LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070824\
Data File : VU059876.D
Acq On : 08 Jul 2024 12:20
Operator : MD/SY
Sample : P3137-17
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZG6

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\SFAMUTR061724WMA.M
Title : TRACE VOA SFAM1.0

Signal : TIC: VU059876.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.345	625	639	666	rVB	267525	537518	1.54%	1.069%
2	4.657	1017	1047	1110	rBV	15766124	34852041	100.00%	69.327%
3	5.718	1354	1377	1388	rBV3	172636	451840	1.30%	0.899%
4	7.570	1938	1953	1982	rBV3	351555	680831	1.95%	1.354%
5	7.891	2043	2053	2068	rVB	215660	370072	1.06%	0.736%
6	8.306	2163	2182	2198	rBV	826881	1402314	4.02%	2.789%
7	8.628	2266	2282	2306	rBV	250557	521426	1.50%	1.037%
8	9.187	2445	2456	2473	rBV	321154	522303	1.50%	1.039%
9	9.412	2515	2526	2537	rBV	278684	472600	1.36%	0.940%
10	9.875	2660	2670	2684	rVB2	251020	407919	1.17%	0.811%
11	10.496	2845	2863	2882	rBV	1355714	2277121	6.53%	4.530%
12	11.245	3081	3096	3112	rBV6	127244	357794	1.03%	0.712%
13	11.492	3160	3173	3184	rBV	310540	520754	1.49%	1.036%
14	11.650	3209	3222	3237	rVV	370530	707753	2.03%	1.408%
15	11.804	3255	3270	3285	rVB3	319521	649723	1.86%	1.292%
16	12.013	3325	3335	3346	rVB3	505664	1063117	3.05%	2.115%
17	12.148	3366	3377	3385	rBV	642278	1080693	3.10%	2.150%
18	12.187	3385	3389	3412	rVB4	260139	649099	1.86%	1.291%
19	12.309	3412	3427	3433	rBV3	202898	399825	1.15%	0.795%
20	12.357	3433	3442	3453	rVB5	238222	545357	1.56%	1.085%
21	12.817	3575	3585	3602	rVB2	210093	356461	1.02%	0.709%
22	14.155	3992	4001	4021	rVB2	333768	588839	1.69%	1.171%
23	14.576	4122	4132	4140	rBV2	215244	385022	1.10%	0.766%
24	15.563	4430	4439	4450	rVB4	268821	471574	1.35%	0.938%

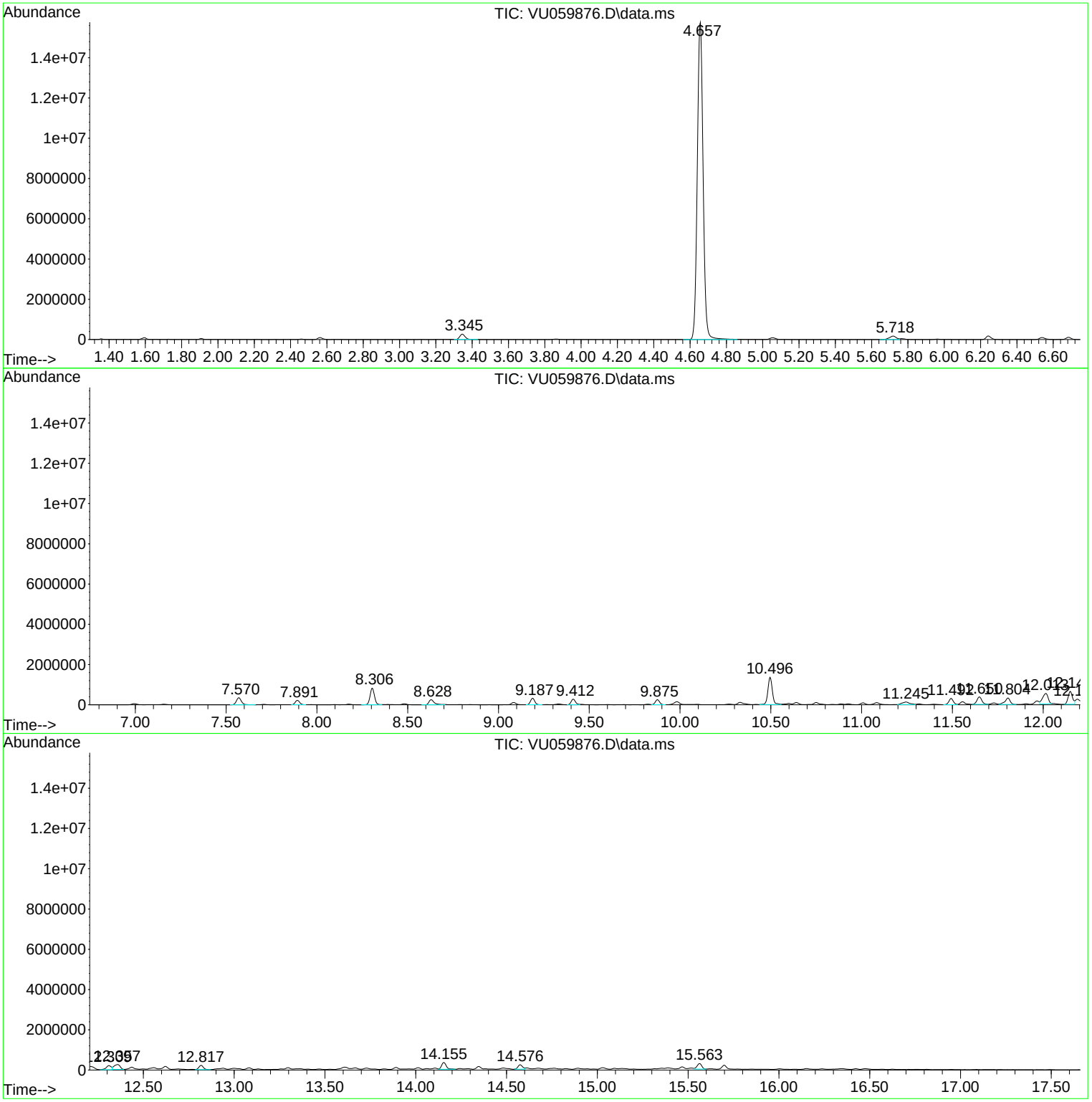
Sum of corrected areas: 50271996

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070824\
Data File : VU059876.D
Acq On : 08 Jul 2024 12:20
Operator : MD/SY
Sample : P3137-17
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZG6

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR061724WMA.M
Quant Title : TRACE VOA SFAM1.0

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070824\
Data File : VU059876.D
Acq On : 08 Jul 2024 12:20
Operator : MD/SY
Sample : P3137-17
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZG6

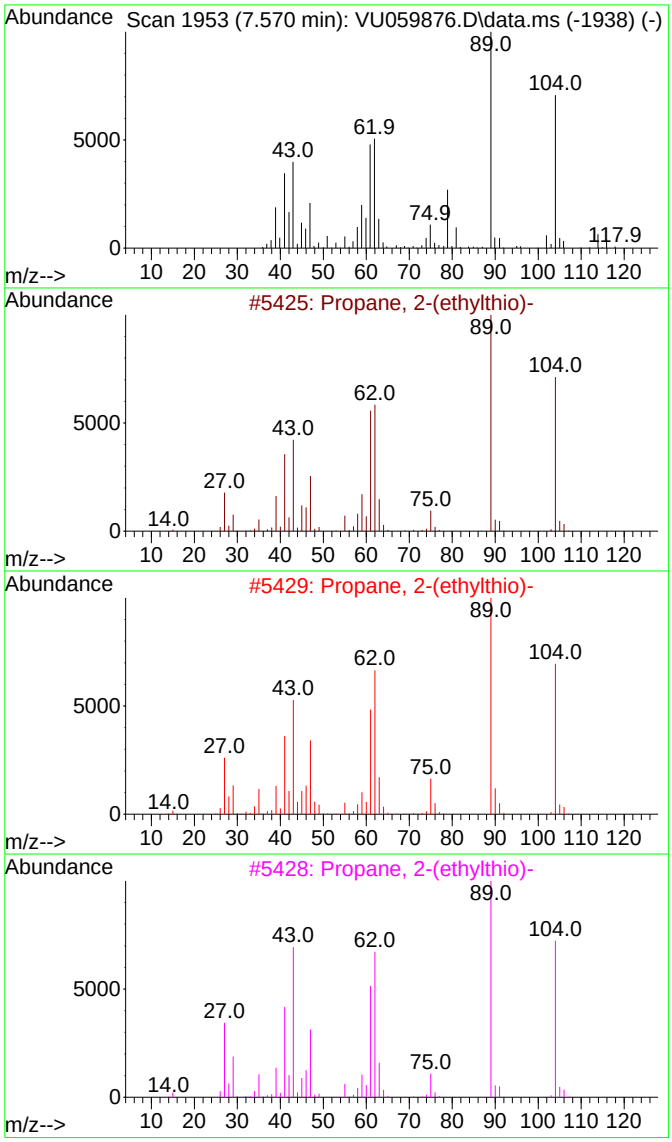
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR061724WMA.M
Quant Title : TRACE VOA SFAM1.0

