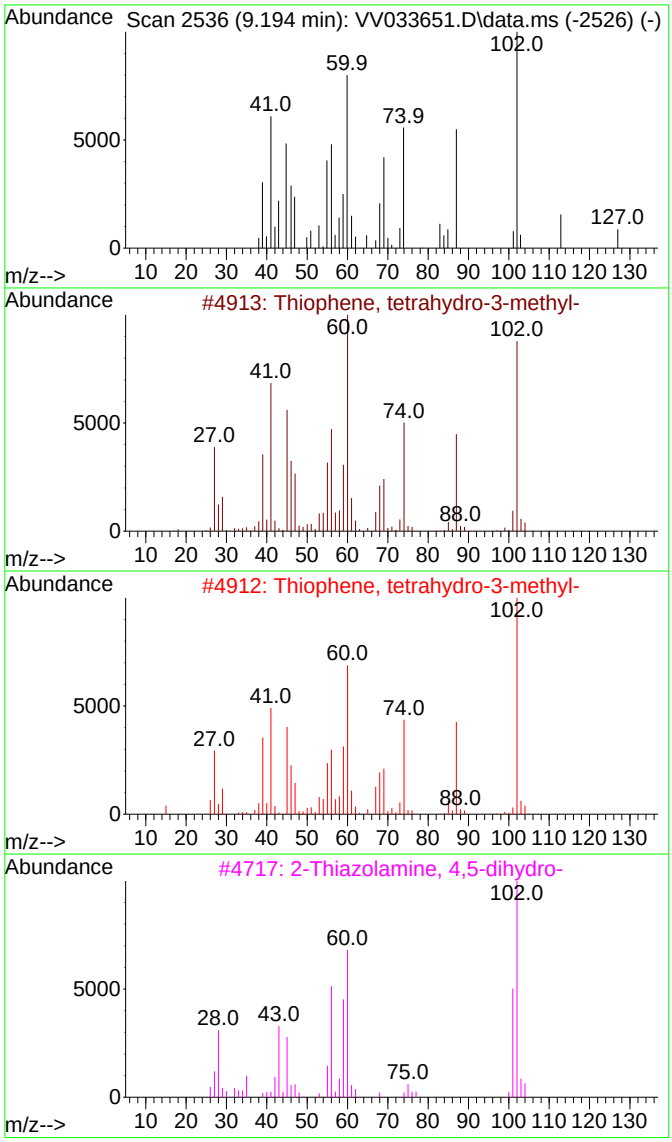
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR011224WMA.M
Quant Title : TRACE VOA SFAM1.0

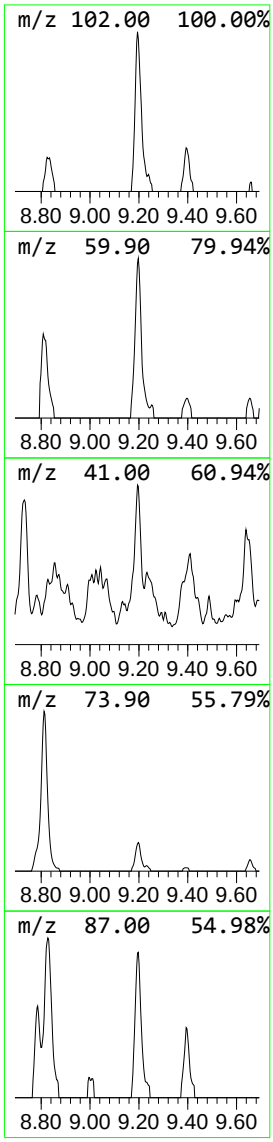
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Thiophene, tetrahydro-3-met... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.194	0.55 ug/L	36475	Chlorobenzene-d5	8.783

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Thiophene, tetrahydro-3-methyl-	102	C5H10S	004740-00-5	96
2	Thiophene, tetrahydro-3-methyl-	102	C5H10S	004740-00-5	87
3	2-Thiazolamine, 4,5-dihydro-	102	C3H6N2S	001779-81-3	43
4	Cyclopentanethiol	102	C5H10S	001679-07-8	38
5	2-Thiazolamine, 4,5-dihydro-	102	C3H6N2S	001779-81-3	37





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011524\
Data File : VV033651.D
Acq On : 15 Jan 2024 18:33
Operator : SY/MD
Sample : P1131-19
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AY1

