LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU111022\
Data File : VU051807.D
Acq On : 10 Nov 2022 17:51
Operator : JC/MD
Sample : N5571-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GBMW1

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\SFAMUTR102822WMA.M
Title : TRACE VOA SFAM1.0

Signal : TIC: VU051807.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.411	15	22	26	rBV	25912	28632	3.43%	0.364%
2	1.437	26	30	45	rVB	22764	32830	3.93%	0.417%
3	1.658	88	99	110	rBV2	89784	138678	16.61%	1.763%
4	1.765	125	132	138	rBV	11752	20358	2.44%	0.259%
5	1.970	189	196	212	rVB	79311	103084	12.35%	1.310%
6	2.636	393	403	419	rVB	151559	259176	31.04%	3.294%
7	4.572	994	1005	1024	rBV4	30652	81099	9.71%	1.031%
8	4.768	1049	1066	1068	rBV4	50689	109180	13.08%	1.388%
9	5.131	1166	1179	1201	rVB	152818	364938	43.70%	4.638%
10	5.832	1374	1397	1424	rBV2	272838	769739	92.18%	9.783%
11	6.353	1545	1559	1578	rBV	243434	456970	54.73%	5.808%
12	6.716	1661	1672	1679	rBV8	4758	10606	1.27%	0.135%
13	6.780	1679	1692	1716	rVB	194035	376024	45.03%	4.779%
14	7.626	1942	1955	1974	rBV	76715	140389	16.81%	1.784%
15	7.948	2041	2055	2089	rBV	305935	559362	66.99%	7.109%
16	8.224	2130	2141	2153	rBV	46092	79470	9.52%	1.010%
17	8.674	2262	2281	2321	rBV	351391	835024	100.00%	10.613%
18	8.822	2321	2327	2345	rVB7	12751	25415	3.04%	0.323%
19	9.086	2398	2409	2416	rBV6	6770	12628	1.51%	0.160%
20	9.440	2508	2519	2543	rBV	318130	567445	67.96%	7.212%
21	9.539	2545	2550	2561	rVB8	8168	12475	1.49%	0.159%
22	10.108	2718	2727	2735	rBV4	13903	25524	3.06%	0.324%
23	10.401	2809	2818	2833	rVB3	7962	19160	2.29%	0.244%
24	10.764	2920	2931	2954	rVB	162552	278049	33.30%	3.534%
25	10.896	2961	2972	2986	rBV9	8439	16191	1.94%	0.206%
26	11.057	3010	3022	3032	rBV3	19904	35604	4.26%	0.453%
27	11.237	3065	3078	3089	rBV3	9193	16815	2.01%	0.214%
28	11.562	3164	3179	3186	rBV8	6377	12744	1.53%	0.162%
29	11.610	3186	3194	3206	rVV3	25157	39831	4.77%	0.506%
30	11.674	3206	3214	3229	rVB9	6799	17728	2.12%	0.225%
31	11.761	3229	3241	3247	rBV9	7246	13972	1.67%	0.178%
32	11.816	3247	3258	3273	rVB	323176	535759	64.16%	6.809%
33	11.935	3286	3295	3301	rBV6	5623	10002	1.20%	0.127%
34	11.980	3305	3309	3322	rVB9	9148	13019	1.56%	0.165%
35	12.060	3323	3334	3347	rBV3	29563	57672	6.91%	0.733%
36	12.195	3365	3376	3395	rVB	353873	576217	69.01%	7.323%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU111022\
Data File : VU051807.D
Acq On : 10 Nov 2022 17:51
Operator : JC/MD
Sample : N5571-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GBMW1

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\SFAMUTR102822WMA.M
Title : TRACE VOA SFAM1.0

37	12.478	3459	3464	3472	rVV5	4688	8424	1.01%	0.107%
38	12.767	3546	3554	3564	rVB2	19286	27829	3.33%	0.354%
39	12.890	3582	3592	3603	rVB2	11735	21393	2.56%	0.272%
40	13.034	3629	3637	3648	rVB2	9197	16708	2.00%	0.212%
41	13.224	3692	3696	3704	rVB9	8305	9858	1.18%	0.125%
42	13.285	3710	3715	3722	rVB6	7761	10137	1.21%	0.129%
43	13.459	3763	3769	3775	rBV10	6317	9356	1.12%	0.119%
44	13.520	3782	3788	3801	rVB2	18250	27435	3.29%	0.349%
45	13.790	3861	3872	3879	rBV5	15918	30148	3.61%	0.383%
46	13.970	3919	3928	3935	rBV4	53732	97319	11.65%	1.237%
47	14.012	3935	3941	3952	rVB3	47818	77407	9.27%	0.984%
48	14.147	3975	3983	3991	rBV2	60145	100234	12.00%	1.274%
49	14.320	4028	4037	4044	rBV2	88294	139814	16.74%	1.777%
50	14.369	4044	4052	4068	rVB4	97390	172604	20.67%	2.194%
51	14.574	4110	4116	4124	rBV4	21244	29899	3.58%	0.380%
52	14.825	4187	4194	4204	rVB6	36262	63902	7.65%	0.812%
53	15.015	4248	4253	4261	rVB9	15036	19301	2.31%	0.245%
54	15.304	4338	4343	4349	rBV3	21691	23121	2.77%	0.294%
55	15.455	4384	4390	4415	rVB7	34798	73894	8.85%	0.939%
56	15.767	4481	4487	4499	rBV7	13870	23414	2.80%	0.298%
57	15.928	4529	4537	4545	rBV4	79074	119276	14.28%	1.516%
58	16.288	4646	4649	4660	rVB4	25533	33589	4.02%	0.427%
59	16.359	4666	4671	4677	rBV3	19151	26712	3.20%	0.339%
60	16.401	4679	4684	4692	rVB4	39418	54660	6.55%	0.695%