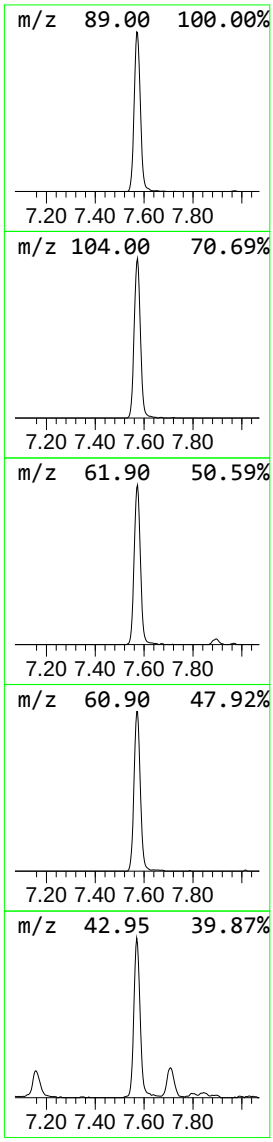
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Propane, 2-(ethylthio)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.570	7.53 ug/L	680831	1,4-Difluorobenzene	6.242

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propane, 2-(ethylthio)-	104	C5H12S	005145-99-3	96
2	Propane, 2-(ethylthio)-	104	C5H12S	005145-99-3	86
3	Propane, 2-(ethylthio)-	104	C5H12S	005145-99-3	83
4	Butane, 2-(ethylthio)-	118	C6H14S	005008-72-0	50
5	1-(1-(Methylthio)propyl)-2-propy...	196	C7H16S3	126876-22-0	47





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070824\
Data File : VU059876.D
Acq On : 08 Jul 2024 12:20
Operator : MD/SY
Sample : P3137-17
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZG6

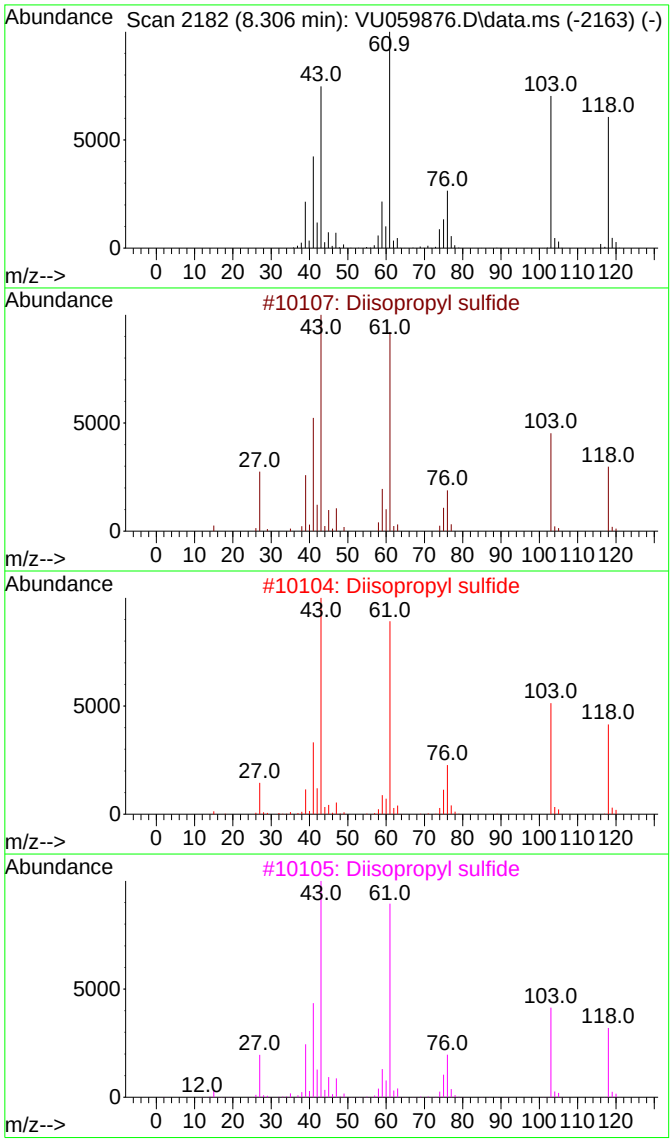
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR061724WMA.M
Quant Title : TRACE VOA SFAM1.0

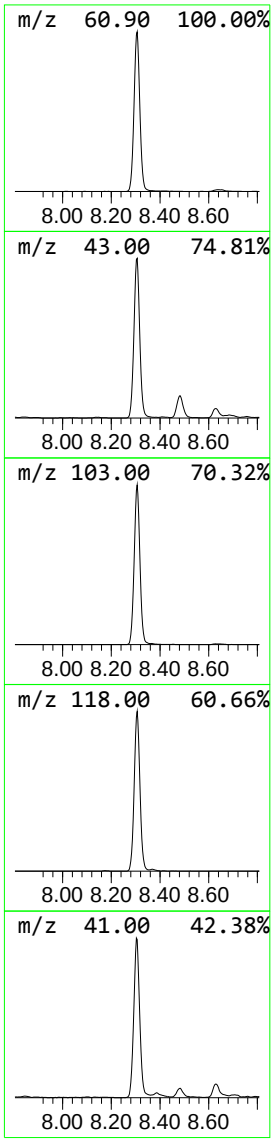
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Diisopropyl sulfide Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.306	14.84 ug/L	1402310	Chlorobenzene-d5	9.412

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diisopropyl sulfide	118	C6H14S	000625-80-9	94
2	Diisopropyl sulfide	118	C6H14S	000625-80-9	93
3	Diisopropyl sulfide	118	C6H14S	000625-80-9	90
4	Diisopropyl sulfide	118	C6H14S	000625-80-9	78
5	N-Methoxy-N-methylacetamide	103	C4H9NO2	078191-00-1	43





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070824\
Data File : VU059876.D
Acq On : 08 Jul 2024 12:20
Operator : MD/SY
Sample : P3137-17
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZG6

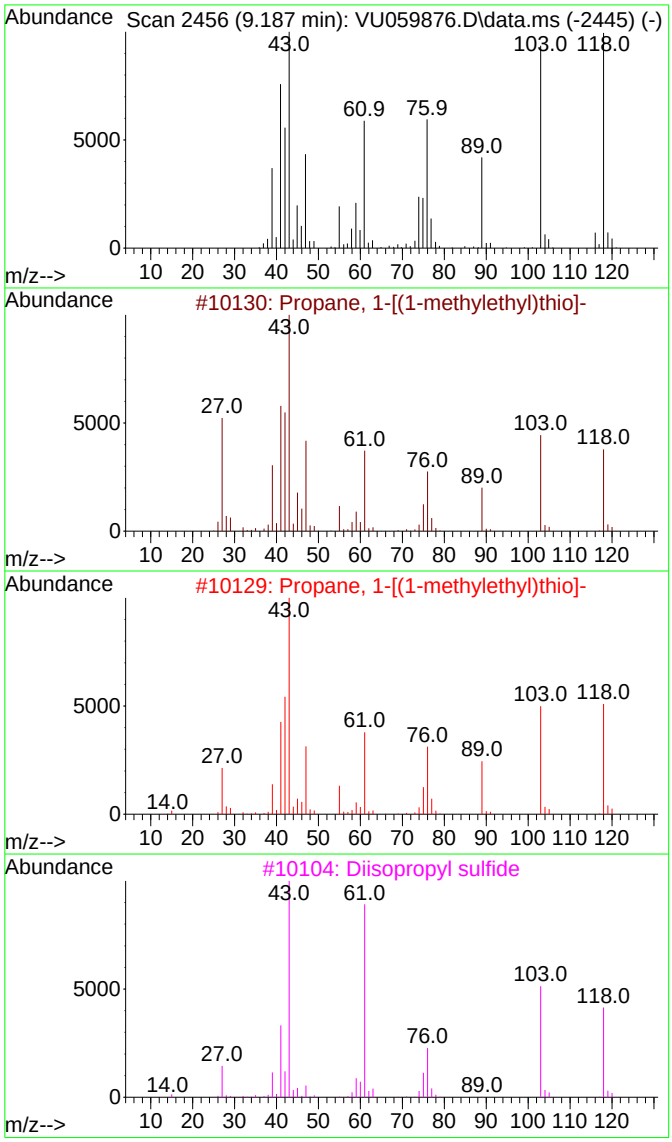
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR061724WMA.M
Quant Title : TRACE VOA SFAM1.0

