LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV100322\
Data File : VV028286.D
Acq On : 03 Oct 2022 18:23
Operator : SY/MD
Sample : N4856-11DL 40X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
F5H36DL

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\SFAMVTR100322WMA.M
Title : TRACE VOA SFAM1.0

Signal : TIC: VV028286.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.301	72	81	90	rBV	303063	312271	4.71%	1.767%
2	1.407	100	114	128	rBV	59314	214702	3.24%	1.215%
3	1.561	156	162	173	rVB	99439	118910	1.79%	0.673%
4	1.629	175	183	199	rVB	560815	688133	10.39%	3.895%
5	1.802	230	237	249	rVB3	93969	143681	2.17%	0.813%
6	2.018	296	304	319	rVV	141566	202571	3.06%	1.147%
7	2.101	321	330	349	rVB	205915	339022	5.12%	1.919%
8	2.490	435	451	469	rBV5	151644	430907	6.50%	2.439%
9	2.571	469	476	497	rVB3	64148	125688	1.90%	0.711%
10	2.754	519	533	548	rBV	61964	125807	1.90%	0.712%
11	3.754	830	844	861	rBV2	69914	171351	2.59%	0.970%
12	3.902	878	890	919	rBV3	65921	269430	4.07%	1.525%
13	4.342	1009	1027	1052	rBV	126489	328763	4.96%	1.861%
14	4.670	1111	1129	1142	rBV5	37617	97682	1.47%	0.553%
15	4.757	1142	1156	1177	rVB3	49559	130589	1.97%	0.739%
16	5.043	1228	1245	1248	rBV2	318500	604145	9.12%	3.419%
17	5.092	1249	1260	1290	rVV	3037114	6624836	100.00%	37.496%
18	5.227	1292	1302	1322	rVB2	56353	160509	2.42%	0.908%
19	5.612	1410	1422	1459	rBV	228471	508739	7.68%	2.879%
20	6.063	1548	1562	1592	rBV	163509	389620	5.88%	2.205%
21	6.770	1769	1782	1799	rBV4	41601	109237	1.65%	0.618%
22	6.982	1830	1848	1869	rBV3	67038	136221	2.06%	0.771%
23	7.310	1940	1950	1965	rBV	339382	595209	8.98%	3.369%
24	7.387	1965	1974	1996	rVB	91434	170280	2.57%	0.964%
25	7.622	2039	2047	2065	rVB	37227	70214	1.06%	0.397%
26	8.085	2181	2191	2224	rBV	311801	622775	9.40%	3.525%
27	8.847	2416	2428	2454	rBV	335642	568966	8.59%	3.220%
28	9.008	2467	2478	2500	rBV	598911	936022	14.13%	5.298%
29	9.136	2507	2518	2543	rVB	302929	511668	7.72%	2.896%
30	10.214	2842	2853	2870	rBV	123806	198330	2.99%	1.123%
31	10.352	2885	2896	2915	rBV	59482	97096	1.47%	0.550%
32	10.445	2915	2925	2931	rBV2	54914	96344	1.45%	0.545%
33	10.911	3059	3070	3087	rBV	160974	243900	3.68%	1.380%
34	11.246	3163	3174	3193	rBV	367943	573686	8.66%	3.247%
35	11.525	3249	3261	3280	rBV2	88104	180923	2.73%	1.024%
36	11.622	3280	3291	3310	rVB	345587	569777	8.60%	3.225%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV100322\
Data File : VV028286.D
Acq On : 03 Oct 2022 18:23
Operator : SY/MD
Sample : N4856-11DL 40X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
F5H36DL

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\SFAMVTR100322WMA.M
Title : TRACE VOA SFAM1.0

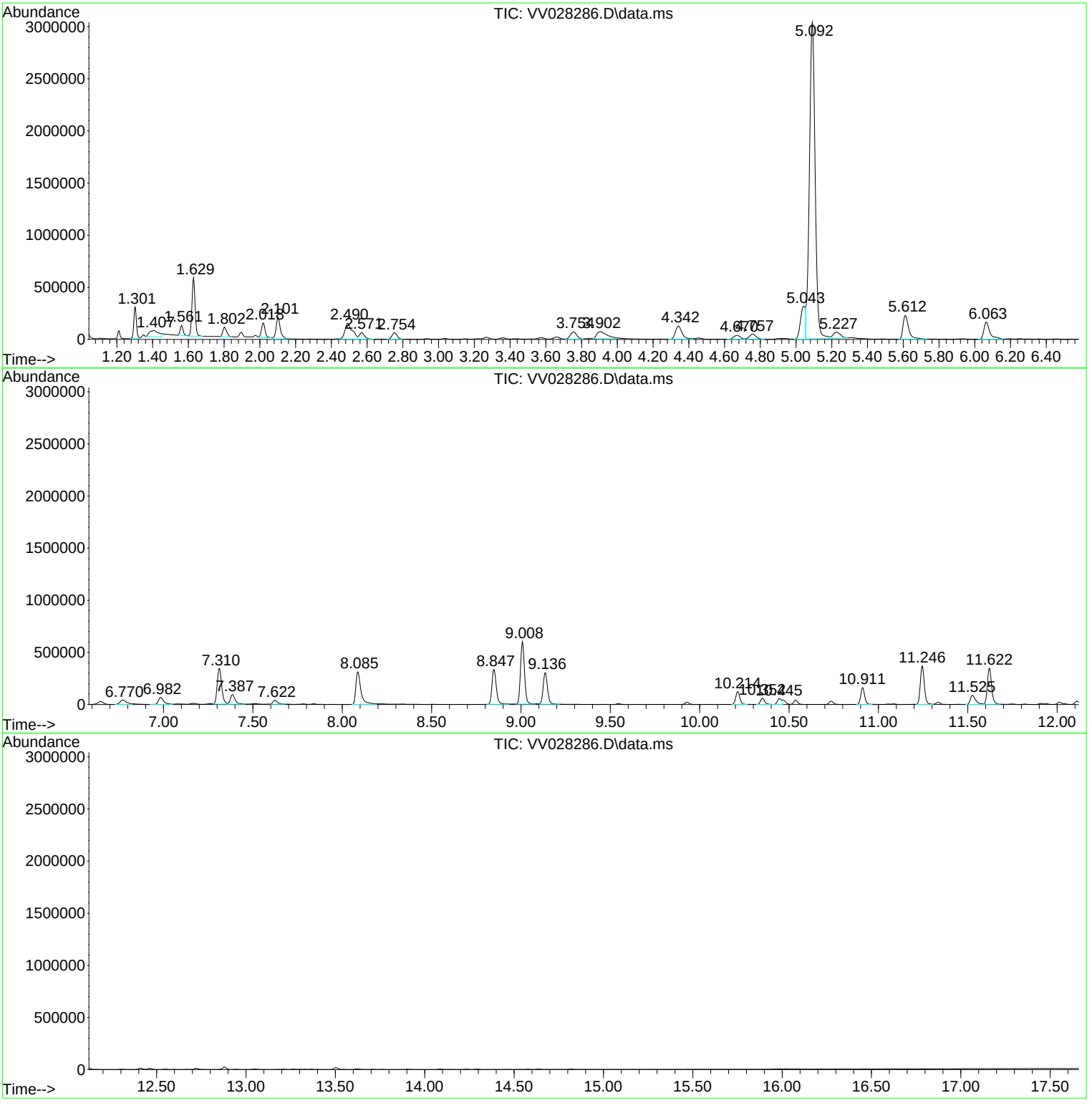
Sum of corrected areas: 17668004

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV100322\
Data File : VV028286.D
Acq On : 03 Oct 2022 18:23
Operator : SY/MD
Sample : N4856-11DL 40X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
F5H36DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR100322WMA.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\VV100322\
Data File : VV028286.D
Acq On : 03 Oct 2022 18:23
Operator : SY/MD
Sample : N4856-11DL 40X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
F5H36DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR100322WMA.M
Quant Title : TRACE VOA SFAM1.0