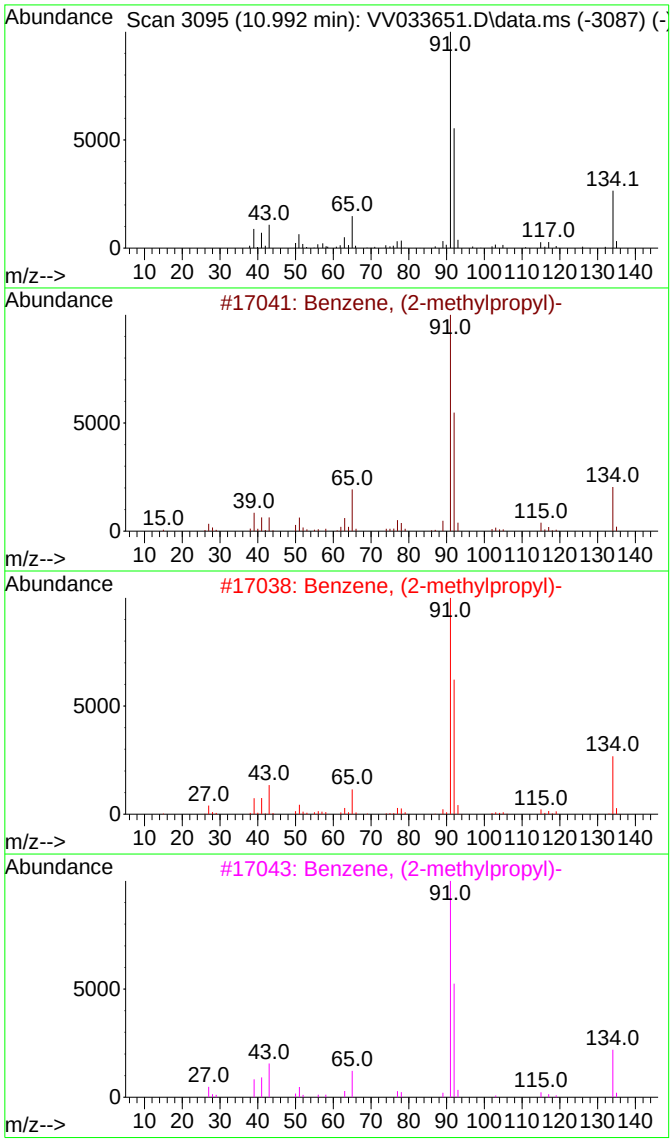
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR011224WMA.M
Quant Title : TRACE VOA SFAM1.0

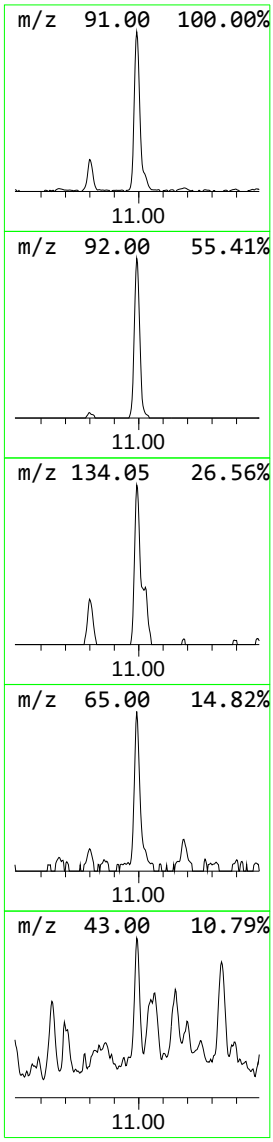
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Benzene, (2-methylpropyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.992	0.64 ug/L	68749	1,4-Dichlorobenzene-d4	11.185

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	91
2	Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	91
3	Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	91
4	Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90
5	Benzene, n-butyl-	134	C10H14	000104-51-8	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011524\
Data File : VV033651.D
Acq On : 15 Jan 2024 18:33
Operator : SY/MD
Sample : P1131-19
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AY1