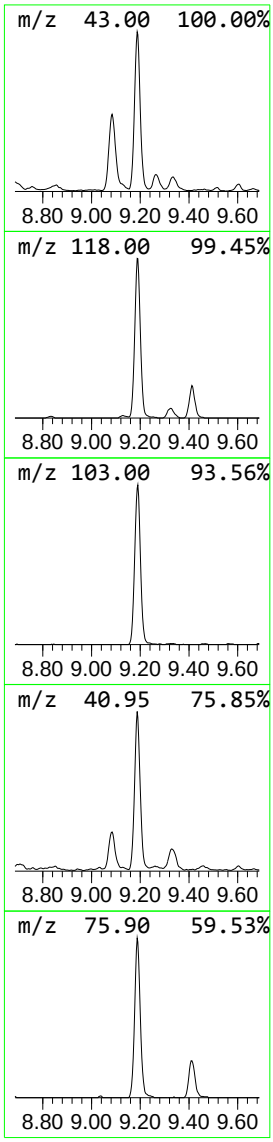
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Propane, 1-[(1-methylethyl)... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.187	5.53 ug/L	522303	Chlorobenzene-d5	9.412

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propane, 1-[(1-methylethyl)thio]-	118	C6H14S	005008-73-1	96
2	Propane, 1-[(1-methylethyl)thio]-	118	C6H14S	005008-73-1	96
3	Diisopropyl sulfide	118	C6H14S	000625-80-9	52
4	Ethanethioic acid, S-(1-methylet...	118	C5H10OS	000926-73-8	41
5	Acetamide, N-(aminothioxomethyl)-	118	C3H6N2OS	000591-08-2	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070824\
Data File : VU059876.D
Acq On : 08 Jul 2024 12:20
Operator : MD/SY
Sample : P3137-17
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZG6

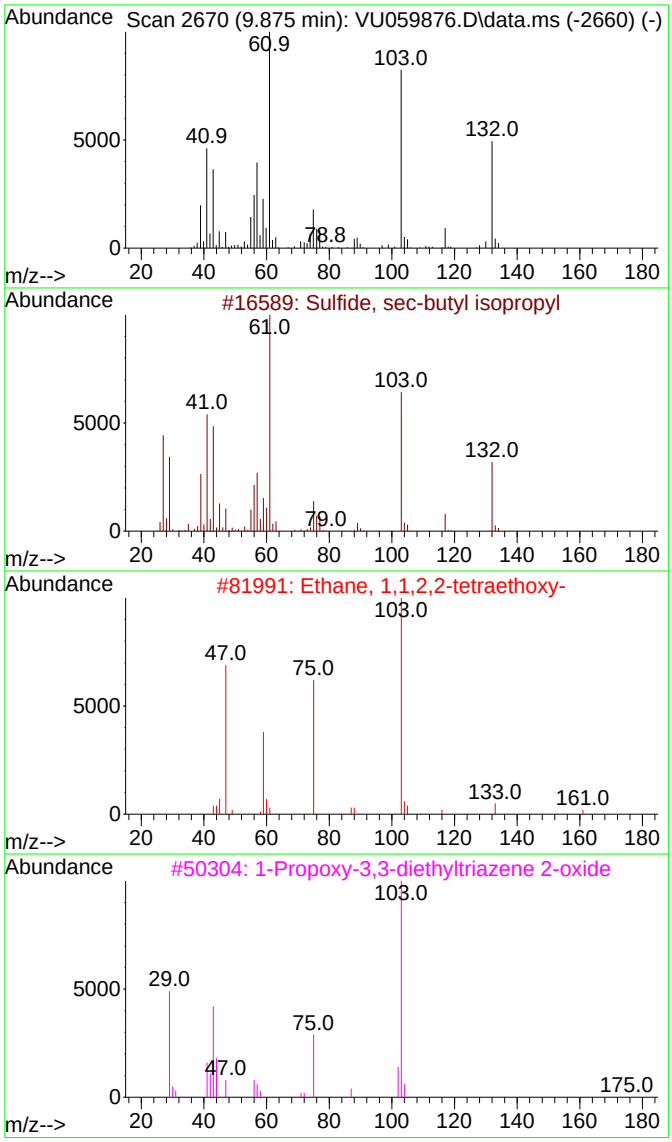
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR061724WMA.M
Quant Title : TRACE VOA SFAM1.0

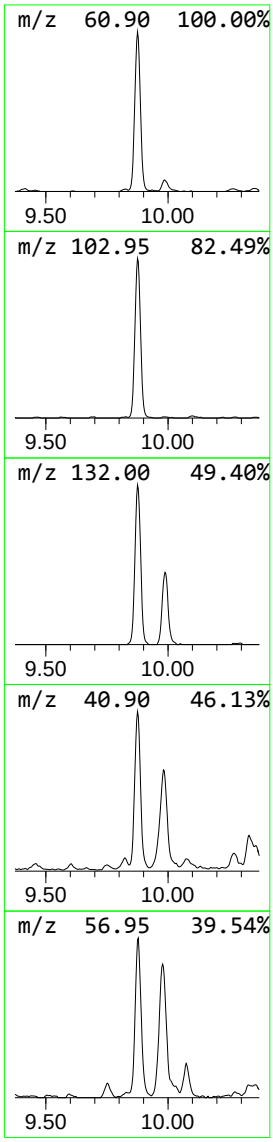
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Sulfide, sec-butyl isopropyl Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.875	4.32 ug/L	407919	Chlorobenzene-d5	9.412

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Sulfide, sec-butyl isopropyl	132	C7H16S	022438-36-4	87
2	Ethane, 1,1,2,2-tetraethoxy-	206	C10H22O4	003975-14-2	27
3	1-Propoxy-3,3-diethyltriazene 2-...	175	C7H17N3O2	112753-64-7	22
4	Butanedioic acid, hydroxy-, dime...	162	C6H10O5	001587-15-1	17
5	Ethanethioamide, N,N-dimethyl-	103	C4H9NS	000631-67-4	16





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070824\
Data File : VU059876.D
Acq On : 08 Jul 2024 12:20
Operator : MD/SY
Sample : P3137-17
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZG6

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR061724WMA.M
Quant Title : TRACE VOA SFAM1.0

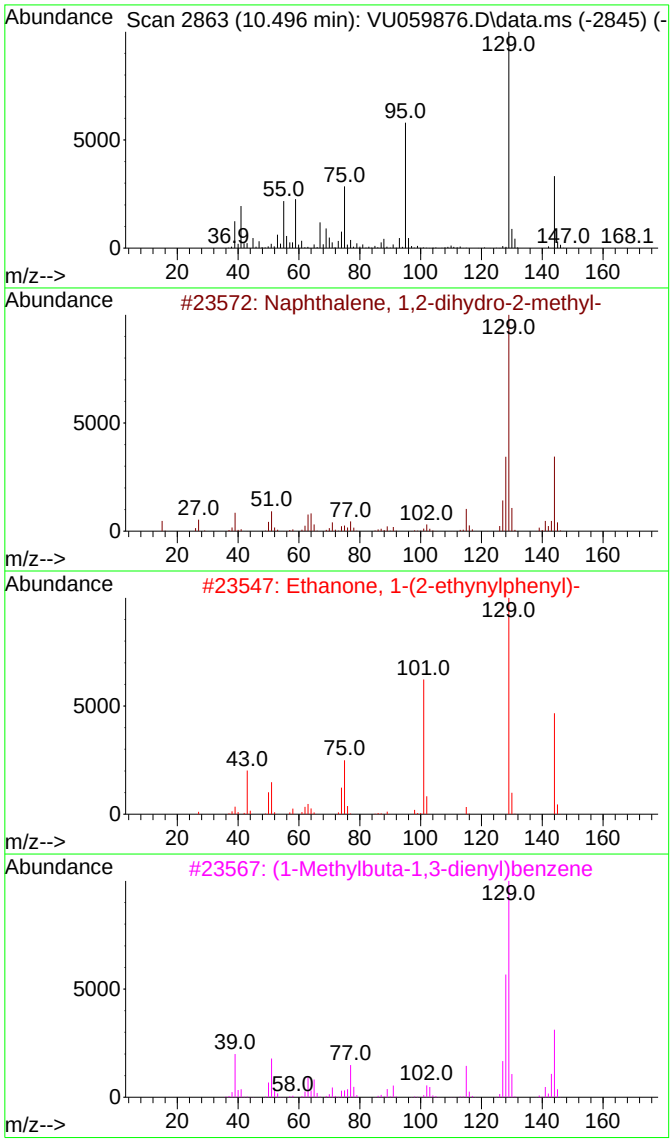
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