Sum of corrected areas: 7868243

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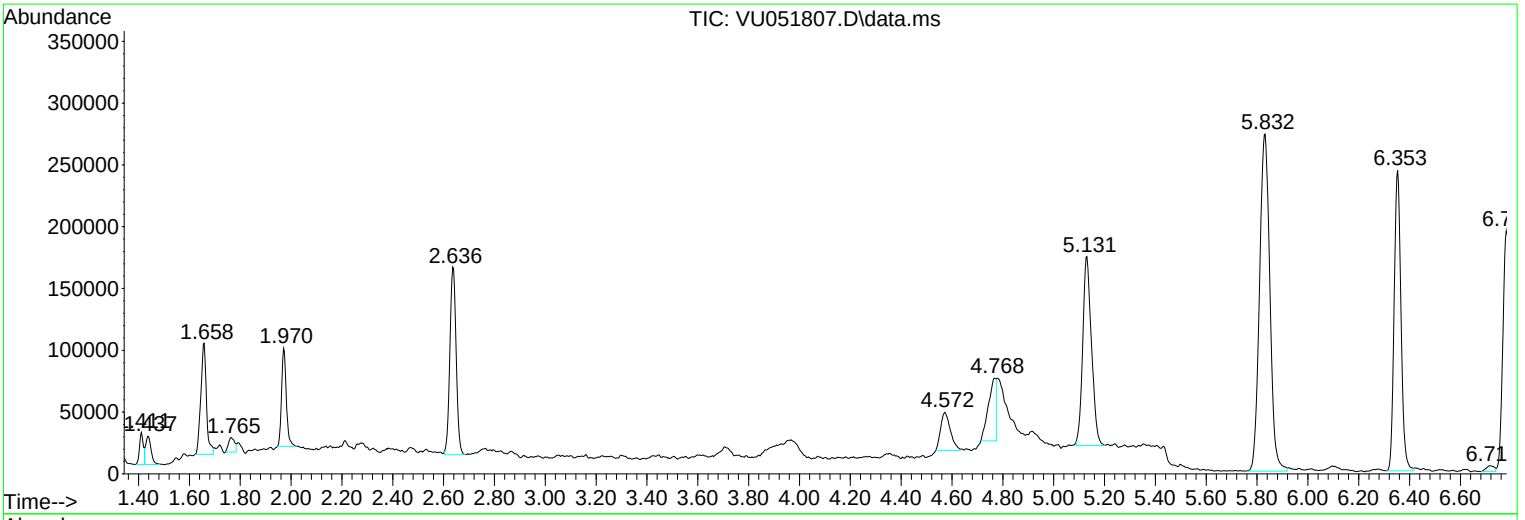
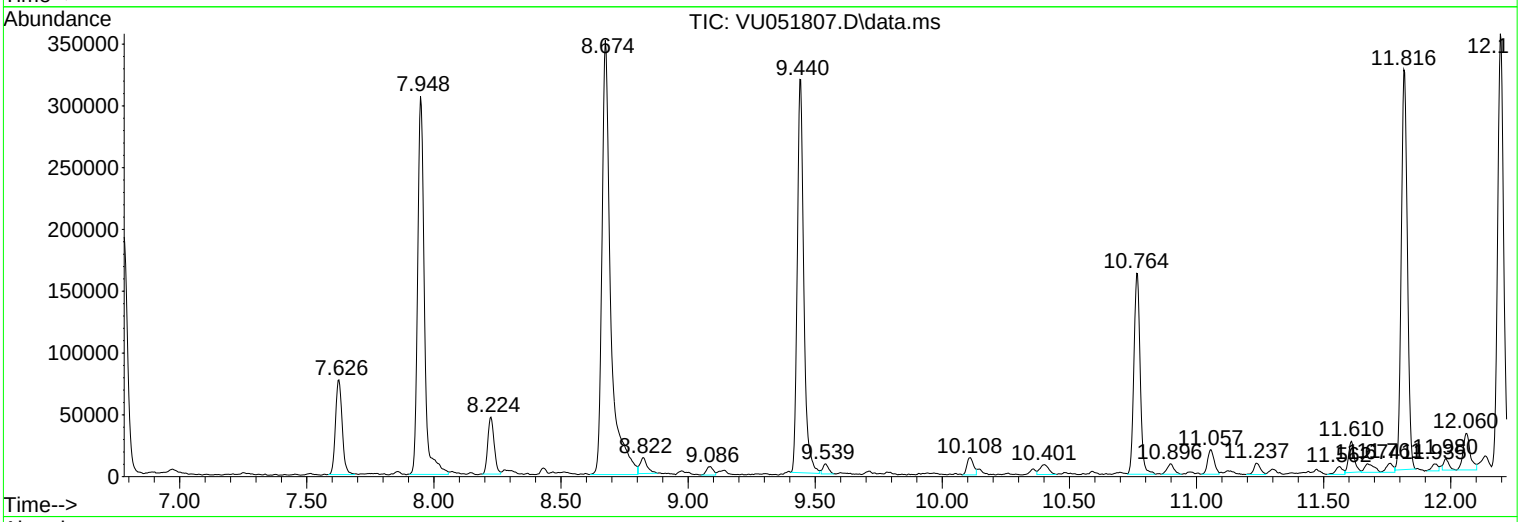
Operator : JC/MD
Sample : N5571-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 12 Sample Multiplier: 1

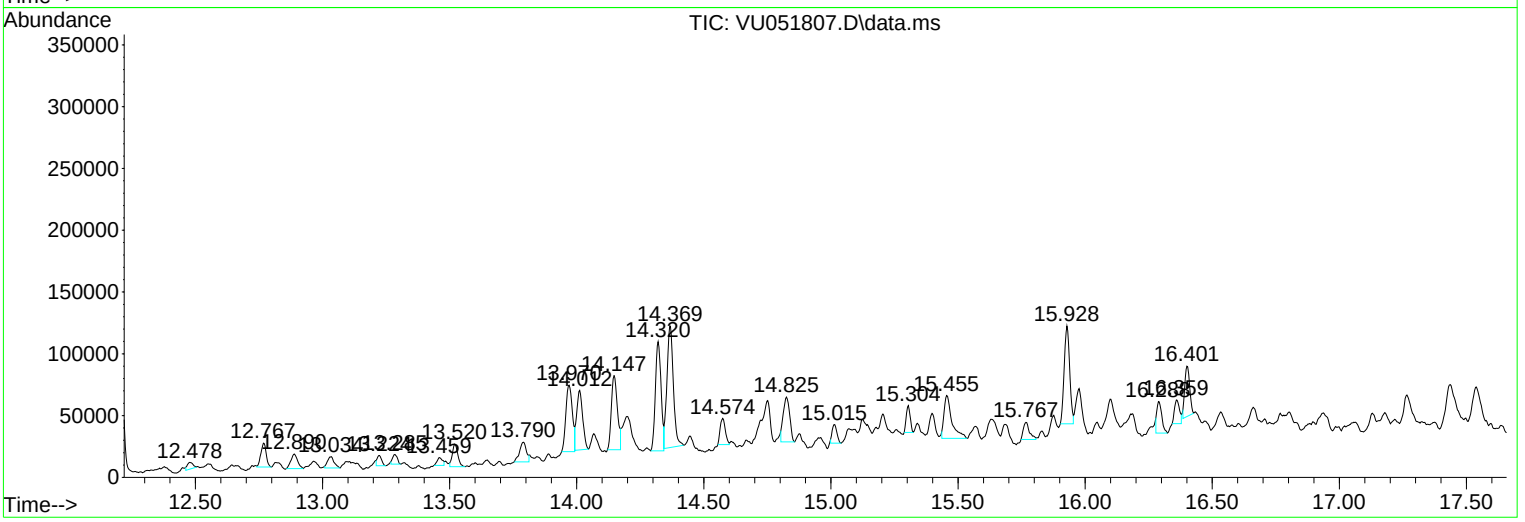
Instrument :
MSVOA_U
ClientSampleId :
GBMW1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102822WMA.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\VU111022\
Data File : VU051807.D
Acq On : 10 Nov 2022 17:51
Operator : JC/MD
Sample : N5571-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GBMW1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102822WMA.M
Quant Title : TRACE VOA SFAM1.0

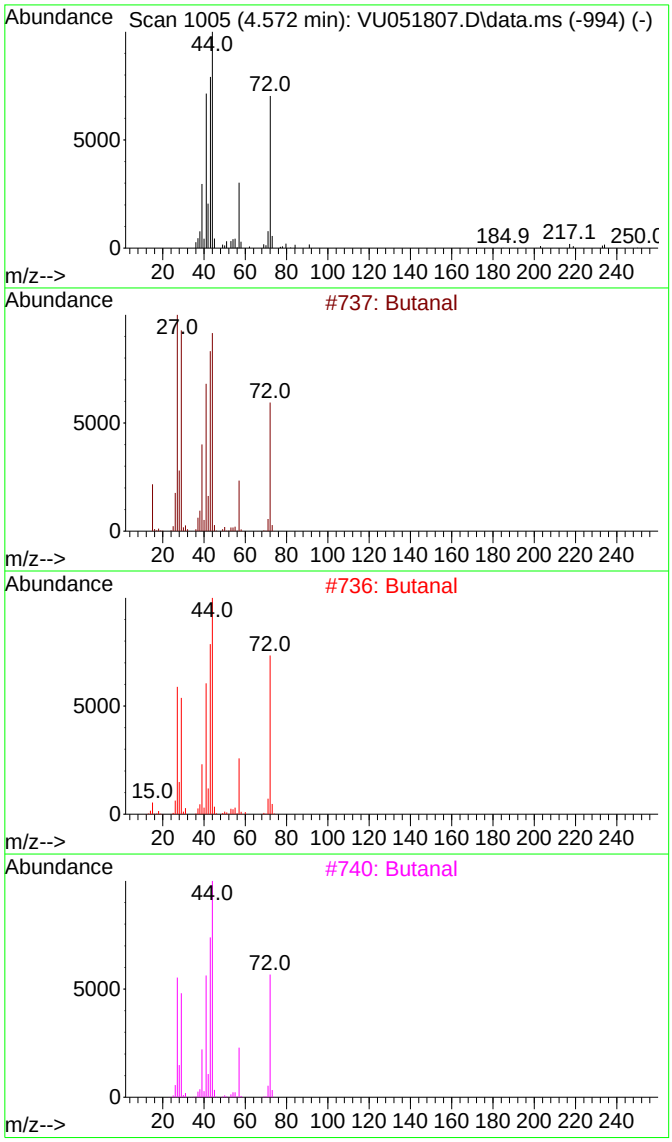
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