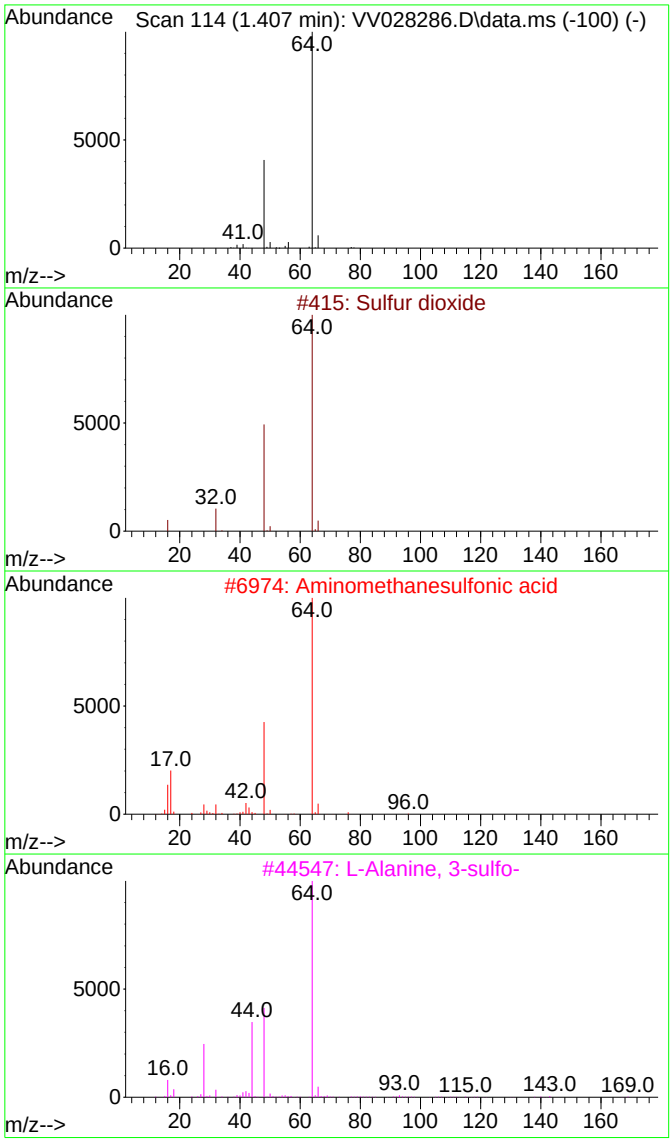
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

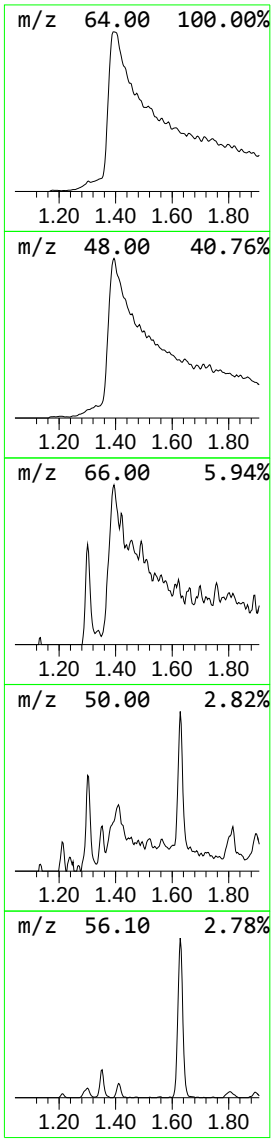
****

Peak Number 1 Sulfur dioxide Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.407	2.11 ug/L	214702	1,4-Difluorobenzene	5.612

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Sulfur dioxide	64	O2S	007446-09-5	90
2	Aminomethanesulfonic acid	111	CH5NO3S	013881-91-9	74
3	L-Alanine, 3-sulfo-	169	C3H7NO5S	000498-40-8	9
4	Aminomethanesulfonic acid	111	CH5NO3S	013881-91-9	9
5	Sulfur dioxide	64	O2S	007446-09-5	7





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV100322\
Data File : VV028286.D
Acq On : 03 Oct 2022 18:23
Operator : SY/MD
Sample : N4856-11DL 40X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
F5H36DL

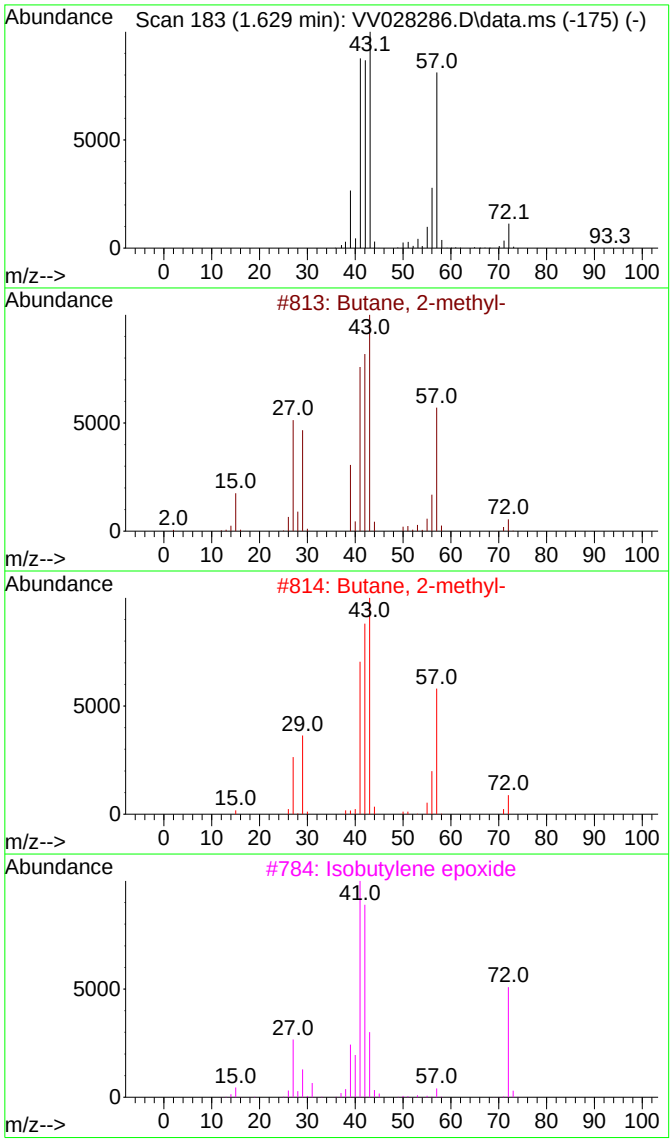
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR100322WMA.M
Quant Title : TRACE VOA SFAM1.0

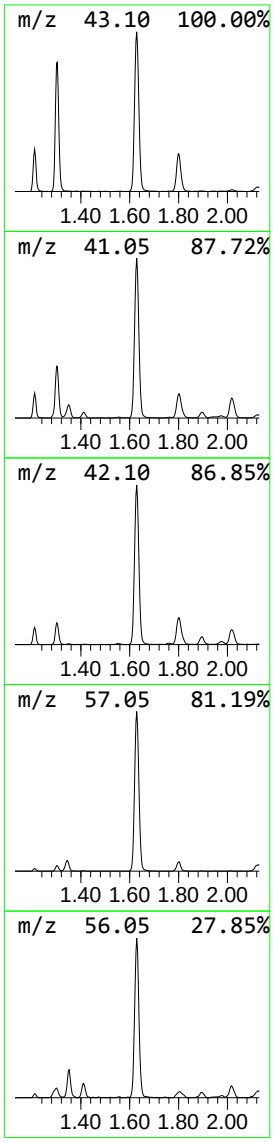
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.629	6.76 ug/L	688133	1,4-Difluorobenzene	5.612