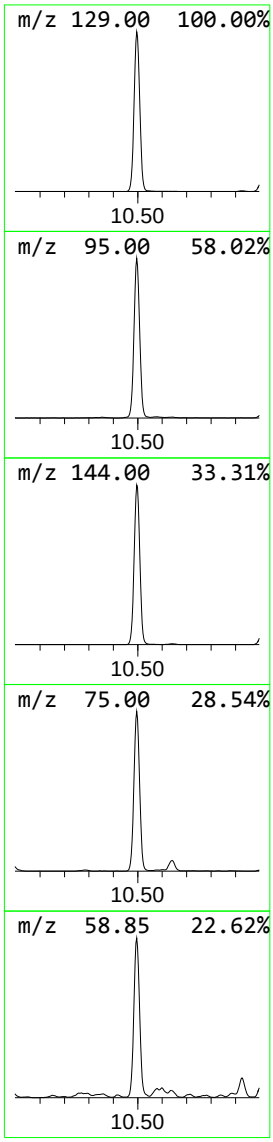
****

Peak Number 5 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.496	24.09 ug/L	2277120	Chlorobenzene-d5	9.412

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2-dihydro-2-methyl-	144	C11H12	021564-79-4	32
2	Ethanone, 1-(2-ethynylphenyl)-	144	C10H8O	104190-22-9	25
3	(1-Methylbuta-1,3-dienyl)benzene	144	C11H12	054758-36-0	23
4	Benzene, (2-cyclopropylethenyl)-	144	C11H12	1000142-01-6	23
5	(1-Methylenebut-2-enyl)benzene	144	C11H12	070588-46-4	23





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070824\
Data File : VU059876.D
Acq On : 08 Jul 2024 12:20
Operator : MD/SY
Sample : P3137-17
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZG6

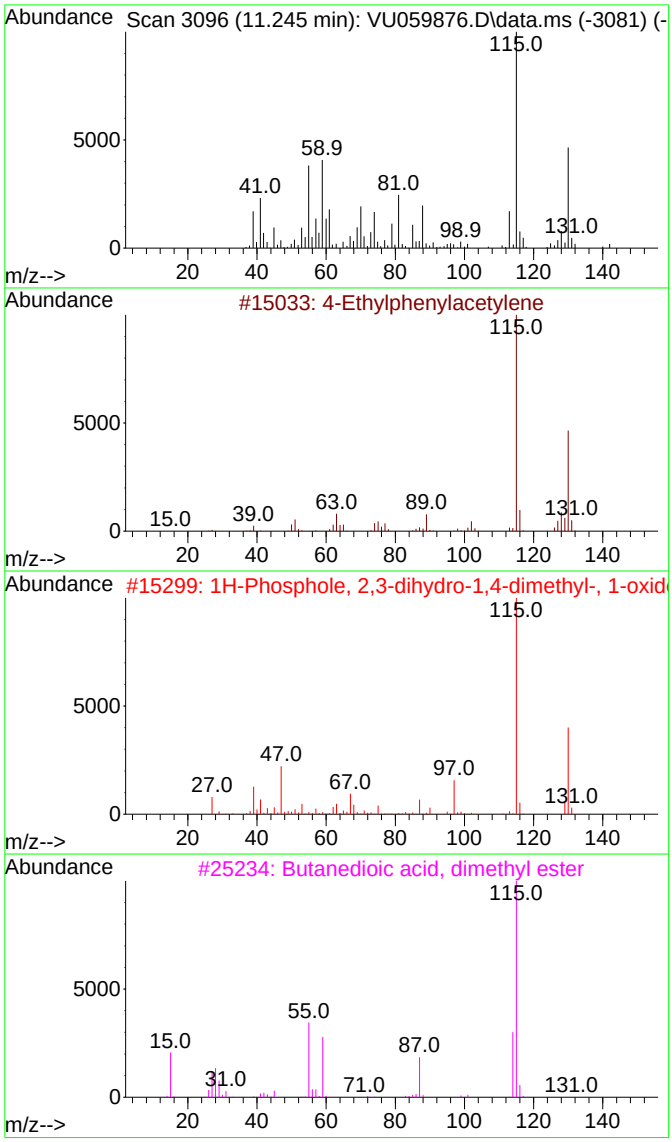
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR061724WMA.M
Quant Title : TRACE VOA SFAM1.0

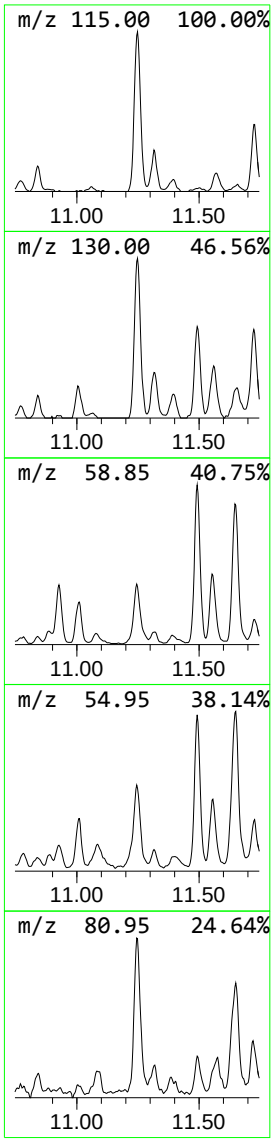
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 unknown-02 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.245	2.75 ug/L	357794	1,4-Dichlorobenzene-d4	11.807

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Ethylphenylacetylene	130	C10H10	040307-11-7	43
2	1H-Phosphole, 2,3-dihydro-1,4-di...	130	C6H11OP	015450-68-7	38
3	Butanedioic acid, dimethyl ester	146	C6H10O4	000106-65-0	38
4	Butanoic acid, 2,3-dimethyl-2-(1...	172	C10H20O2	055519-33-0	37
5	2,6-Difluoropyridine	115	C5H3F2N	001513-65-1	35





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070824\
Data File : VU059876.D
Acq On : 08 Jul 2024 12:20
Operator : MD/SY
Sample : P3137-17
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZG6

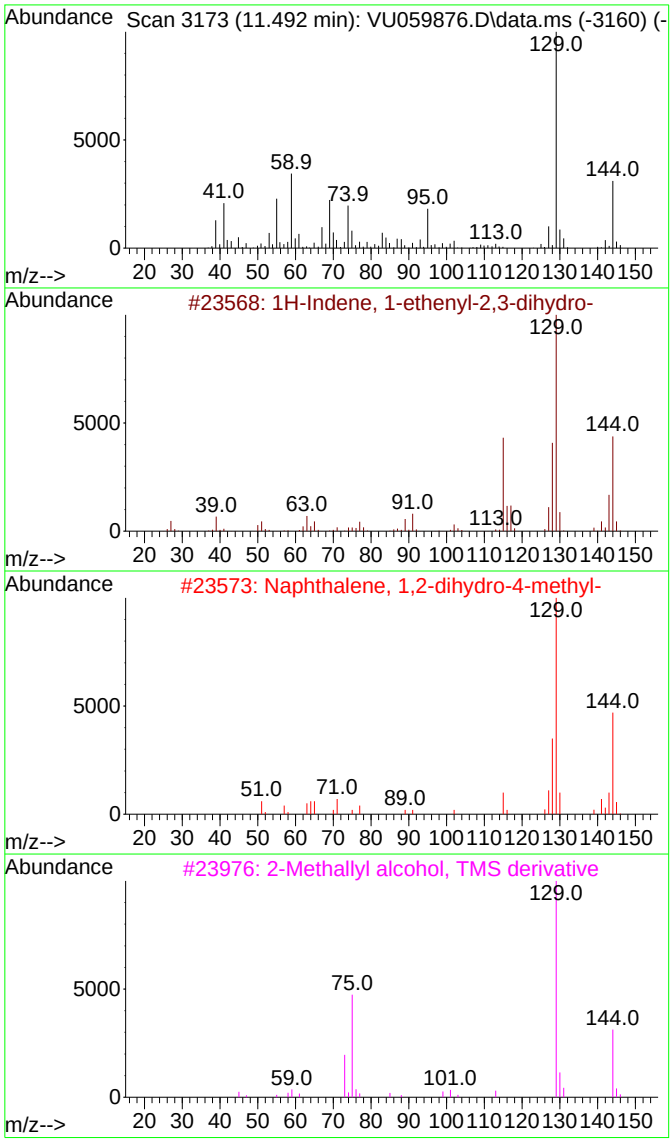
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR061724WMA.M
Quant Title : TRACE VOA SFAM1.0

