

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041222\
 Data File : VV025482.D
 Acq On : 12 Apr 2022 18:16
 Operator : SY/MD
 Sample : N2356-08
 Misc : 25mL/MSVOA_V/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 GBDS6

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\SFAMVTR040822WMA.M
 Title : TRACE VOA SFAM1.0

Signal : TIC: VV025482.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.217	48	55	70	rBV	543257	487983	1.48%	0.840%
2	1.310	75	84	94	rBV	3724028	3528877	10.68%	6.076%
3	1.638	175	186	209	rVB	7084960	8625541	26.11%	14.852%
4	1.809	230	239	254	rBV	300078	393829	1.19%	0.678%
5	2.027	299	307	323	rVB	235269	331713	1.00%	0.571%
6	2.111	324	333	356	rVB2	162000	391941	1.19%	0.675%
7	2.500	423	454	461	rBV2	567323	1243640	3.76%	2.141%
8	2.580	471	479	499	rVB	490179	916011	2.77%	1.577%
9	2.764	519	536	561	rBV	393506	776974	2.35%	1.338%
10	3.767	832	848	866	rVV	378900	935932	2.83%	1.612%
11	4.680	1112	1132	1146	rBV	370296	914456	2.77%	1.575%
12	4.767	1146	1159	1188	rVB	225255	589689	1.79%	1.015%
13	5.101	1233	1263	1282	rBV	15557280	33034105	100.00%	56.880%
14	5.233	1294	1304	1322	rVB5	279830	771238	2.33%	1.328%
15	5.619	1413	1424	1462	rBV	156004	367043	1.11%	0.632%
16	6.654	1732	1746	1764	rVB	176737	367497	1.11%	0.633%
17	6.776	1769	1784	1802	rBV2	316805	755702	2.29%	1.301%
18	7.317	1942	1952	1966	rVB	209794	359047	1.09%	0.618%
19	8.088	2180	2192	2211	rBV	172412	340779	1.03%	0.587%
20	8.853	2405	2430	2442	rBV	244787	414849	1.26%	0.714%
21	11.246	3164	3174	3192	rBV	266631	419898	1.27%	0.723%
22	11.525	3248	3261	3282	rBV	976261	1611338	4.88%	2.774%
23	11.625	3282	3292	3313	rVB2	294168	499015	1.51%	0.859%