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	86
2	Butane, 2-methyl-	72	C5H12	000078-78-4	72
3	Isobutylene epoxide	72	C4H8O	000558-30-5	38
4	Pentane	72	C5H12	000109-66-0	33
5	Pentane	72	C5H12	000109-66-0	28





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV100322\
Data File : VV028286.D
Acq On : 03 Oct 2022 18:23
Operator : SY/MD
Sample : N4856-11DL 40X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
F5H36DL

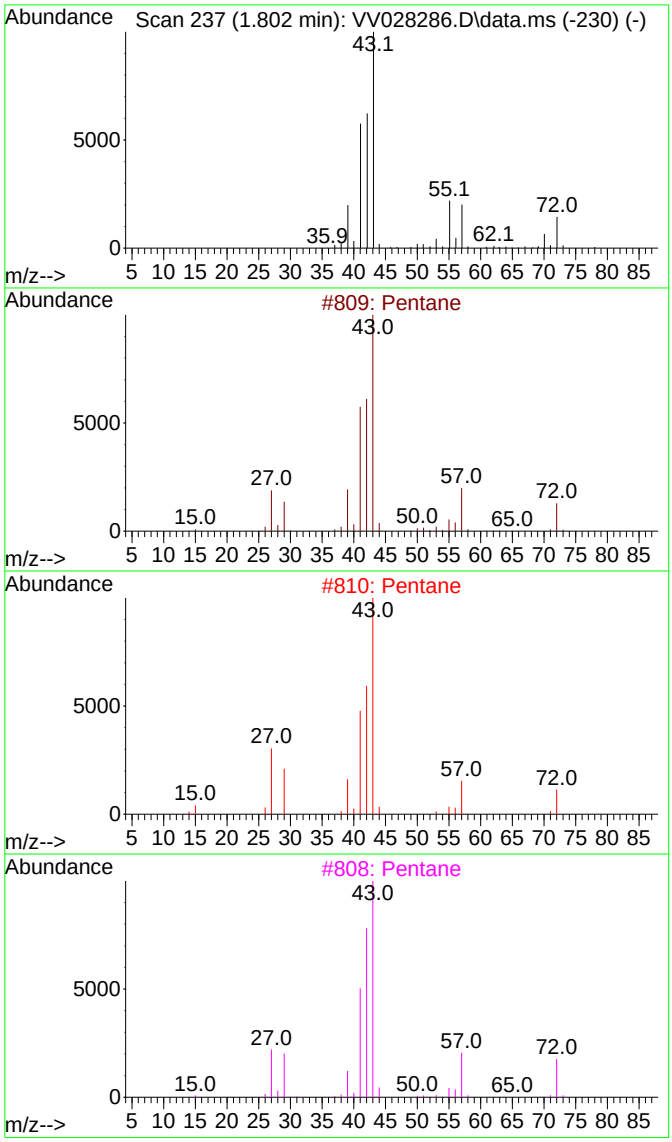
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR100322WMA.M
Quant Title : TRACE VOA SFAM1.0

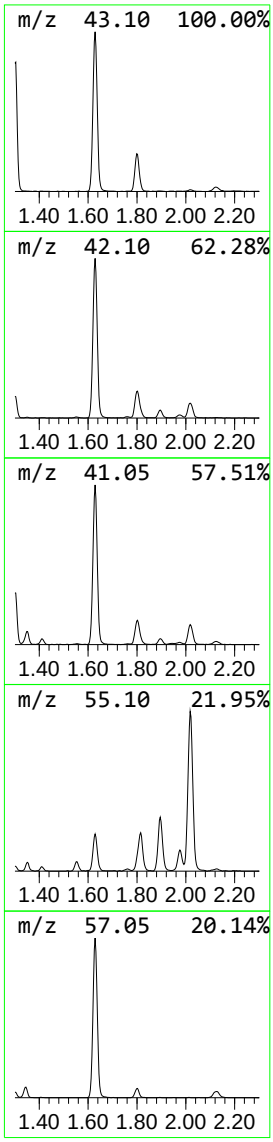
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.802	1.41 ug/L	143681	1,4-Difluorobenzene	5.612

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane	72	C5H12	000109-66-0	87
2	Pentane	72	C5H12	000109-66-0	80
3	Pentane	72	C5H12	000109-66-0	80
4	Pentane	72	C5H12	000109-66-0	78
5	Pentane	72	C5H12	000109-66-0	72





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV100322\
Data File : VV028286.D
Acq On : 03 Oct 2022 18:23
Operator : SY/MD
Sample : N4856-11DL 40X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
F5H36DL

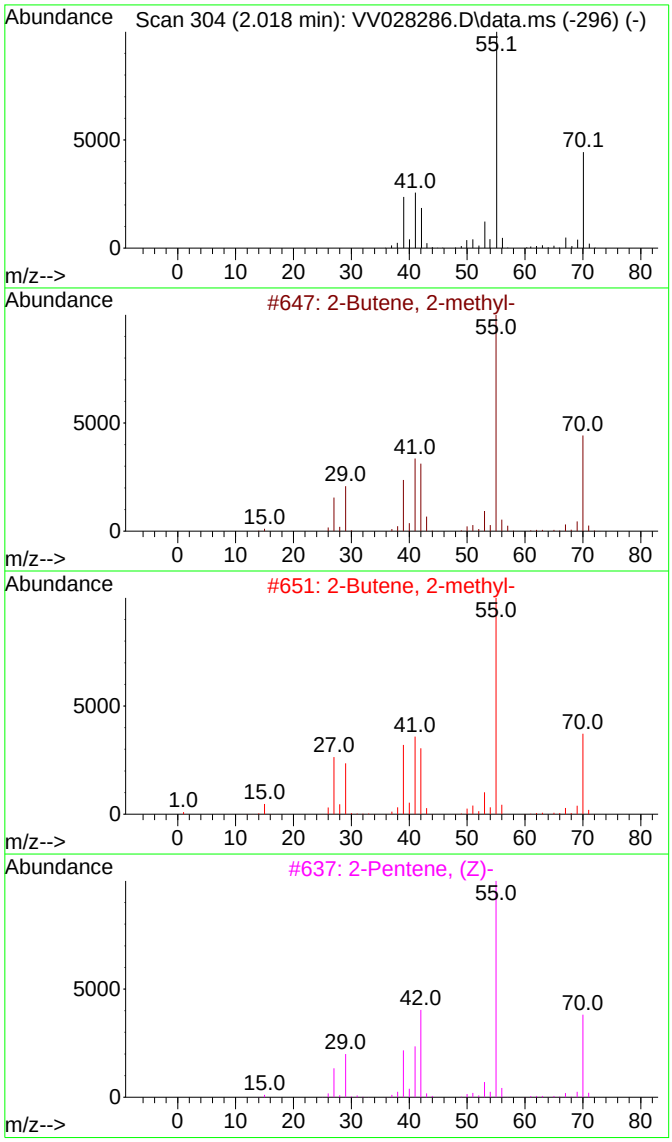
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR100322WMA.M
Quant Title : TRACE VOA SFAM1.0