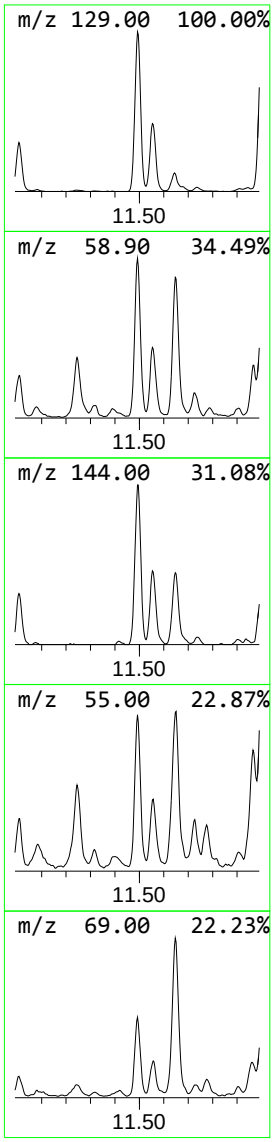
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 1H-Indene, 1-ethenyl-2,3-di... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.492	4.01 ug/L	520754	1,4-Dichlorobenzene-d4	11.807

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 1-ethenyl-2,3-dihydro-	144	C11H12	051783-46-1	52
2	Naphthalene, 1,2-dihydro-4-methyl-	144	C11H12	004373-13-1	50
3	2-Methallyl alcohol, TMS derivative	144	C7H16OSi	025195-85-1	43
4	1,3-Disilacyclobutane, 1,1,3,3-t...	144	C6H16Si2	001627-98-1	40
5	Naphthalene, 1,2-dihydro-2-methyl-	144	C11H12	021564-79-4	40





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070824\
Data File : VU059876.D
Acq On : 08 Jul 2024 12:20
Operator : MD/SY
Sample : P3137-17
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZG6

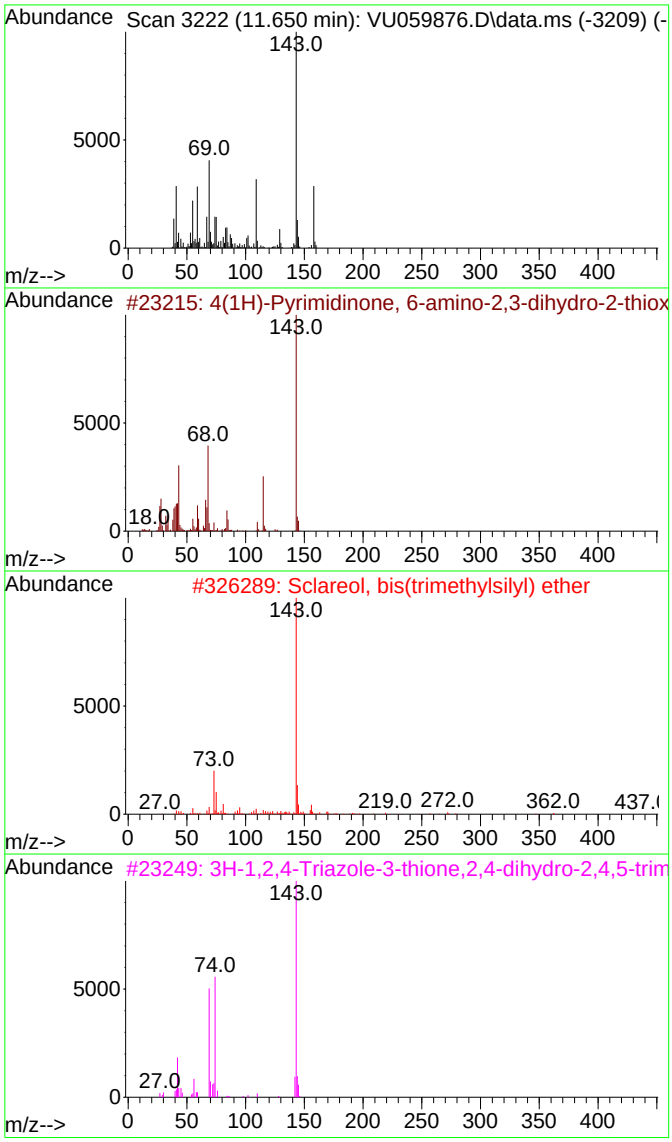
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR061724WMA.M
Quant Title : TRACE VOA SFAM1.0

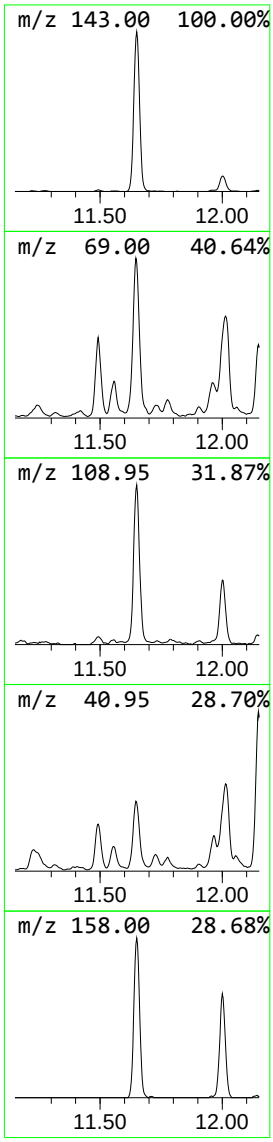
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 4(1H)-Pyrimidinone, 6-amino... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.650	5.45 ug/L	707753	1,4-Dichlorobenzene-d4	11.807

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4(1H)-Pyrimidinone, 6-amino-2,3-...	143	C4H5N3OS	001004-40-6	50
2	Sclareol, bis(trimethylsilyl) ether	452	C26H52O2Si2	1000463-82-9	43
3	3H-1,2,4-Triazole-3-thione,2,4-d...	143	C5H9N3S	037526-42-4	43
4	5-Nonanol, 5-butyl-	200	C13H28O	000597-93-3	38
5	Quinoline, 2-methyl-	143	C10H9N	000091-63-4	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070824\
Data File : VU059876.D
Acq On : 08 Jul 2024 12:20
Operator : MD/SY
Sample : P3137-17
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZG6

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR061724WMA.M
Quant Title : TRACE VOA SFAM1.0

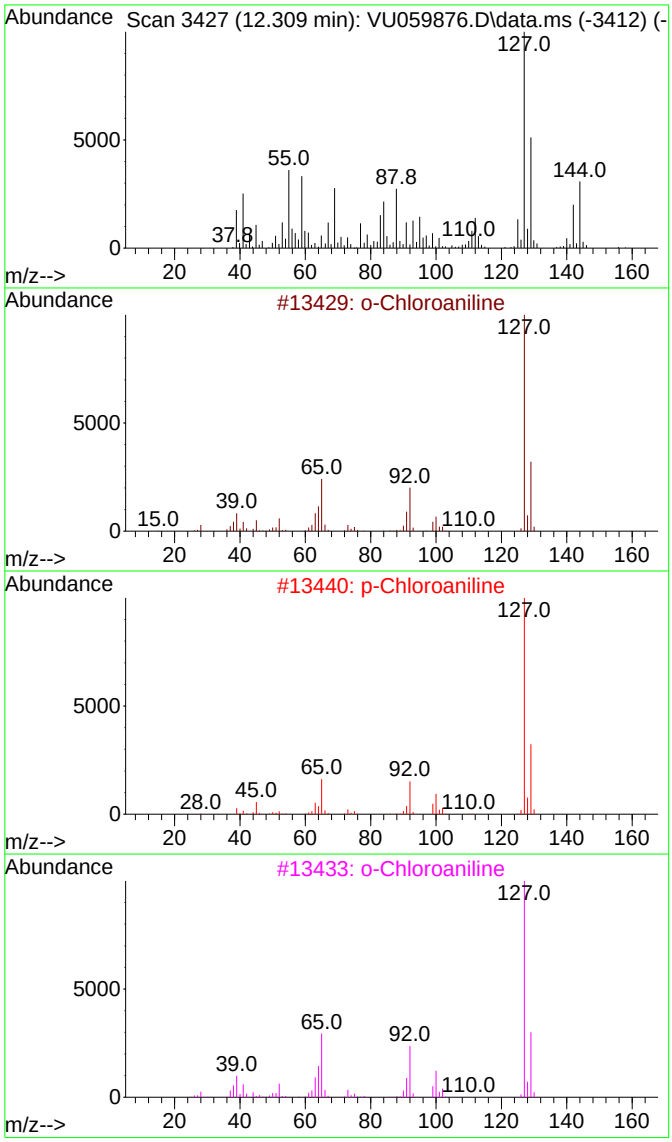
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