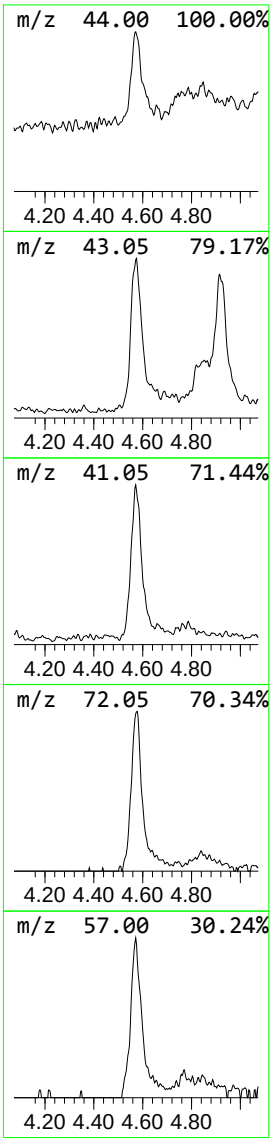
****

Peak Number 1 Butanal Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.572	0.89 ug/L	81099	1,4-Difluorobenzene	6.353

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanal	72	C4H8O	000123-72-8	91
2	Butanal	72	C4H8O	000123-72-8	87
3	Butanal	72	C4H8O	000123-72-8	87
4	Butanal	72	C4H8O	000123-72-8	83
5	Butanal	72	C4H8O	000123-72-8	74





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU111022\
Data File : VU051807.D
Acq On : 10 Nov 2022 17:51
Operator : JC/MD
Sample : N5571-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GBMW1

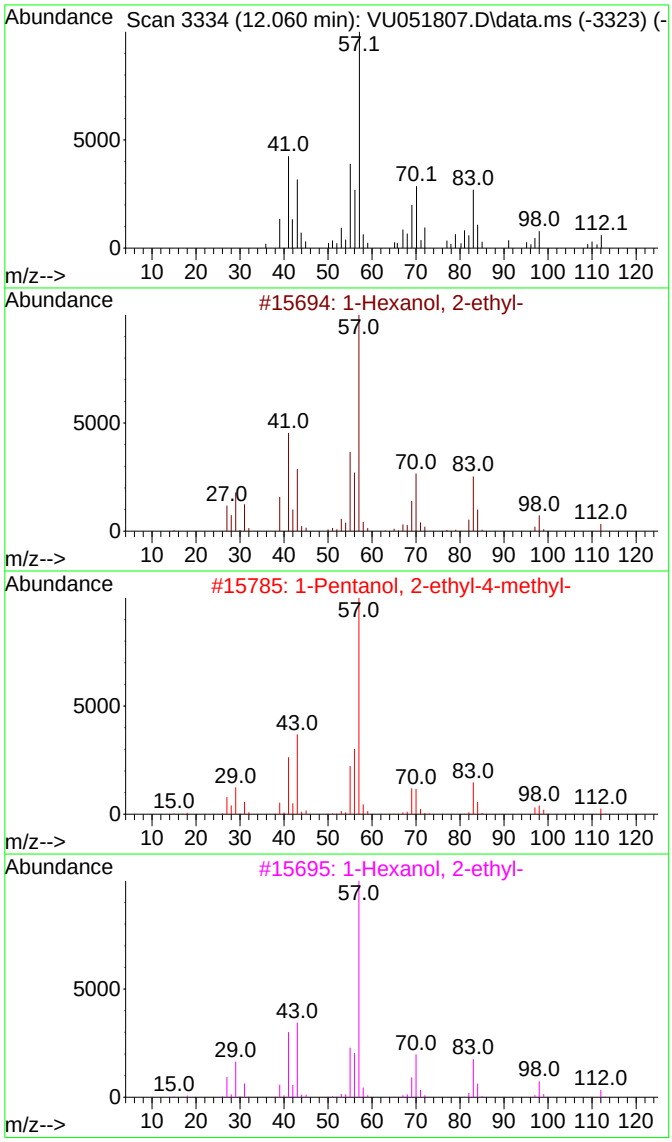
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102822WMA.M
Quant Title : TRACE VOA SFAM1.0

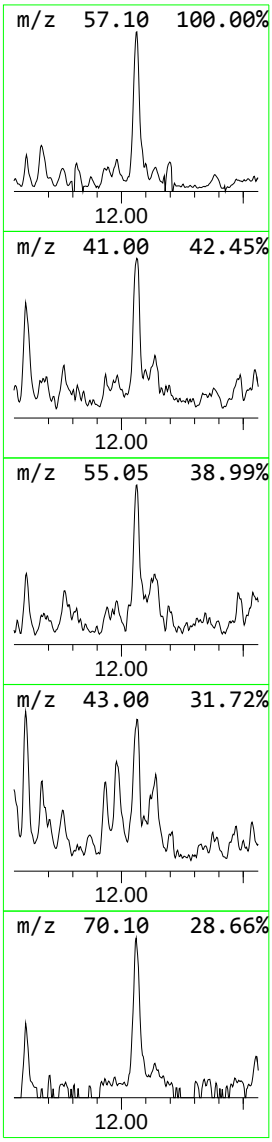
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 1-Hexanol, 2-ethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.060	0.54 ug/L	57672	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	64
2	1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	59
3	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	59
4	2-Propyl-1-pentanol	130	C8H18O	058175-57-8	53
5	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	53





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU111022\
Data File : VU051807.D
Acq On : 10 Nov 2022 17:51
Operator : JC/MD
Sample : N5571-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GBMW1