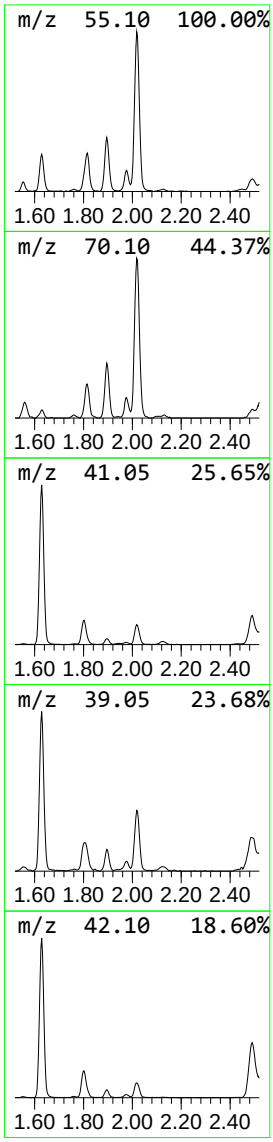
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.018	1.99 ug/L	202571	1,4-Difluorobenzene	5.612

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butene, 2-methyl-	70	C5H10	000513-35-9	90
2	2-Butene, 2-methyl-	70	C5H10	000513-35-9	90
3	2-Pentene, (Z)-	70	C5H10	000627-20-3	83
4	2-Methyl-1-butene	70	C5H10	000563-46-2	83
5	2-Pentene	70	C5H10	000109-68-2	83





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV100322\
Data File : VV028286.D
Acq On : 03 Oct 2022 18:23
Operator : SY/MD
Sample : N4856-11DL 40X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
F5H36DL

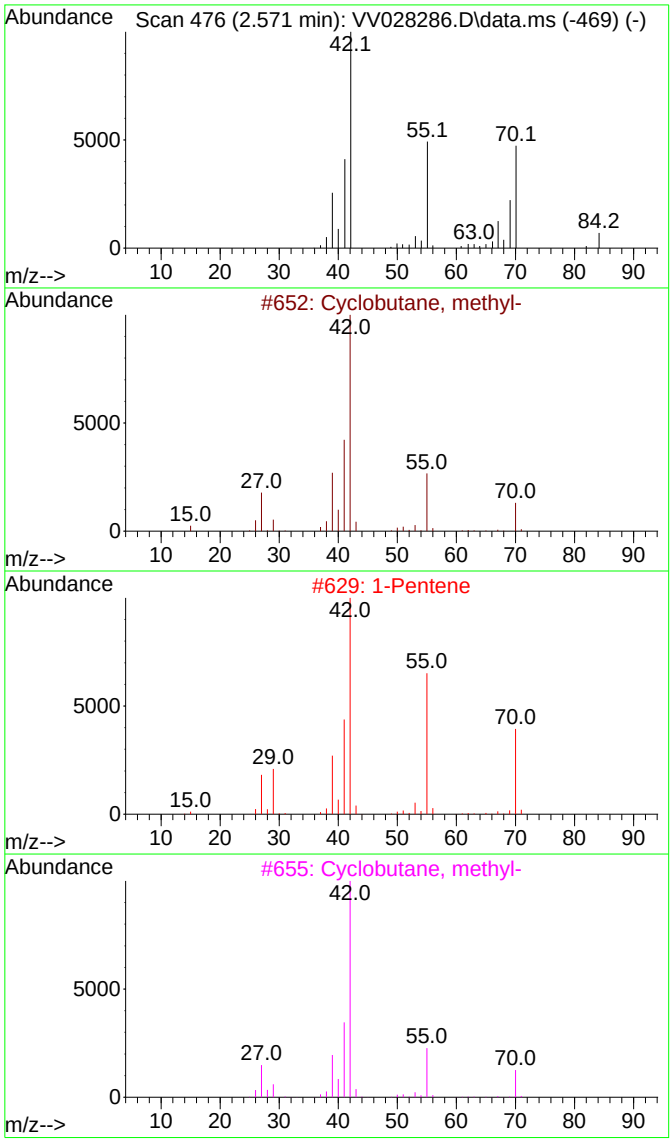
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR100322WMA.M
Quant Title : TRACE VOA SFAM1.0

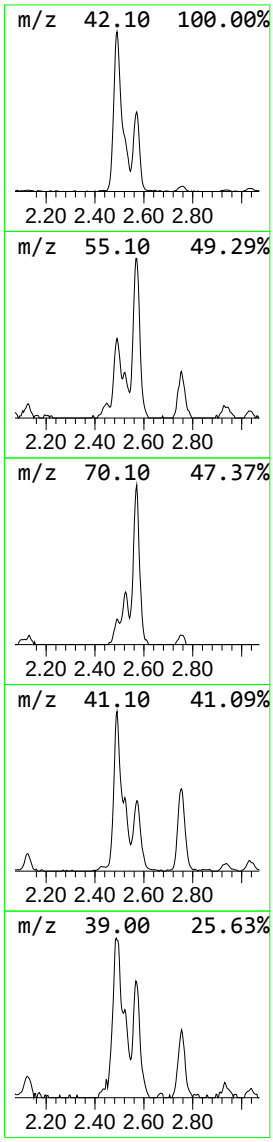
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.571	1.24 ug/L	125688	1,4-Difluorobenzene	5.612

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclobutane, methyl-	70	C5H10	000598-61-8	64
2	1-Pentene	70	C5H10	000109-67-1	64
3	Cyclobutane, methyl-	70	C5H10	000598-61-8	59
4	Cyclopropane, ethyl-	70	C5H10	001191-96-4	47
5	1-Pentene	70	C5H10	000109-67-1	45





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV100322\
Data File : VV028286.D
Acq On : 03 Oct 2022 18:23
Operator : SY/MD
Sample : N4856-11DL 40X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
F5H36DL

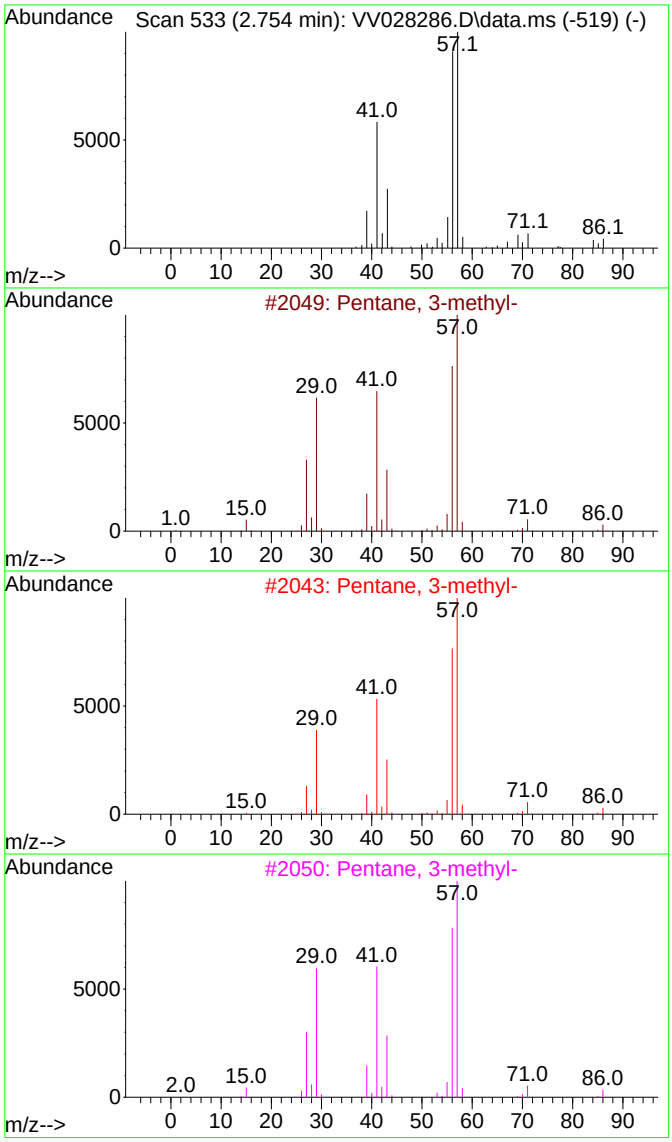
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR100322WMA.M
Quant Title : TRACE VOA SFAM1.0