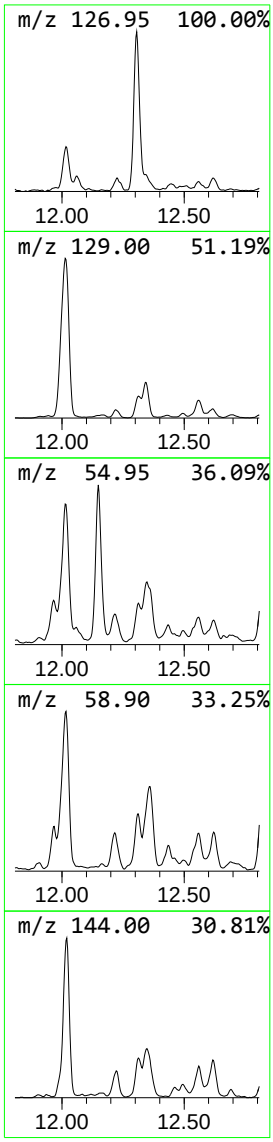
****

Peak Number 10 unknown-03 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.309	3.08 ug/L	399825	1,4-Dichlorobenzene-d4	11.807

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Chloroaniline	127	C6H6ClN	000095-51-2	35
2	p-Chloroaniline	127	C6H6ClN	000106-47-8	35
3	o-Chloroaniline	127	C6H6ClN	000095-51-2	30
4	o-Chloroaniline	127	C6H6ClN	000095-51-2	30
5	p-Chloroaniline	127	C6H6ClN	000106-47-8	30





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070824\
Data File : VU059876.D
Acq On : 08 Jul 2024 12:20
Operator : MD/SY
Sample : P3137-17
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZG6

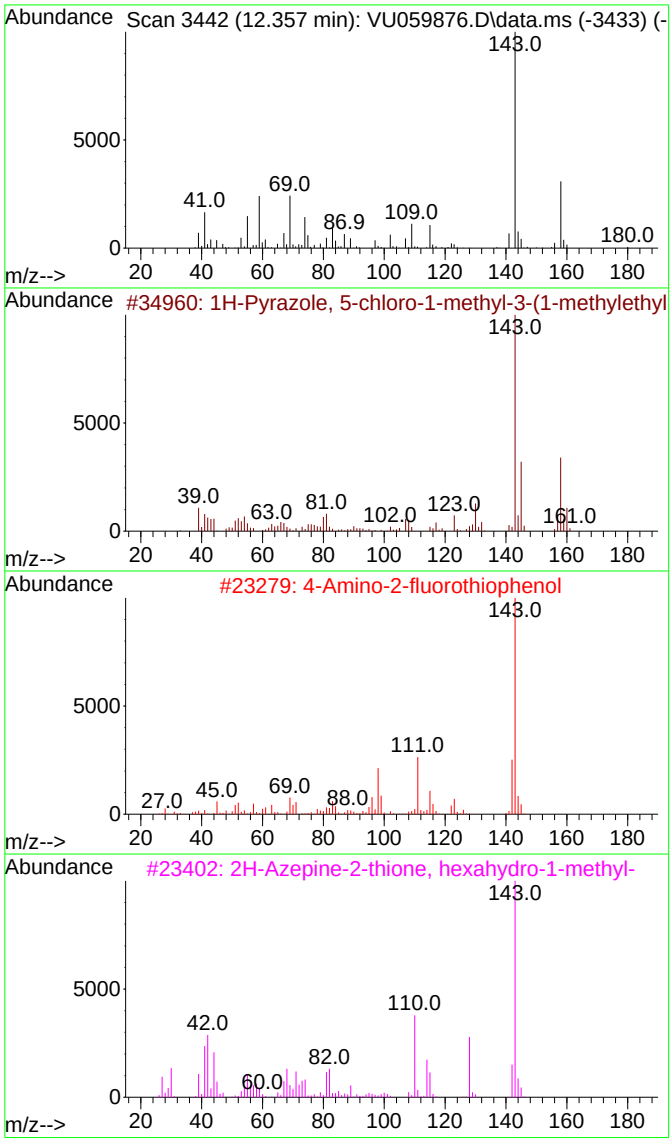
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR061724WMA.M
Quant Title : TRACE VOA SFAM1.0

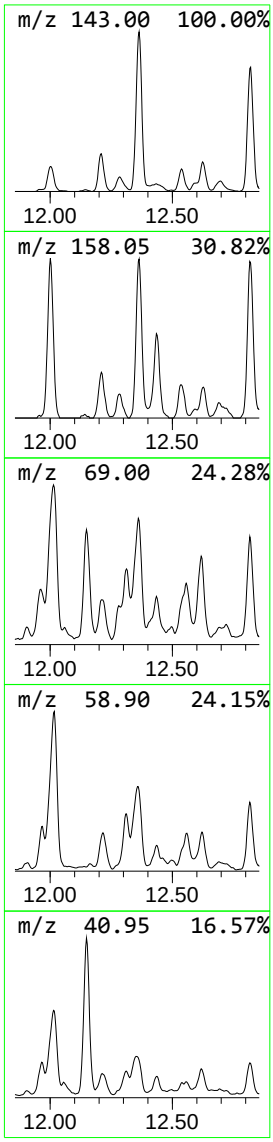
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 1H-Pyrazole, 5-chloro-1-met... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.357	4.20 ug/L	545357	1,4-Dichlorobenzene-d4	11.807

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Pyrazole, 5-chloro-1-methyl-3...	158	C7H11ClN2	1000319-90-4	50
2	4-Amino-2-fluorothiophenol	143	C6H6FNS	015178-52-6	47
3	2H-Azepine-2-thione, hexahydro-1...	143	C7H13NS	034949-15-0	47
4	Quinoline, 2-methyl-	143	C10H9N	000091-63-4	43
5	5-Methyl-2-N-methylethylamino-2-...	158	C7H14N2S	077158-37-3	42





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070824\
Data File : VU059876.D
Acq On : 08 Jul 2024 12:20
Operator : MD/SY
Sample : P3137-17
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZG6

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR061724WMA.M
Quant Title : TRACE VOA SFAM1.0

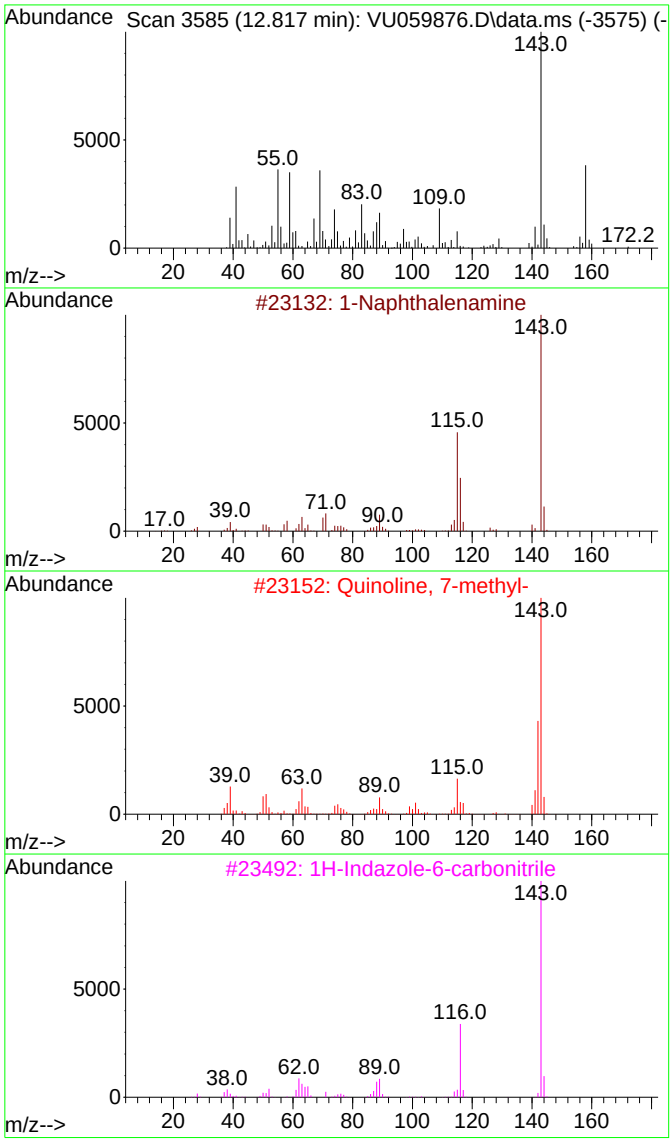
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