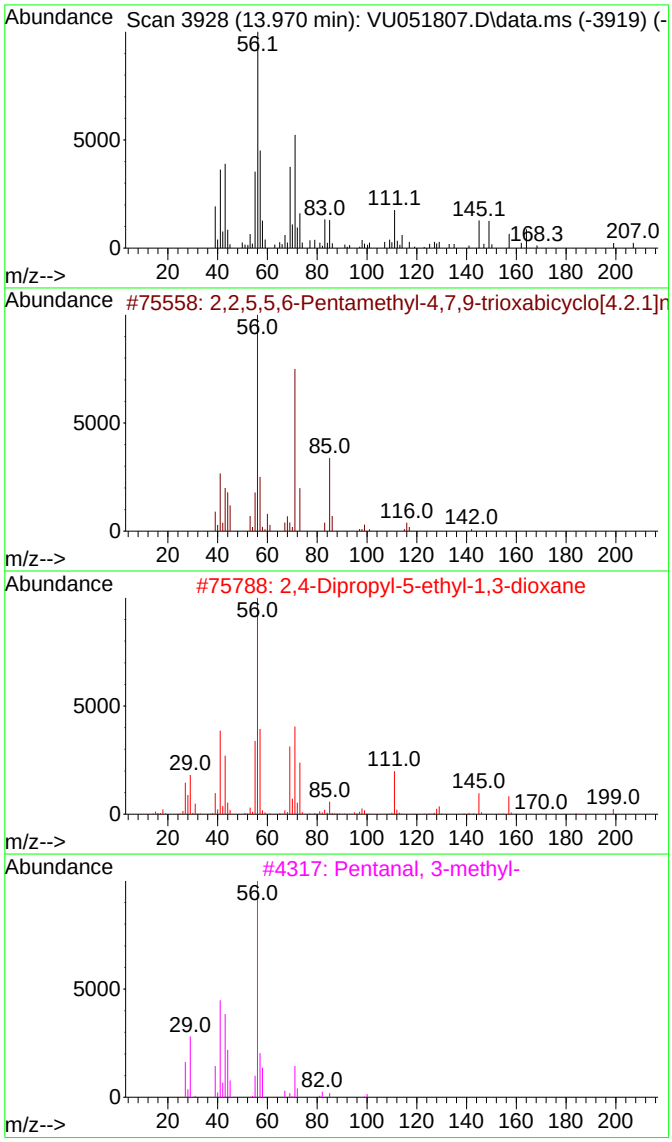
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102822WMA.M
Quant Title : TRACE VOA SFAM1.0

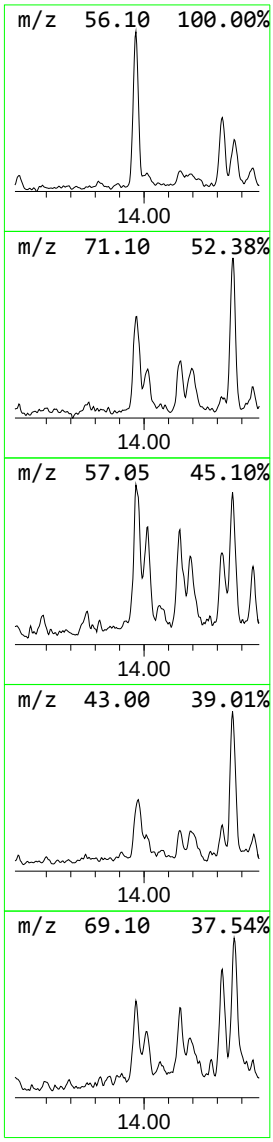
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 2,2,5,5,6-Pentamethyl-4,7,9... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.970	0.91 ug/L	97319	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2,5,5,6-Pentamethyl-4,7,9-trio...	200	C11H20O3	062759-70-0	53
2	2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	38
3	Pentanal, 3-methyl-	100	C6H12O	015877-57-3	38
4	1-Undecene, 10-methyl-	168	C12H24	022370-55-4	35
5	3-Undecene, 7-methyl-, (E)-	168	C12H24	074630-53-8	35





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU111022\
Data File : VU051807.D
Acq On : 10 Nov 2022 17:51
Operator : JC/MD
Sample : N5571-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GBMW1

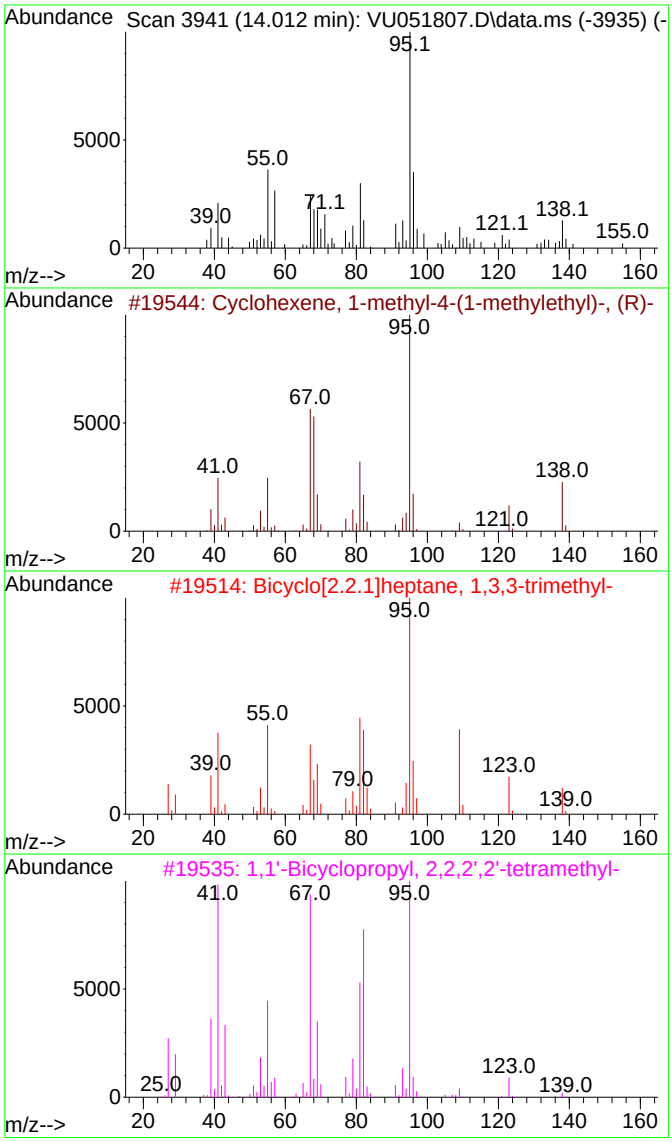
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102822WMA.M
Quant Title : TRACE VOA SFAM1.0

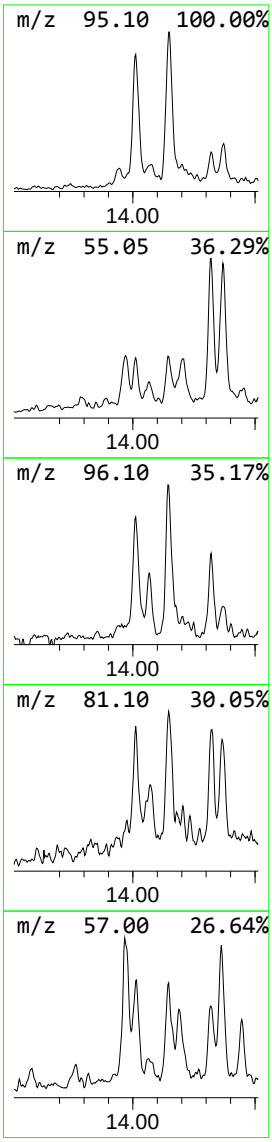
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Cyclohexene, 1-methyl-4-(1-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.012	0.72 ug/L	77407	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexene, 1-methyl-4-(1-methy...	138	C10H18	001195-31-9	59
2	Bicyclo[2.2.1]heptane, 1,3,3-tri...	138	C10H18	006248-88-0	58
3	1,1'-Bicyclopropyl, 2,2,2',2'-te...	138	C10H18	068998-20-9	53
4	Cyclohexene, 1-methyl-3-(1-methy...	138	C10H18	013828-31-4	53
5	Cyclopentene, 1,4-dimethyl-5-(1-...	138	C10H18	061142-33-4	53





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU111022\
Data File : VU051807.D
Acq On : 10 Nov 2022 17:51
Operator : JC/MD
Sample : N5571-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GBMW1