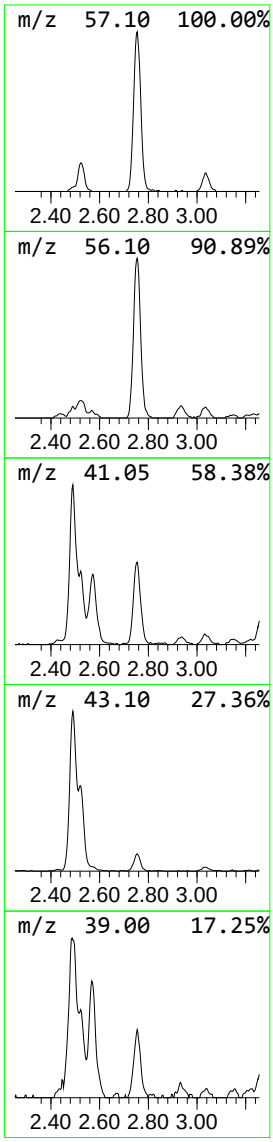
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.754	1.24 ug/L	125807	1,4-Difluorobenzene	5.612

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	83
2	Pentane, 3-methyl-	86	C6H14	000096-14-0	80
3	Pentane, 3-methyl-	86	C6H14	000096-14-0	80
4	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	72
5	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV100322\
Data File : VV028286.D
Acq On : 03 Oct 2022 18:23
Operator : SY/MD
Sample : N4856-11DL 40X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
F5H36DL

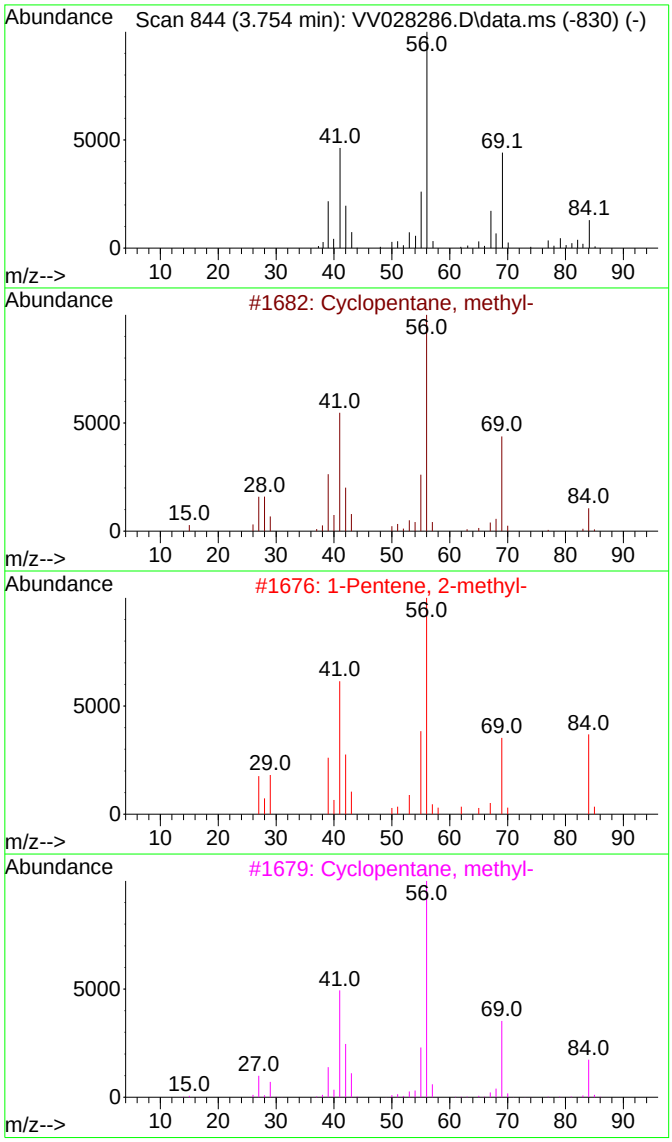
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR100322WMA.M
Quant Title : TRACE VOA SFAM1.0

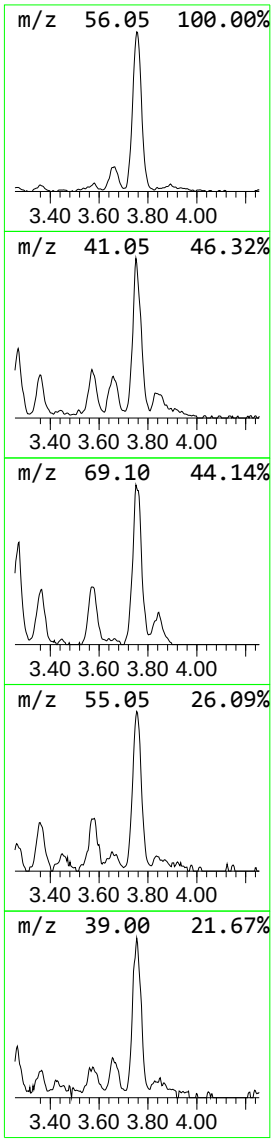
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 (DEL) Alkane: Cyclic3.754 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.754	1.68 ug/L	171351	1,4-Difluorobenzene	5.612

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	81
2	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	64
3	Cyclopentane, methyl-	84	C6H12	000096-37-7	64
4	Cyclopentane, methyl-	84	C6H12	000096-37-7	64
5	Cyclopentane, methyl-	84	C6H12	000096-37-7	59





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV100322\
Data File : VV028286.D
Acq On : 03 Oct 2022 18:23
Operator : SY/MD
Sample : N4856-11DL 40X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
F5H36DL

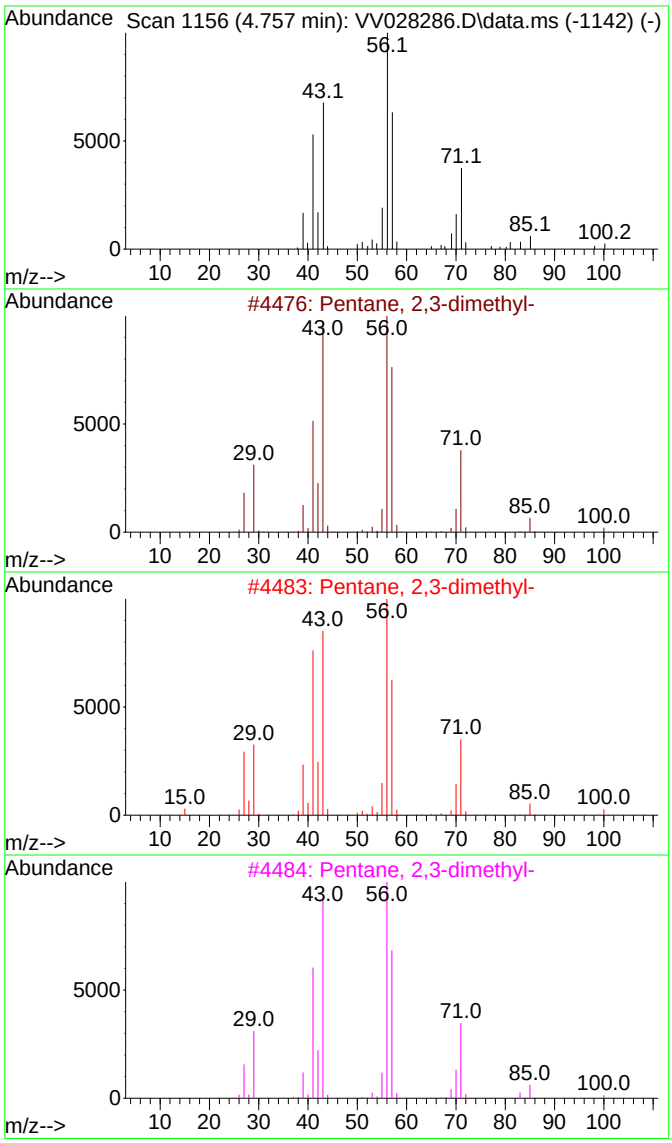
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR100322WMA.M
Quant Title : TRACE VOA SFAM1.0