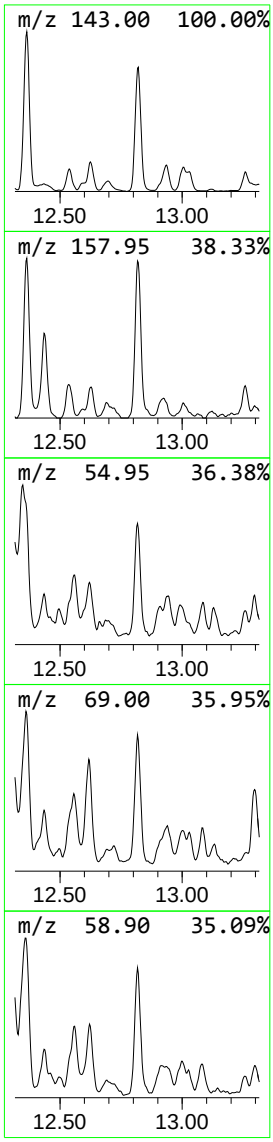
****

Peak Number 12 unknown-04 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.817	2.74 ug/L	356461	1,4-Dichlorobenzene-d4	11.807

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Naphthalenamine	143	C10H9N	000134-32-7	38
2	Quinoline, 7-methyl-	143	C10H9N	000612-60-2	35
3	1H-Indazole-6-carbonitrile	143	C8H5N3	141290-59-7	35
4	5-Nonanol, 5-butyl-	200	C13H28O	000597-93-3	35
5	Quinoline, 2-methyl-	143	C10H9N	000091-63-4	35





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070824\
Data File : VU059876.D
Acq On : 08 Jul 2024 12:20
Operator : MD/SY
Sample : P3137-17
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZG6

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR061724WMA.M
Quant Title : TRACE VOA SFAM1.0

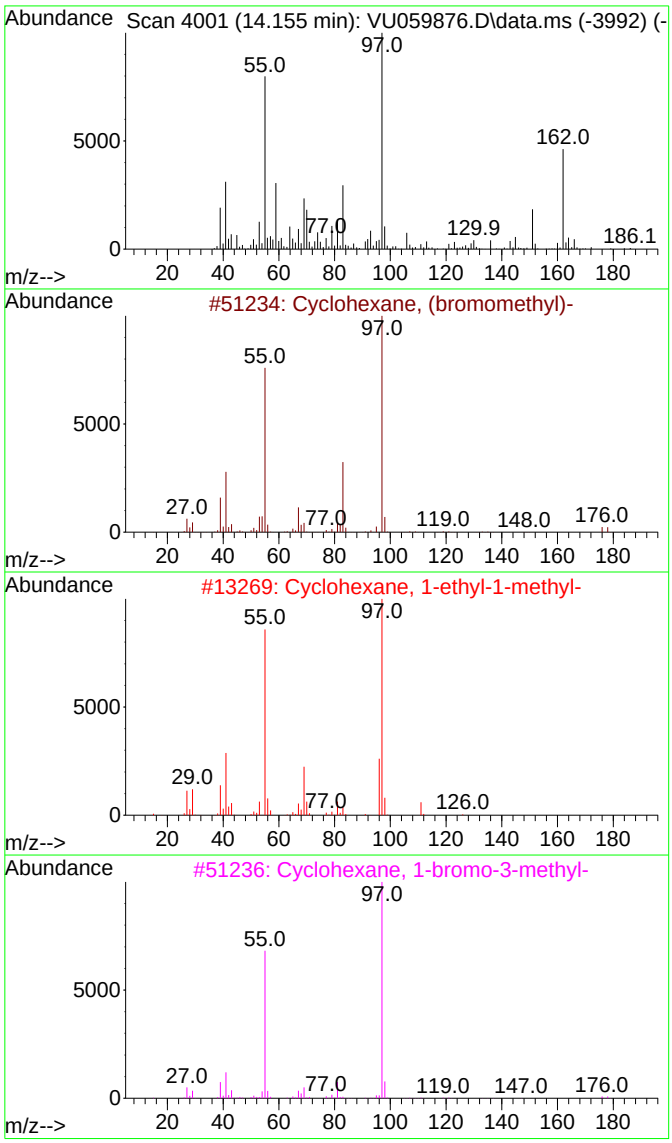
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

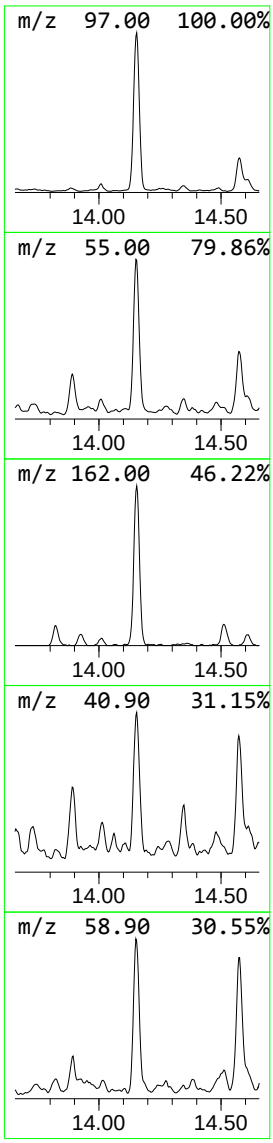
****

Peak Number 13 unknown-05 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.155	4.53 ug/L	588839	1,4-Dichlorobenzene-d4	11.807

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, (bromomethyl)-	176	C7H13Br	002550-36-9	47
2	Cyclohexane, 1-ethyl-1-methyl-	126	C9H18	004926-90-3	38
3	Cyclohexane, 1-bromo-3-methyl-	176	C7H13Br	013905-48-1	38
4	4-Octen-3-one	126	C8H14O	014129-48-7	38
5	Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070824\
Data File : VU059876.D
Acq On : 08 Jul 2024 12:20
Operator : MD/SY
Sample : P3137-17
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZG6

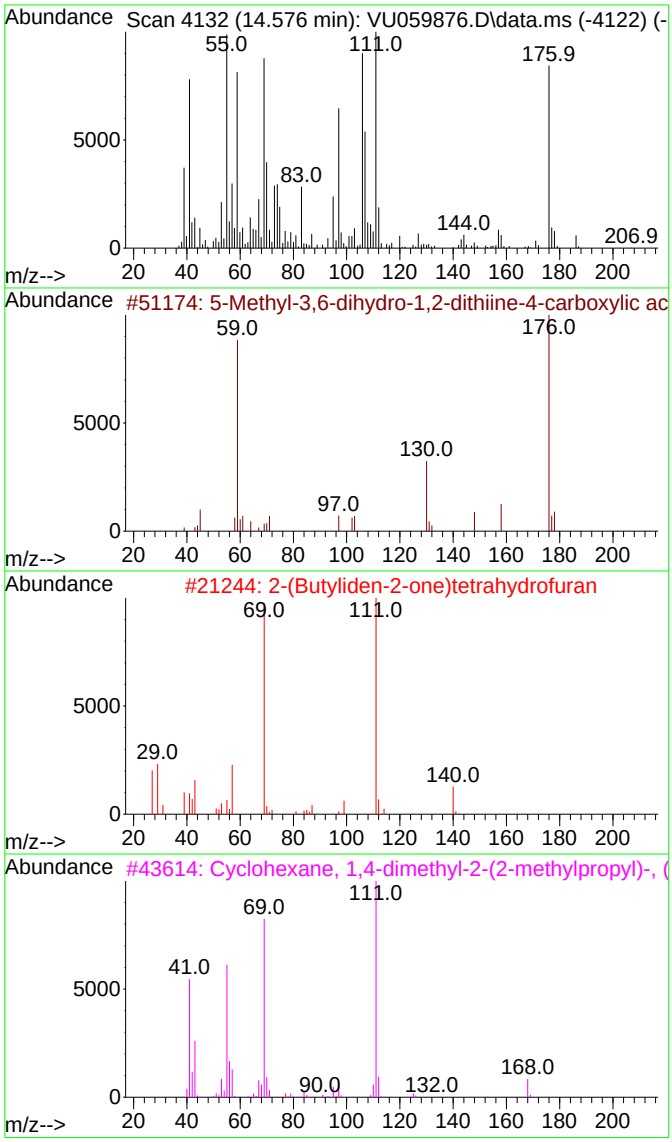
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR061724WMA.M
Quant Title : TRACE VOA SFAM1.0