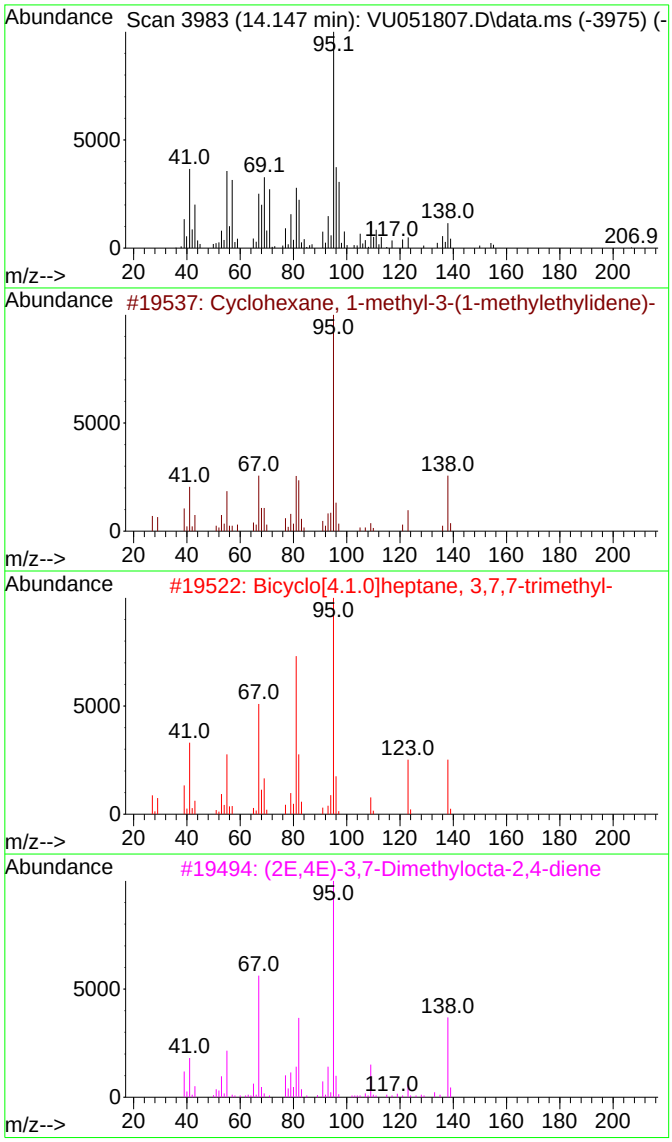
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102822WMA.M
Quant Title : TRACE VOA SFAM1.0

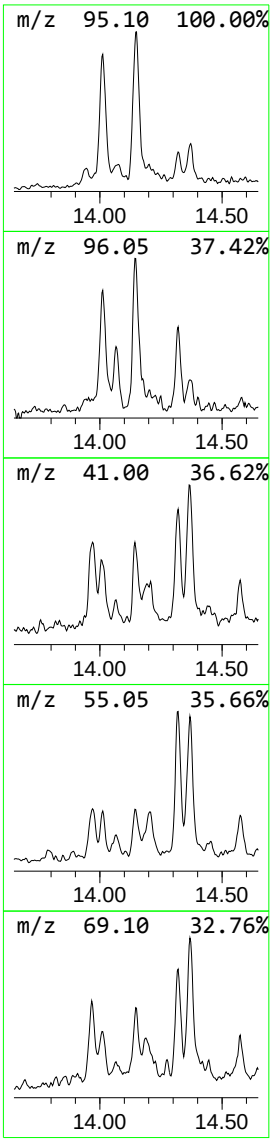
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Cyclohexane, 1-methyl-3-(1-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.147	0.94 ug/L	100234	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1-methyl-3-(1-methy...	138	C10H18	013828-34-7	58
2	Bicyclo[4.1.0]heptane, 3,7,7-tri...	138	C10H18	000554-59-6	53
3	(2E,4E)-3,7-Dimethylocta-2,4-diene	138	C10H18	006874-39-1	52
4	Cyclohexene, 1-methyl-3-vinyloxy-	138	C9H14O	100144-30-7	49
5	1-(1,2-Dimethyl-cyclopent-2-enyl...	138	C9H14O	070987-82-5	46





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU111022\
Data File : VU051807.D
Acq On : 10 Nov 2022 17:51
Operator : JC/MD
Sample : N5571-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GBMW1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102822WMA.M
Quant Title : TRACE VOA SFAM1.0

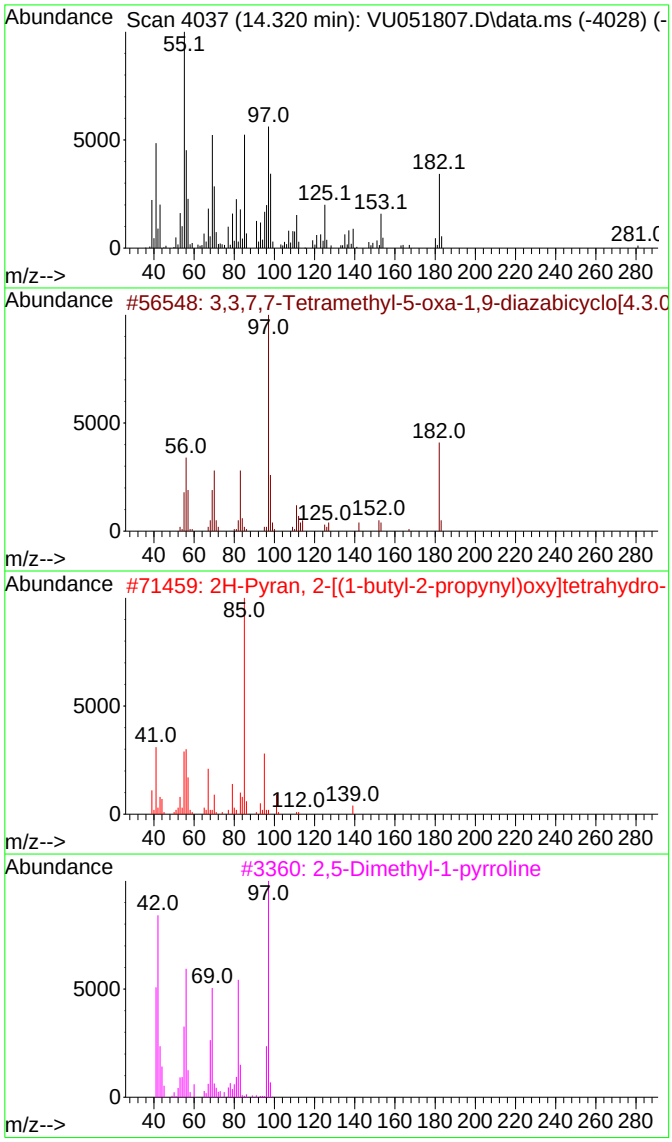
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

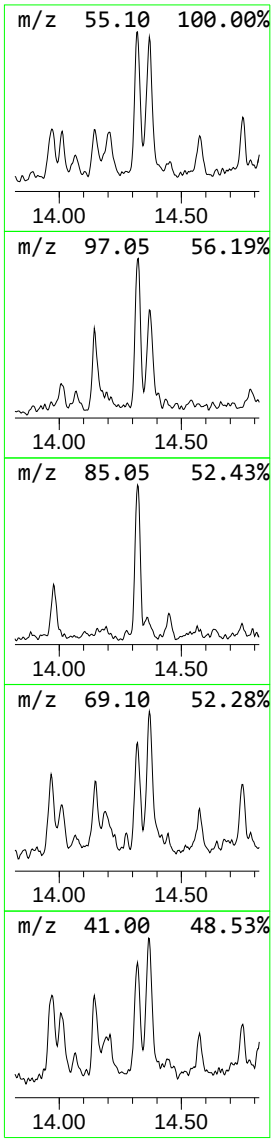
****

Peak Number 7 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.320	1.30 ug/L	139814	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,3,7,7-Tetramethyl-5-oxa-1,9-di...	182	C10H18N2O	090832-25-0	45
2	2H-Pyran, 2-[(1-butyl-2-propynyl...	196	C12H20O2	000831-83-4	27
3	2,5-Dimethyl-1-pyrroline	97	C6H11N	000822-64-0	27
4	Cyclohexane, 2-butyl-1,1,3-trime...	182	C13H26	054676-39-0	27
5	Cyclohexane, 1,1-dimethyl-	112	C8H16	000590-66-9	22