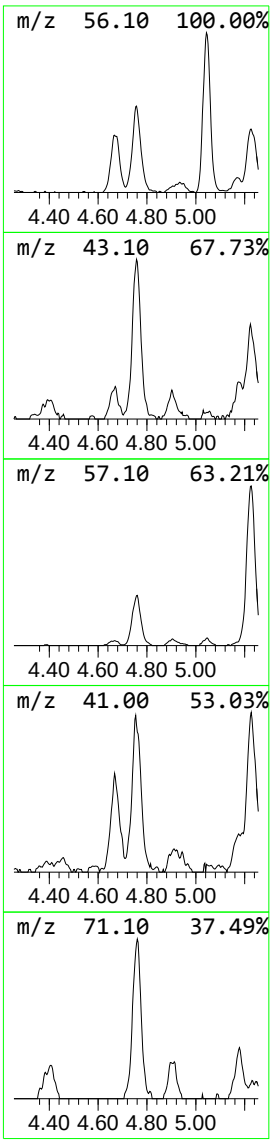
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.757	1.28 ug/L	130589	1,4-Difluorobenzene	5.612

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
2	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	90
3	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	86
4	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	83
5	Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	56





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV100322\
Data File : VV028286.D
Acq On : 03 Oct 2022 18:23
Operator : SY/MD
Sample : N4856-11DL 40X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
F5H36DL

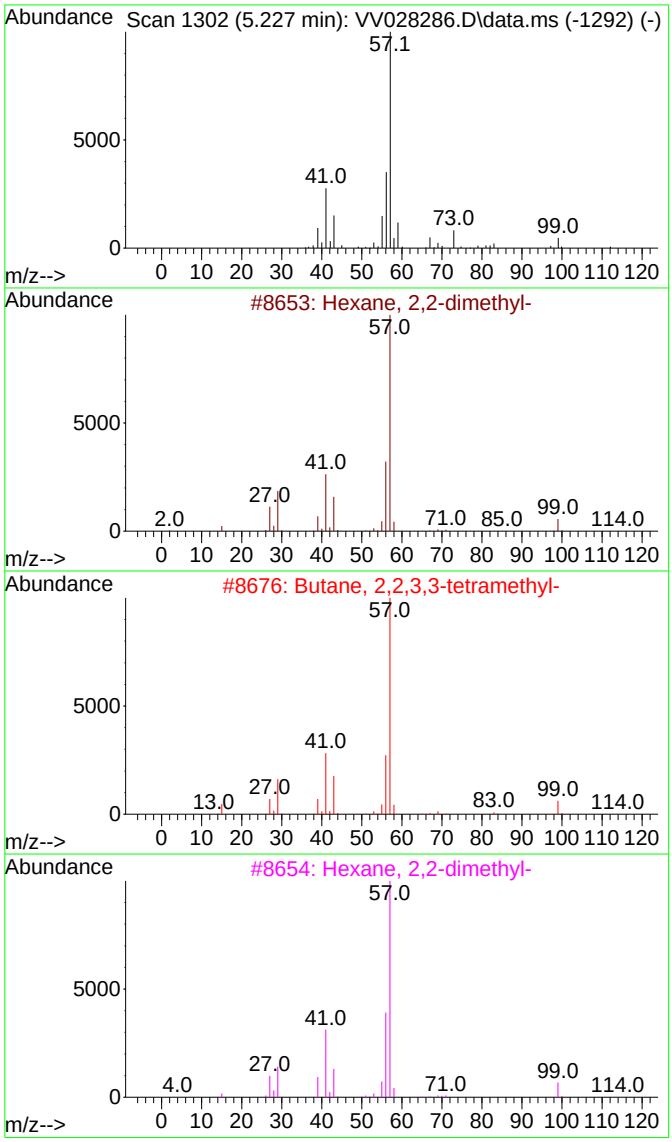
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR100322WMA.M
Quant Title : TRACE VOA SFAM1.0

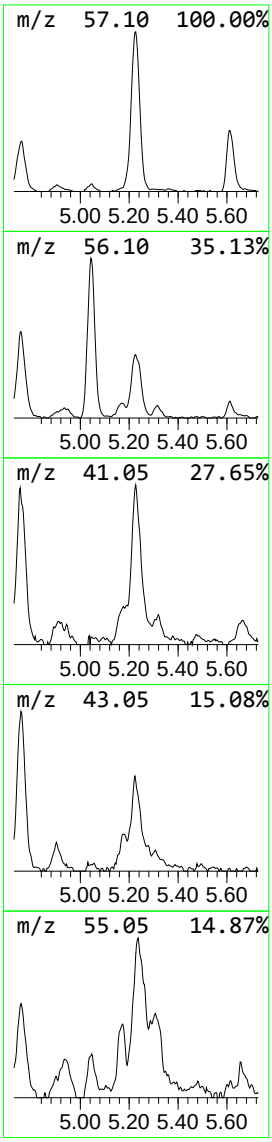
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.227	1.58 ug/L	160509	1,4-Difluorobenzene	5.612

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	50
2	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	50
3	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	50
4	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	40
5	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	40





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV100322\
Data File : VV028286.D
Acq On : 03 Oct 2022 18:23
Operator : SY/MD
Sample : N4856-11DL 40X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
F5H36DL

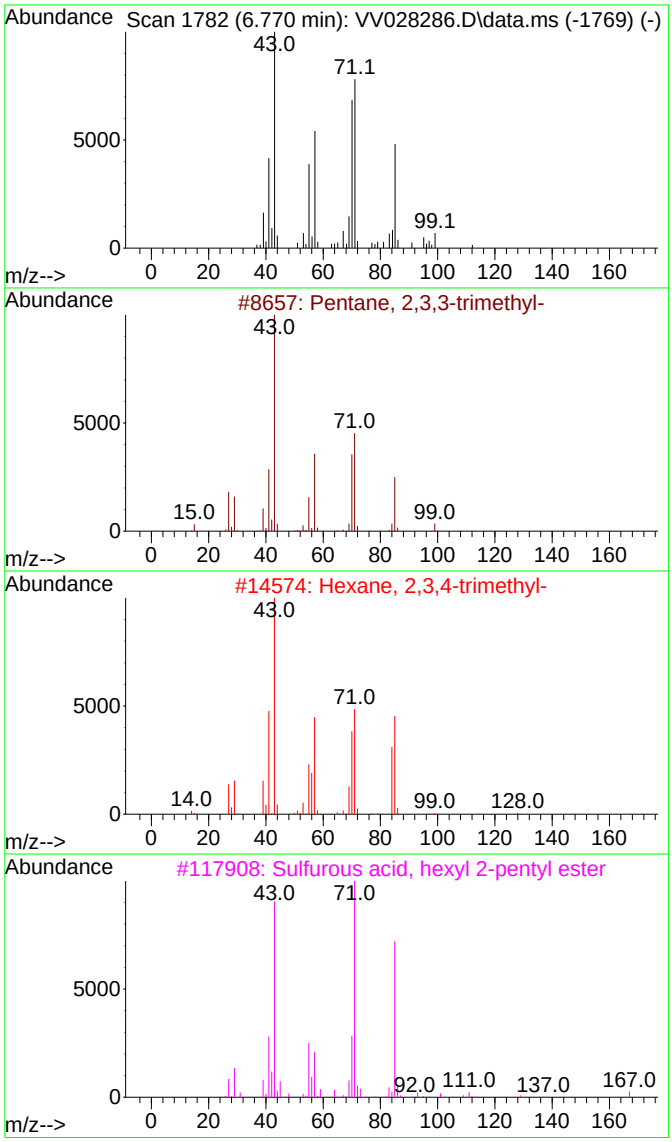
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR100322WMA.M
Quant Title : TRACE VOA SFAM1.0