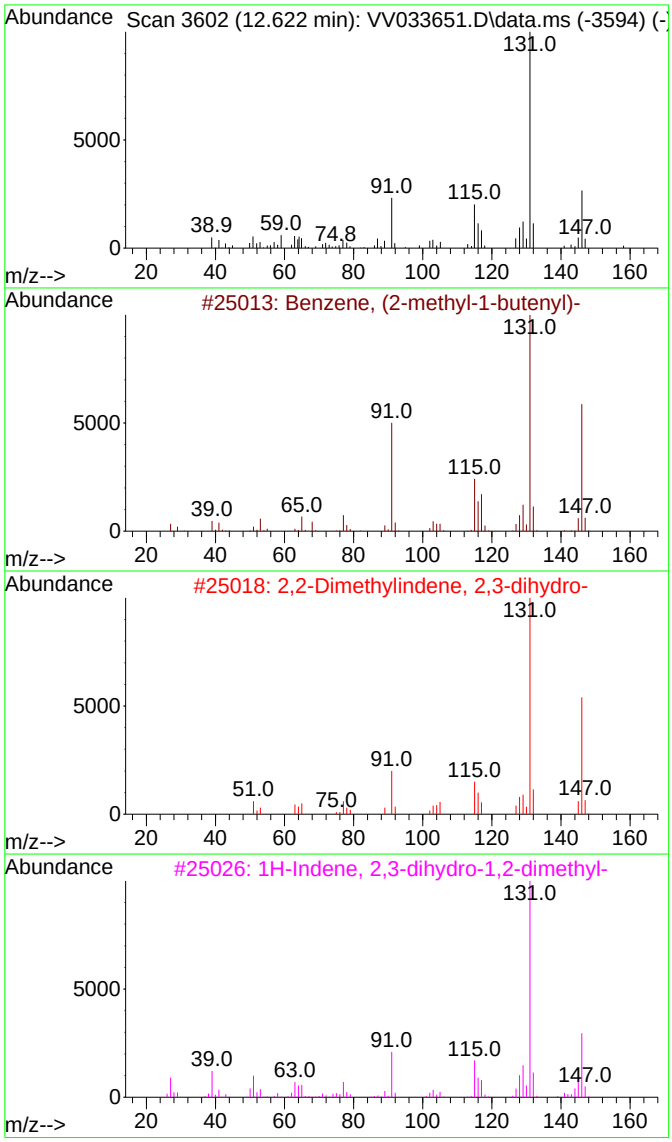
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR011224WMA.M
Quant Title : TRACE VOA SFAM1.0

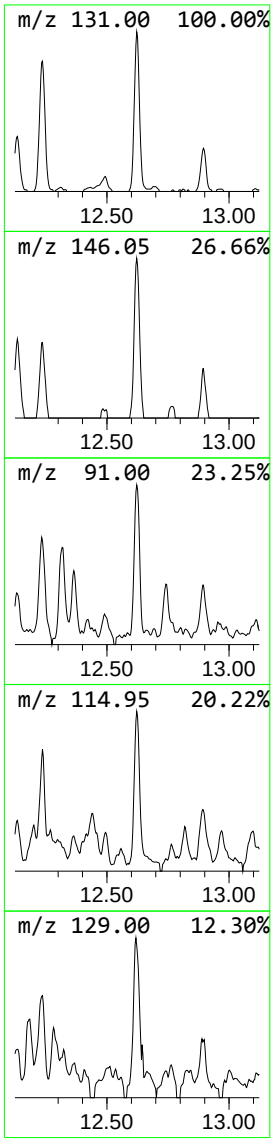
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Benzene, (2-methyl-1-butenyl)- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.622	0.54 ug/L	58138	1,4-Dichlorobenzene-d4	11.185

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	94
2	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	94
3	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	94
4	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	91
5	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011524\
Data File : VV033651.D
Acq On : 15 Jan 2024 18:33
Operator : SY/MD
Sample : P1131-19
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AY1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR011224WMA.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...	1.082	1.1	ug/L	63615	1	5.532	300440	5.0
(DEL) Alkane: S...	1.185	3.4	ug/L	205088	1	5.532	300440	5.0
(DEL) Alkane: S...	1.590	1.5	ug/L	92509	1	5.532	300440	5.0
(DEL) Alkane: S...	2.462	0.6	ug/L	36965	1	5.532	300440	5.0
(DEL) Alkane: S...	2.510	1.2	ug/L	72301	1	5.532	300440	5.0
(DEL) Alkane: C...	3.654	0.8	ug/L	46859	1	5.532	300440	5.0
Butane, 1-chloro-	4.670	0.6	ug/L	38684	1	5.532	300440	5.0
unknown-01	5.127	0.9	ug/L	51954	1	5.532	300440	5.0
Cyclohexene, 3-...	7.088	1.0	ug/L	59390	1	5.532	300440	5.0
Thiophene, tetr...	9.194	0.6	ug/L	36475	2	8.783	334579	5.0
Benzene, (2-met...	10.992	0.6	ug/L	68749	3	11.185	538978	5.0
Benzene, (2-met...	12.622	0.5	ug/L	58138	3	11.185	538978	5.0