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU111022\
Data File : VU051807.D
Acq On : 10 Nov 2022 17:51
Operator : JC/MD
Sample : N5571-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GBMW1

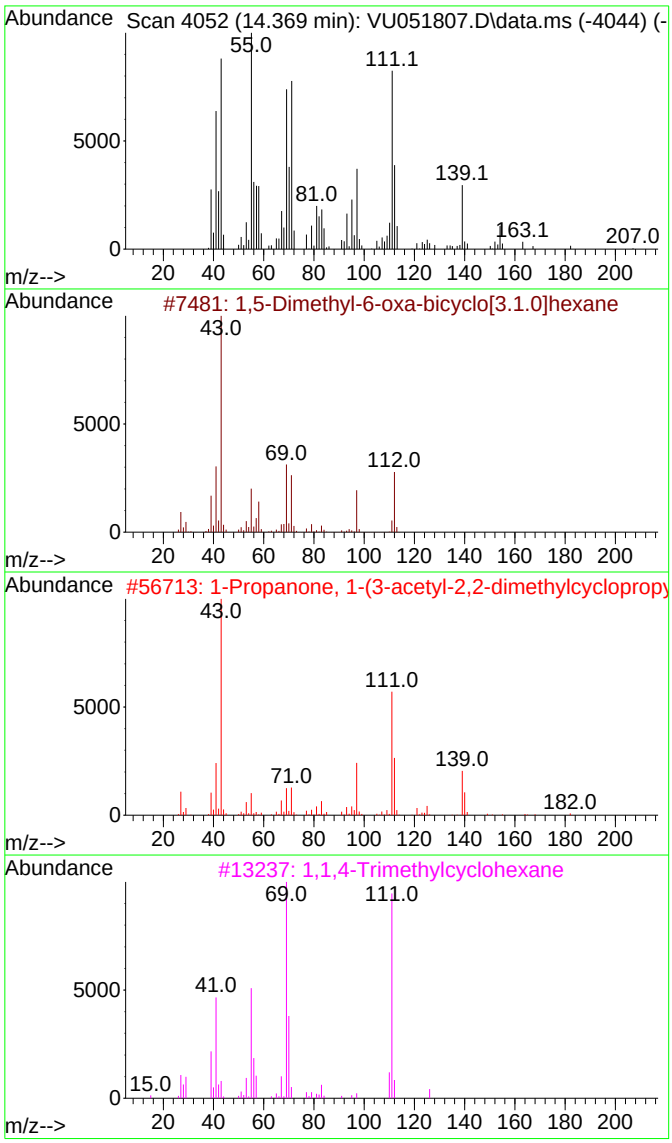
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102822WMA.M
Quant Title : TRACE VOA SFAM1.0

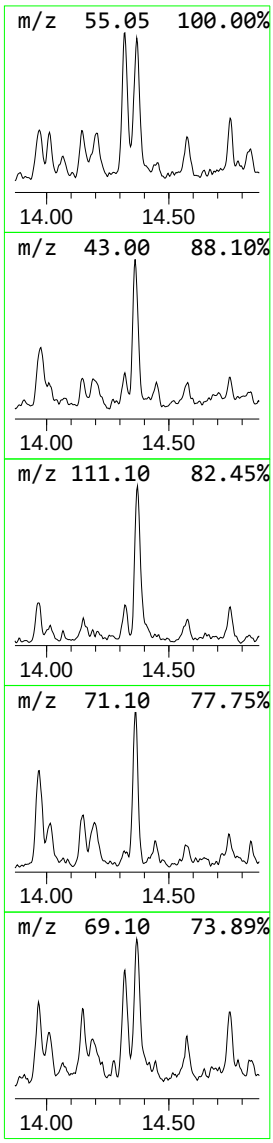
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 1,5-Dimethyl-6-oxa-bicyclo[... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.369	1.61 ug/L	172604	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,5-Dimethyl-6-oxa-bicyclo[3.1.0...	112	C7H12O	082461-31-2	60
2	1-Propanone, 1-(3-acetyl-2,2-dim...	182	C11H18O2	077142-84-8	43
3	1,1,4-Trimethylcyclohexane	126	C9H18	007094-27-1	35
4	1,1,4-Trimethylcyclohexane	126	C9H18	007094-27-1	35
5	2-t-Butyl-5-isopropyl-6-methyl[1...	214	C12H22O3	129287-66-7	35





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Data File : VU051807.D
Acq On : 10 Nov 2022 17:51
Operator : JC/MD
Sample : N5571-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GBMW1

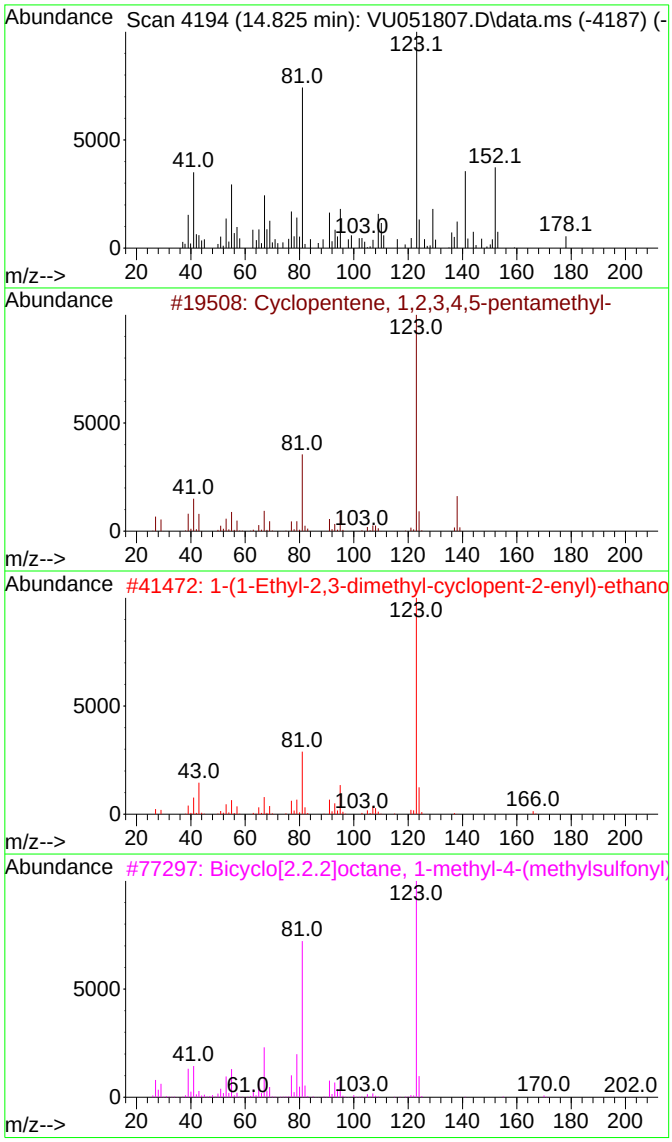
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102822WMA.M
Quant Title : TRACE VOA SFAM1.0