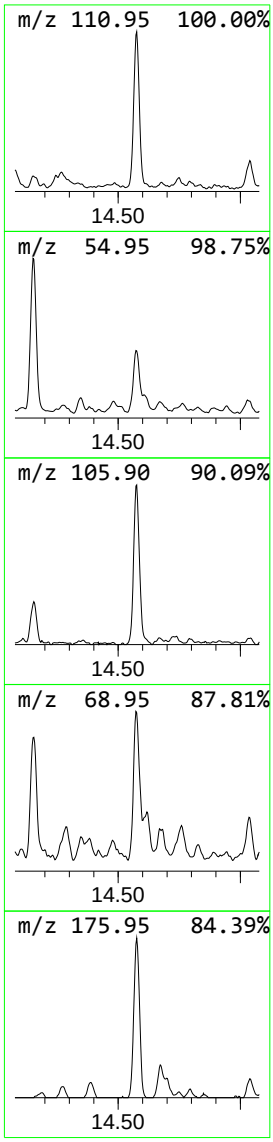
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 unknown-06 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.576	2.96 ug/L	385022	1,4-Dichlorobenzene-d4	11.807

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Methyl-3,6-dihydro-1,2-dithiin...	176	C6H8O2S2	959252-25-6	20
2	2-(Butyliden-2-one)tetrahydrofuran	140	C8H12O2	343270-50-8	14
3	Cyclohexane, 1,4-dimethyl-2-(2-m...	168	C12H24	054411-17-5	14
4	1,3-Dimethyl-5-n-decylcyclohexane	252	C18H36	086710-36-3	12
5	1,1,4-Trimethylcyclohexane	126	C9H18	007094-27-1	12





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070824\
Data File : VU059876.D
Acq On : 08 Jul 2024 12:20
Operator : MD/SY
Sample : P3137-17
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZG6

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR061724WMA.M
Quant Title : TRACE VOA SFAM1.0

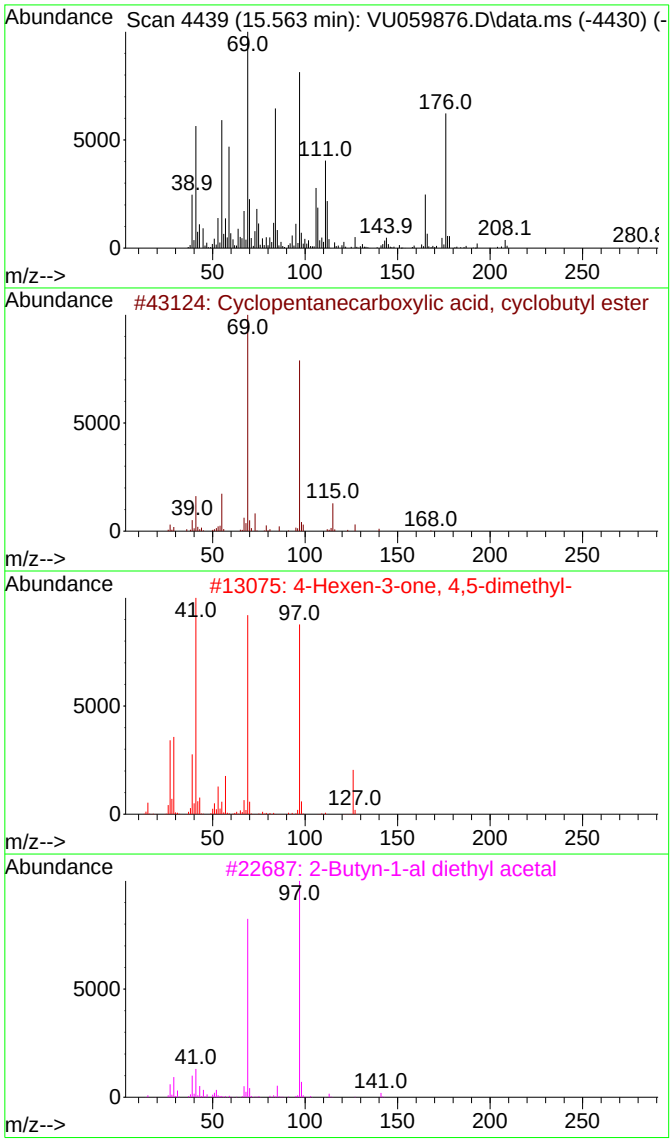
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

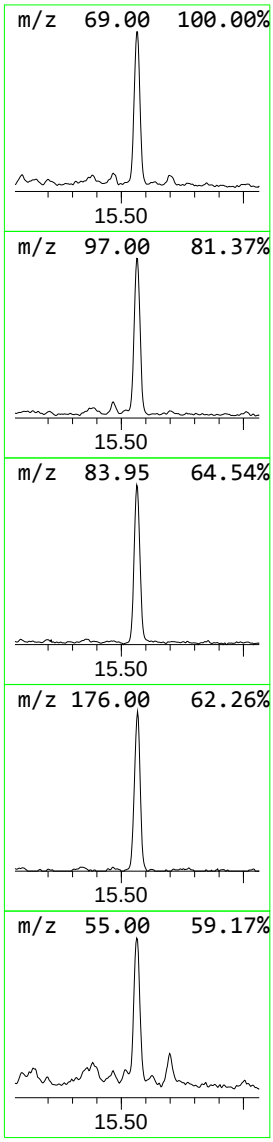
****

Peak Number 15 unknown-07 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.563	3.63 ug/L	471574	1,4-Dichlorobenzene-d4	11.807

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentanecarboxylic acid, cyc...	168	C10H16O2	1000282-76-8	30
2	4-Hexen-3-one, 4,5-dimethyl-	126	C8H14O	017325-90-5	27
3	2-Butyn-1-al diethyl acetal	142	C8H14O2	002806-97-5	27
4	Cyclopentane, 1-butyl-2-ethyl-	154	C11H22	072993-32-9	22
5	Cyclohexane, 1,2-dimethyl-, trans-	112	C8H16	006876-23-9	18





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070824\
Data File : VU059876.D
Acq On : 08 Jul 2024 12:20
Operator : MD/SY
Sample : P3137-17
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZG6

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR061724WMA.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
Propane, 2-(eth...	7.570	7.5	ug/L	680831	1	6.242	451840	5.0
Diisopropyl sul...	8.306	14.8	ug/L	1402310	2	9.412	472600	5.0
Propane, 1-[(1-...	9.187	5.5	ug/L	522303	2	9.412	472600	5.0
Sulfide, sec-bu...	9.875	4.3	ug/L	407919	2	9.412	472600	5.0
unknown-01	10.496	24.1	ug/L	2277120	2	9.412	472600	5.0
unknown-02	11.245	2.8	ug/L	357794	3	11.807	649723	5.0
1H-Indene, 1-et...	11.492	4.0	ug/L	520754	3	11.807	649723	5.0
4(1H)-Pyrimidin...	11.650	5.5	ug/L	707753	3	11.807	649723	5.0
unknown-03	12.309	3.1	ug/L	399825	3	11.807	649723	5.0
1H-Pyrazole, 5-...	12.357	4.2	ug/L	545357	3	11.807	649723	5.0
unknown-04	12.817	2.7	ug/L	356461	3	11.807	649723	5.0
unknown-05	14.155	4.5	ug/L	588839	3	11.807	649723	5.0
unknown-06	14.576	3.0	ug/L	385022	3	11.807	649723	5.0
unknown-07	15.563	3.6	ug/L	471574	3	11.807	649723	5.0