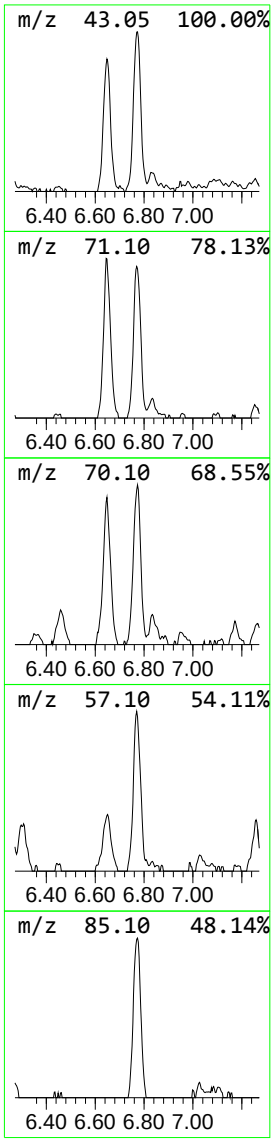
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.770	1.07 ug/L	109237	1,4-Difluorobenzene	5.612

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	72
2	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	64
3	Sulfurous acid, hexyl 2-pentyl e...	236	C11H24O3S	1000309-15-6	59
4	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	59
5	Hexane, 3-methyl-	100	C7H16	000589-34-4	53





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV100322\
Data File : VV028286.D
Acq On : 03 Oct 2022 18:23
Operator : SY/MD
Sample : N4856-11DL 40X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
F5H36DL

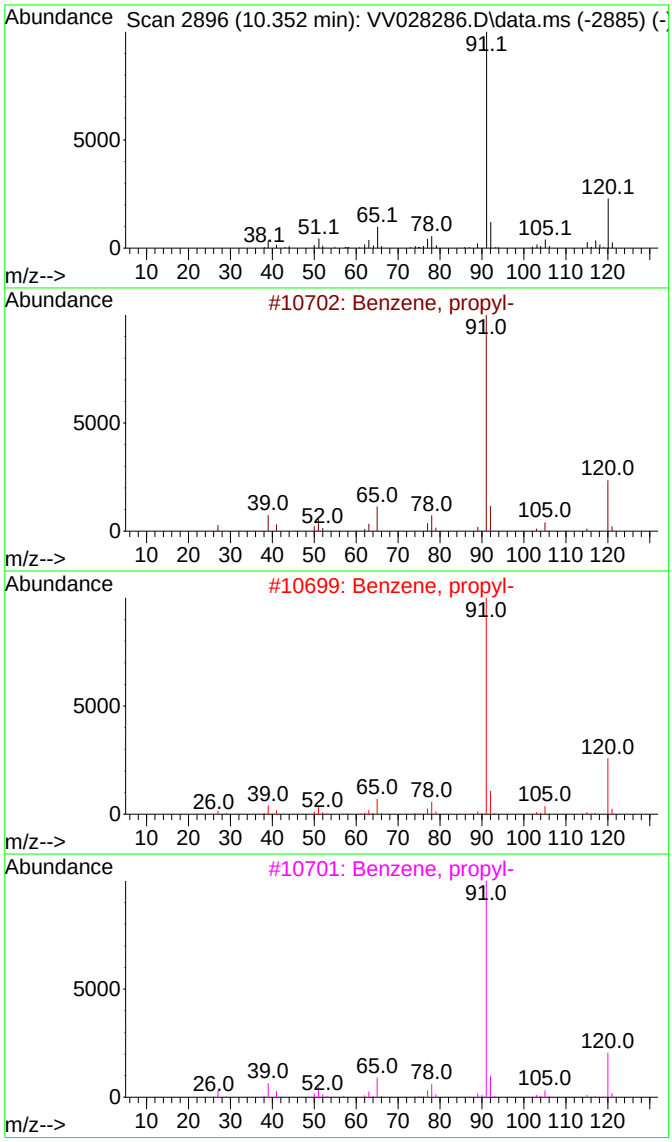
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR100322WMA.M
Quant Title : TRACE VOA SFAM1.0

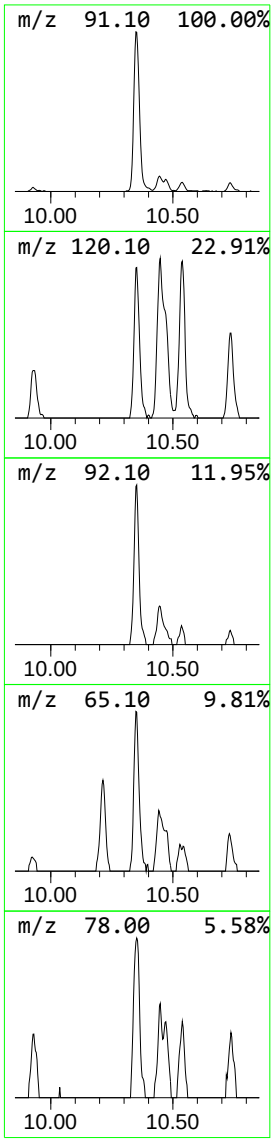
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Benzene, propyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.352	0.85 ug/L	97096	1,4-Dichlorobenzene-d4	11.246

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	90
2	Benzene, propyl-	120	C9H12	000103-65-1	87
3	Benzene, propyl-	120	C9H12	000103-65-1	87
4	Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	72
5	N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV100322\
Data File : VV028286.D
Acq On : 03 Oct 2022 18:23
Operator : SY/MD
Sample : N4856-11DL 40X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
F5H36DL

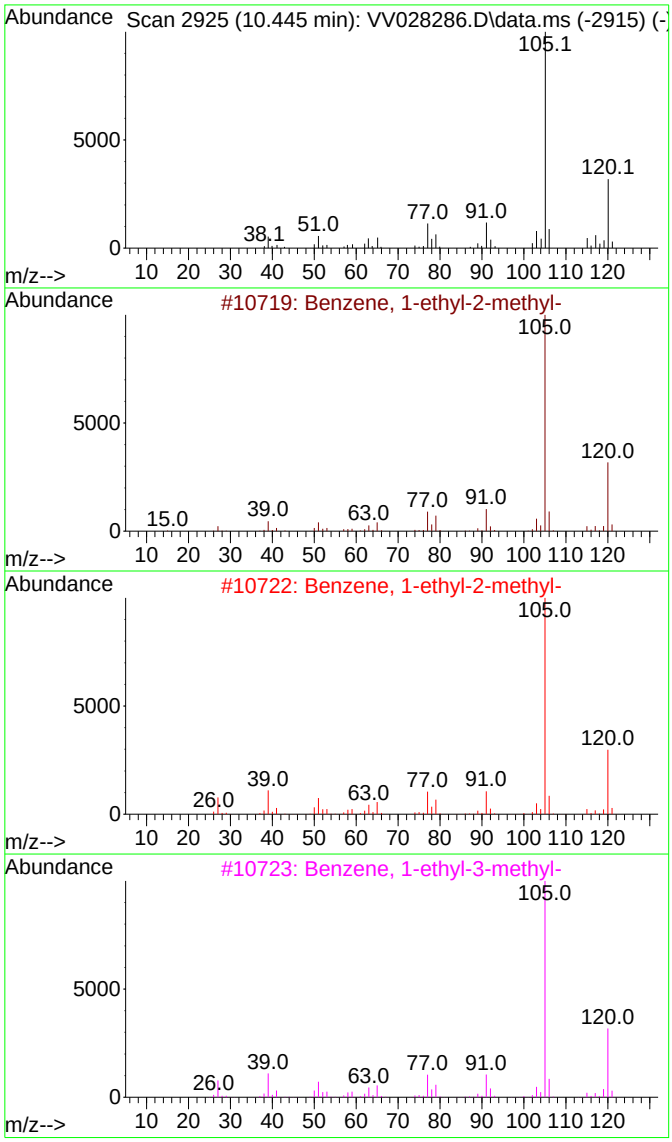
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR100322WMA.M
Quant Title : TRACE VOA SFAM1.0