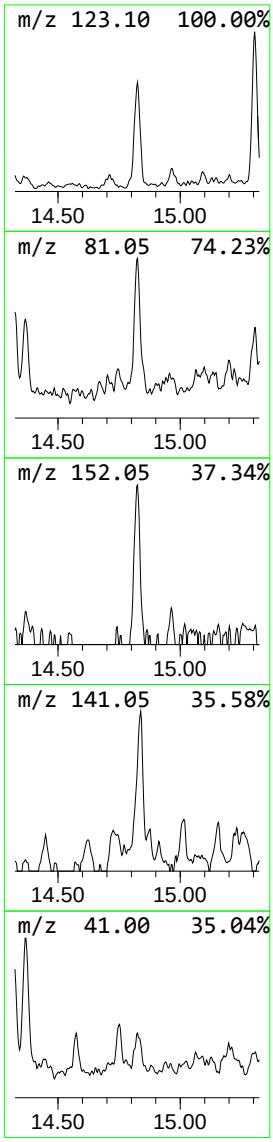
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Cyclopentene, 1,2,3,4,5-pen... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.825	0.60 ug/L	63902	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentene, 1,2,3,4,5-pentamet...	138	C10H18	1000154-28-6	59
2	1-(1-Ethyl-2,3-dimethyl-cyclopen...	166	C11H18O	1000185-18-6	53
3	Bicyclo[2.2.2]octane, 1-methyl-4...	202	C10H18O2S	069855-48-7	50
4	Furan, 2-(1,1-dimethylethyl)-4-m...	138	C9H14O	006141-68-0	42
5	Acetaldehyde, (phenylthio)-	152	C8H8OS	066303-55-7	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU111022\
Data File : VU051807.D
Acq On : 10 Nov 2022 17:51
Operator : JC/MD
Sample : N5571-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GBMW1

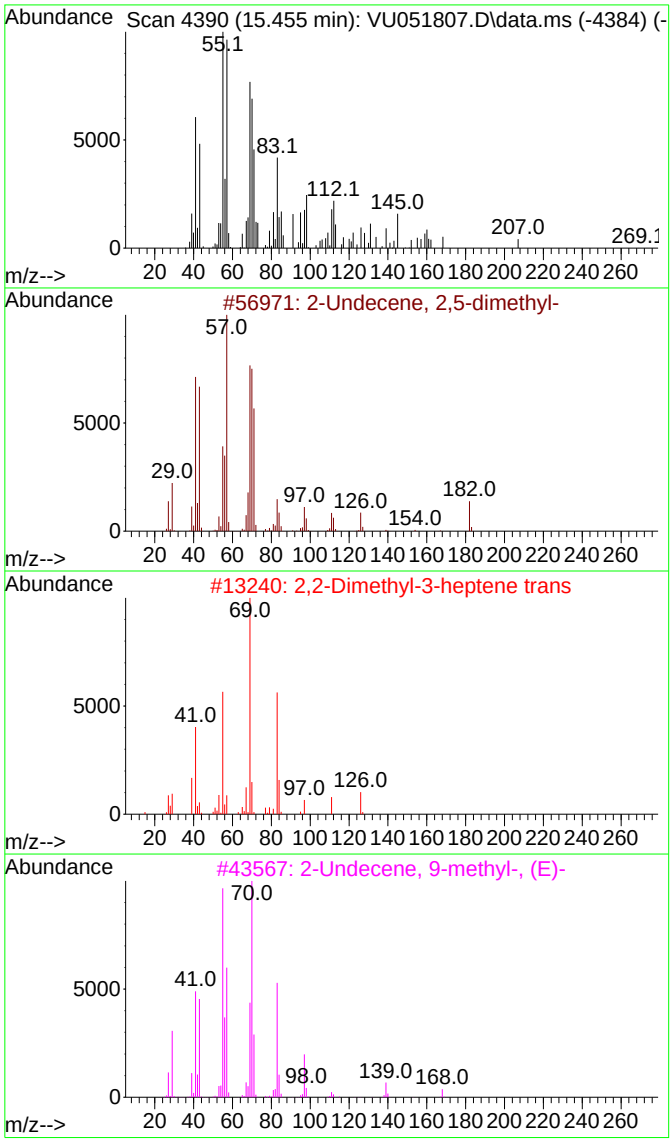
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102822WMA.M
Quant Title : TRACE VOA SFAM1.0

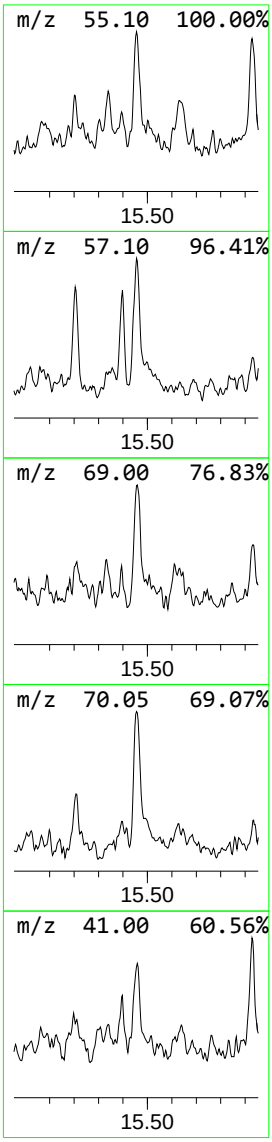
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.455	0.69 ug/L	73894	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Undecene, 2,5-dimethyl-	182	C13H26	049622-16-4	50
2	2,2-Dimethyl-3-heptene trans	126	C9H18	019550-75-5	43
3	2-Undecene, 9-methyl-, (E)-	168	C12H24	074630-46-9	42
4	3-Heptene, 4-methyl-	112	C8H16	004485-16-9	38
5	3-Heptene, 4-methyl-	112	C8H16	004485-16-9	35





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU111022\
Data File : VU051807.D
Acq On : 10 Nov 2022 17:51
Operator : JC/MD
Sample : N5571-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GBMW1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102822WMA.M
Quant Title : TRACE VOA SFAM1.0

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

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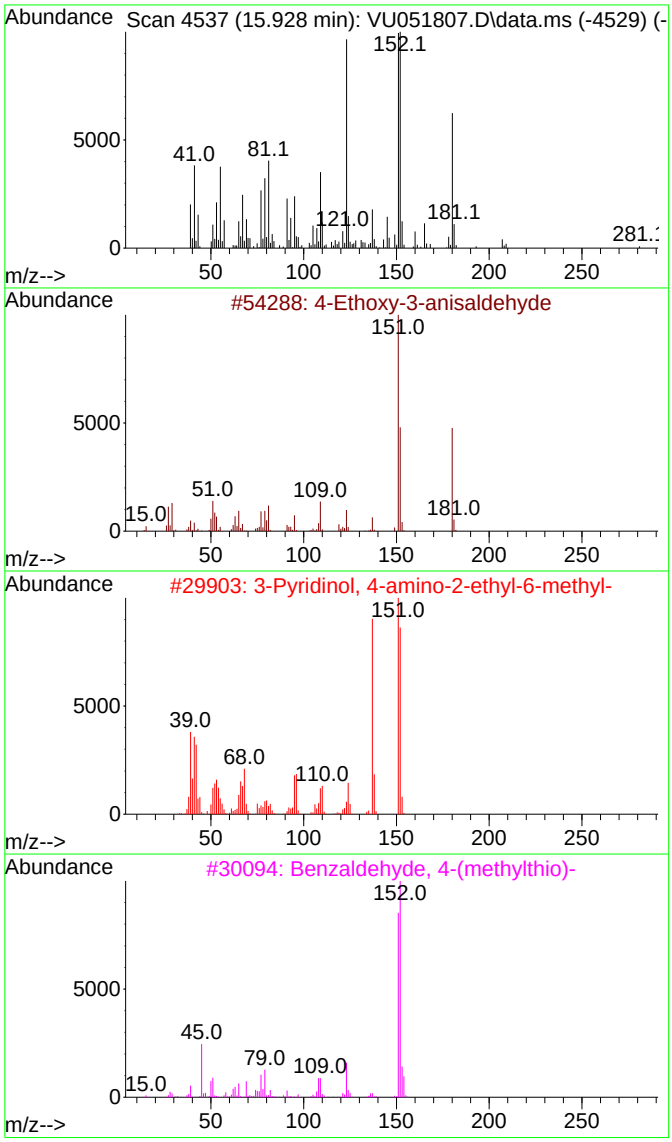
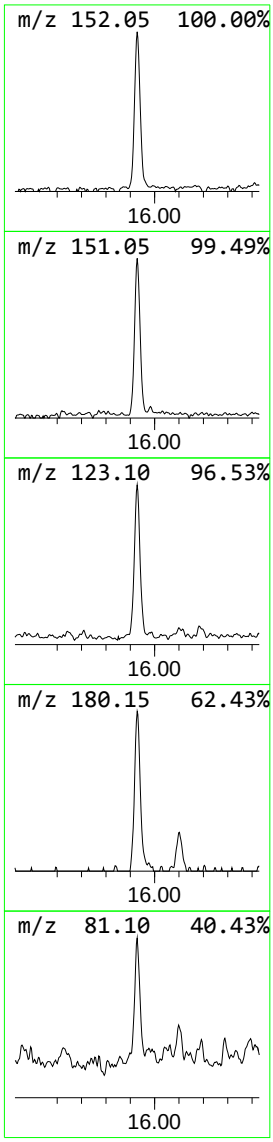
Peak Number 11 unknown-02 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.928	1.11 ug/L	119276	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Ethoxy-3-anisaldehyde	180	C10H12O3	000120-25-2	49
2	3-Pyridinol, 4-amino-2-ethyl-6-m...	152	C8H12N2O	1010351-35-8	46
3	Benzaldehyde, 4-(methylthio)-	152	C8H8OS	003446-89-7	46
4	2-Amino-4-acetamino anisole	180	C9H12N2O2	006375-47-9	46
5	Vanillin	152	C8H8O3	000121-33-5	46





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU111022\
Data File : VU051807.D
Acq On : 10 Nov 2022 17:51
Operator : JC/MD
Sample : N5571-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GBMW1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102822WMA.M
Quant Title : TRACE VOA SFAM1.0

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

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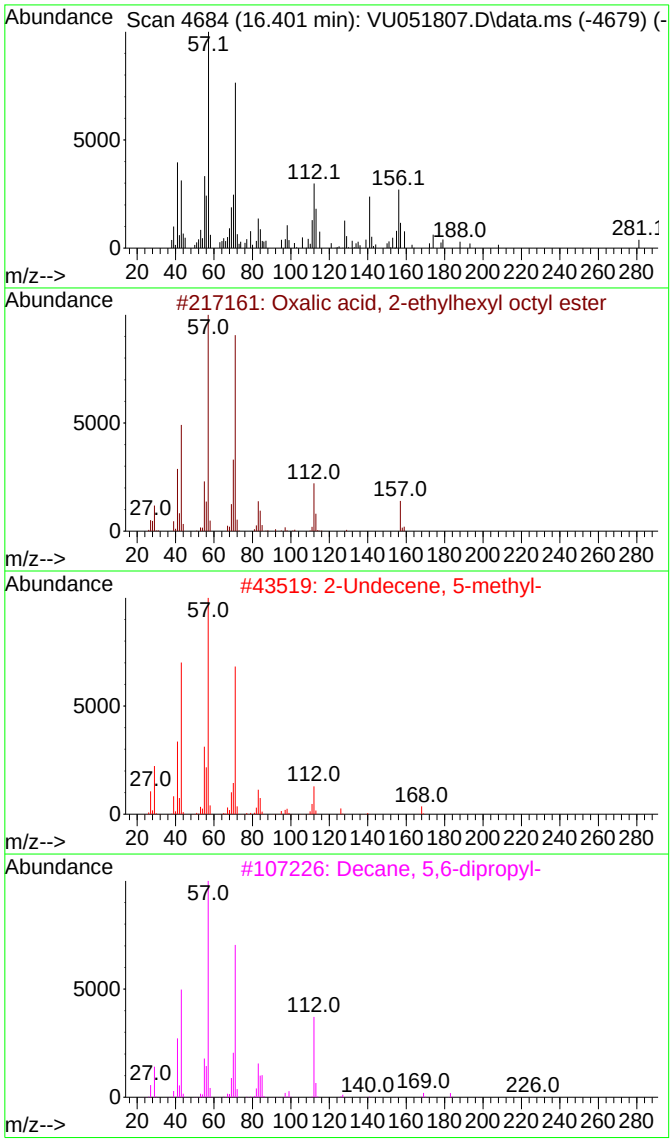
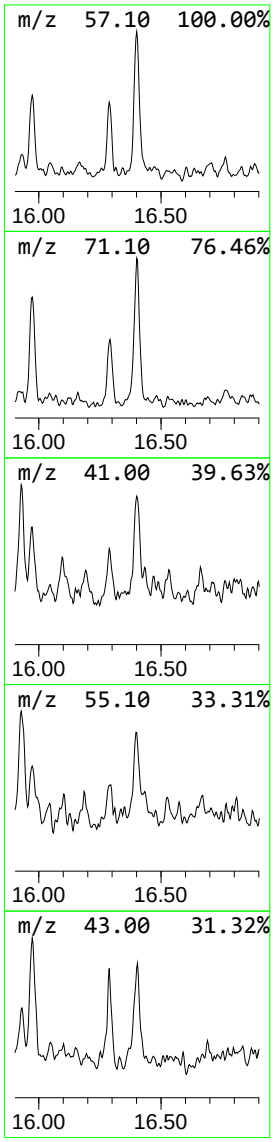
Peak Number 12 unknown-03 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.401	0.51 ug/L	54660	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxalic acid, 2-ethylhexyl octyl ...	314	C18H34O4	1000309-39-1	43
2	2-Undecene, 5-methyl-	168	C12H24	056851-34-4	38
3	Decane, 5,6-dipropyl-	226	C16H34	119209-20-0	38
4	Oxalic acid, 2-ethylhexyl isobut...	258	C14H26O4	1010309-38-5	38
5	Methoxyacetic acid, 2-octyl ester	202	C11H22O3	1000282-69-6	38





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU111022\
Data File : VU051807.D
Acq On : 10 Nov 2022 17:51
Operator : JC/MD
Sample : N5571-01
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GBMW1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102822WMA.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
Butanal	4.572	0.9	ug/L	81099	1	6.353	456970	5.0
1-Hexanol, 2-et...	12.060	0.5	ug/L	57672	3	11.816	535759	5.0
2,2,5,5,6-Penta...	13.970	0.9	ug/L	97319	3	11.816	535759	5.0
Cyclohexene, 1-...	14.012	0.7	ug/L	77407	3	11.816	535759	5.0
Cyclohexane, 1-...	14.147	0.9	ug/L	100234	3	11.816	535759	5.0
unknown-01	14.320	1.3	ug/L	139814	3	11.816	535759	5.0
1,5-Dimethyl-6-...	14.369	1.6	ug/L	172604	3	11.816	535759	5.0
Cyclopentene, 1...	14.825	0.6	ug/L	63902	3	11.816	535759	5.0
(DEL) Alkane: S...	15.455	0.7	ug/L	73894	3	11.816	535759	5.0
unknown-02	15.928	1.1	ug/L	119276	3	11.816	535759	5.0
unknown-03	16.401	0.5	ug/L	54660	3	11.816	535759	5.0