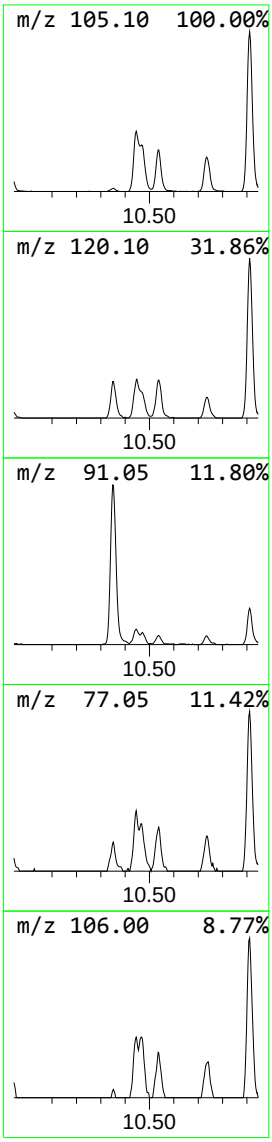
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Benzene, 1-ethyl-2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.445	0.84 ug/L	96344	1,4-Dichlorobenzene-d4	11.246

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV100322\
Data File : VV028286.D
Acq On : 03 Oct 2022 18:23
Operator : SY/MD
Sample : N4856-11DL 40X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
F5H36DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR100322WMA.M
Quant Title : TRACE VOA SFAM1.0

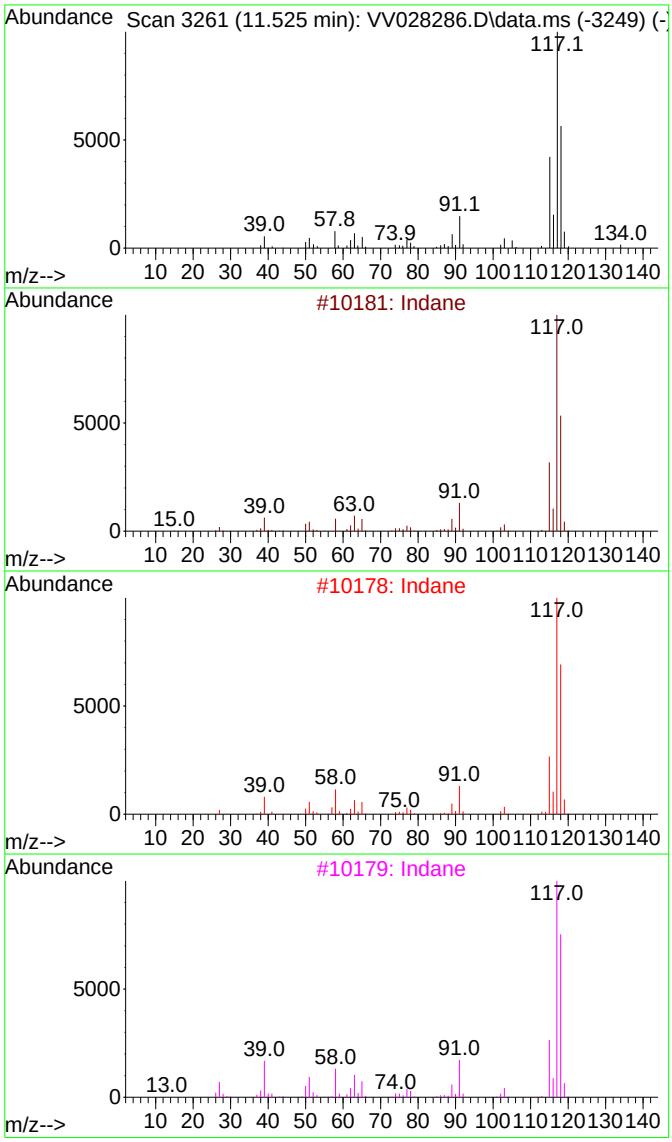
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

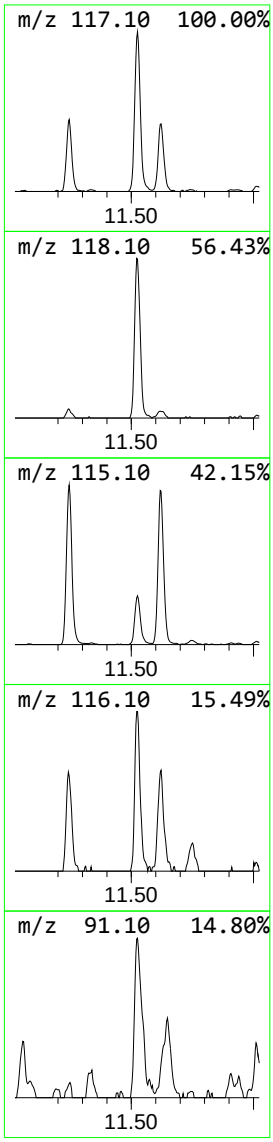
****

Peak Number 14 Indane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.525	1.58 ug/L	180923	1,4-Dichlorobenzene-d4	11.246

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	93
2	Indane	118	C9H10	000496-11-7	93
3	Indane	118	C9H10	000496-11-7	76
4	Benzene, 1-propenyl-	118	C9H10	000637-50-3	76
5	Benzene, cyclopropyl-	118	C9H10	000873-49-4	70





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV100322\
Data File : VV028286.D
Acq On : 03 Oct 2022 18:23
Operator : SY/MD
Sample : N4856-11DL 40X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
F5H36DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR100322WMA.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
Sulfur dioxide	1.407	2.1	ug/L	214702	1	5.612	508739	5.0
(DEL) Alkane: S...	1.629	6.8	ug/L	688133	1	5.612	508739	5.0
(DEL) Alkane: S...	1.802	1.4	ug/L	143681	1	5.612	508739	5.0
(DEL) Alkane: S...	2.018	2.0	ug/L	202571	1	5.612	508739	5.0
(DEL) Alkane: S...	2.571	1.2	ug/L	125688	1	5.612	508739	5.0
(DEL) Alkane: S...	2.754	1.2	ug/L	125807	1	5.612	508739	5.0
(DEL) Alkane: C...	3.754	1.7	ug/L	171351	1	5.612	508739	5.0
(DEL) Alkane: S...	4.757	1.3	ug/L	130589	1	5.612	508739	5.0
(DEL) Alkane: S...	5.227	1.6	ug/L	160509	1	5.612	508739	5.0
(DEL) Alkane: S...	6.770	1.1	ug/L	109237	1	5.612	508739	5.0
Benzene, propyl-	10.352	0.8	ug/L	97096	3	11.246	573686	5.0
Benzene, 1-ethy...	10.445	0.8	ug/L	96344	3	11.246	573686	5.0
Indane	11.525	1.6	ug/L	180923	3	11.246	573686	5.0