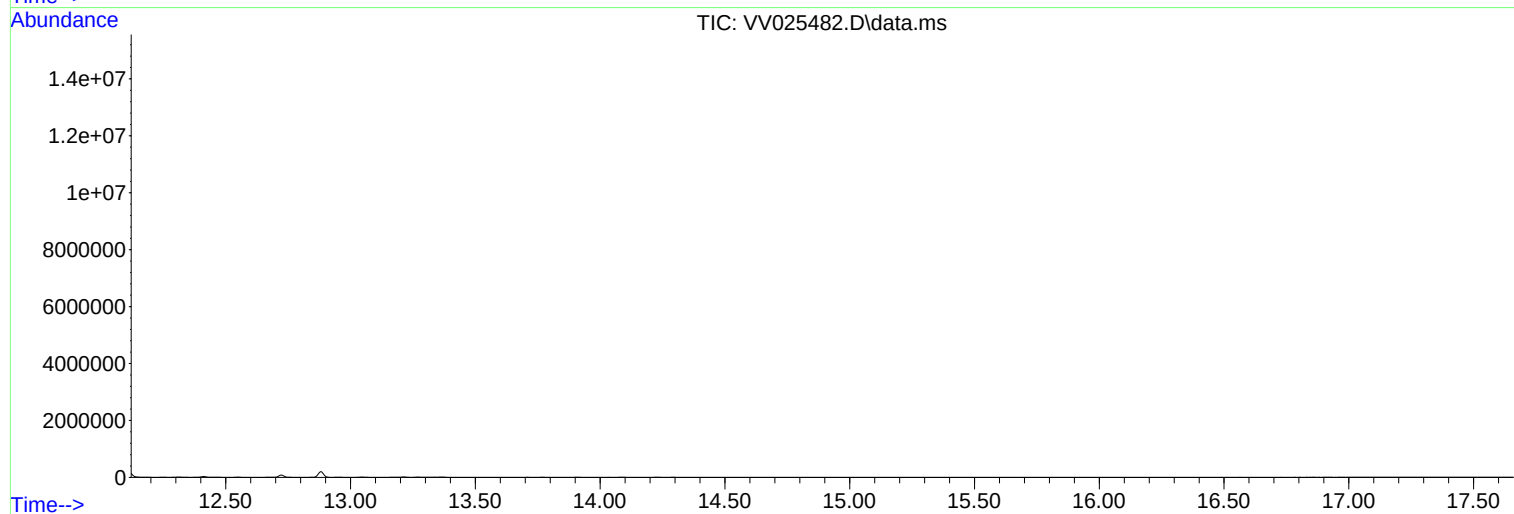
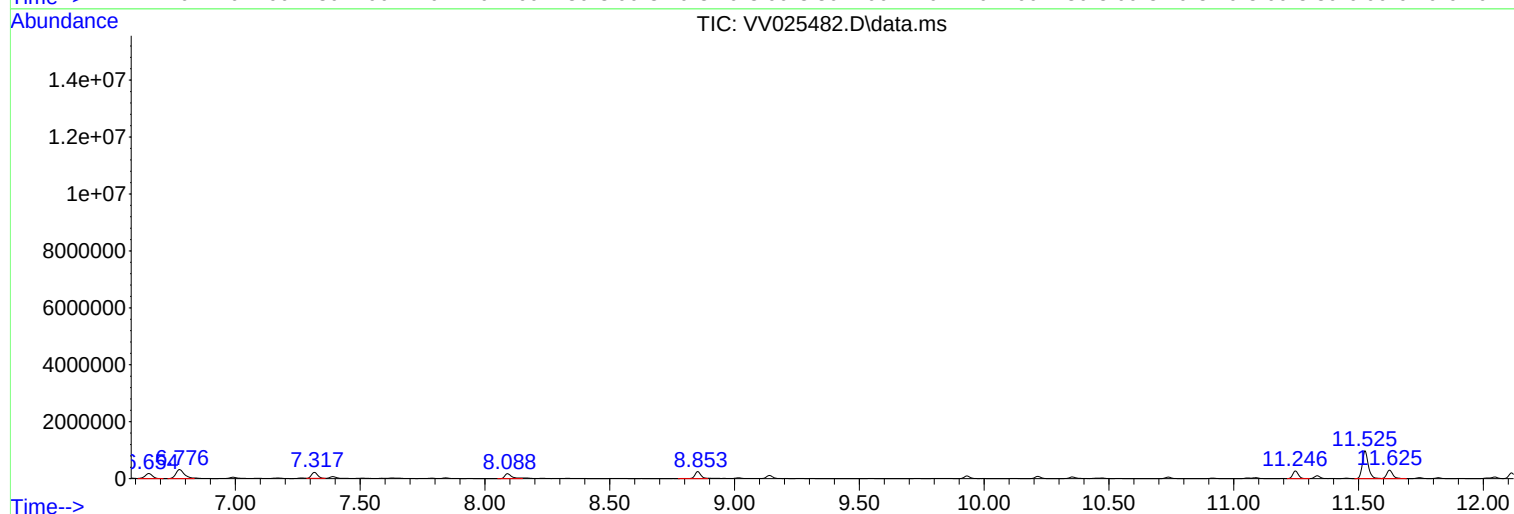
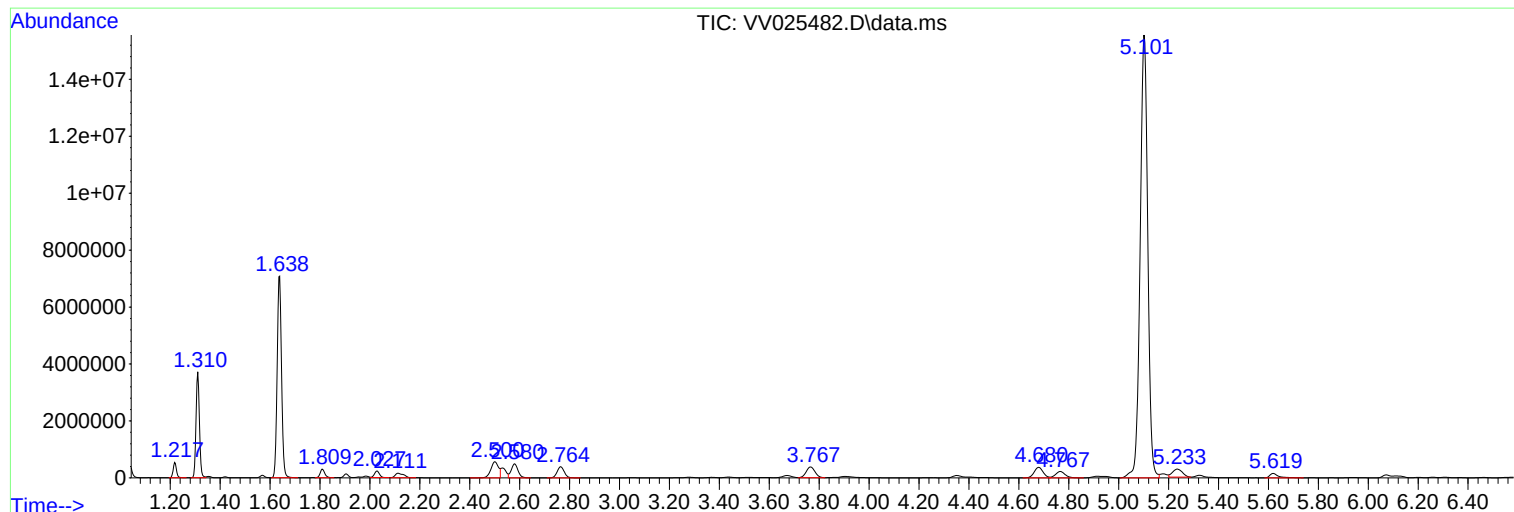
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR040822WMA.M
Quant Title : TRACE VOA SFAM1.0

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



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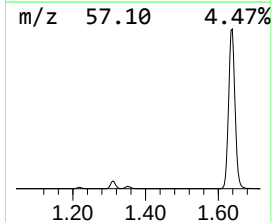
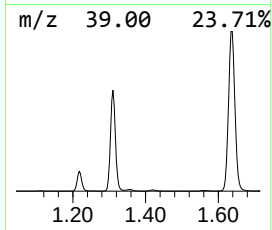
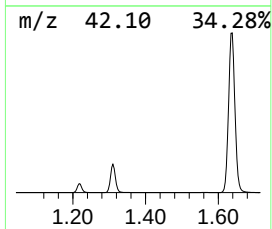
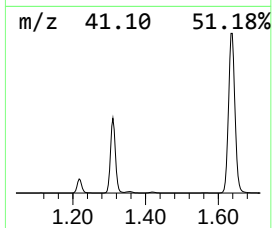
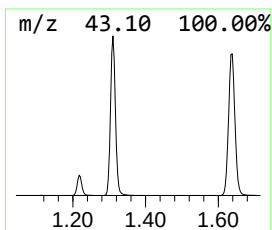
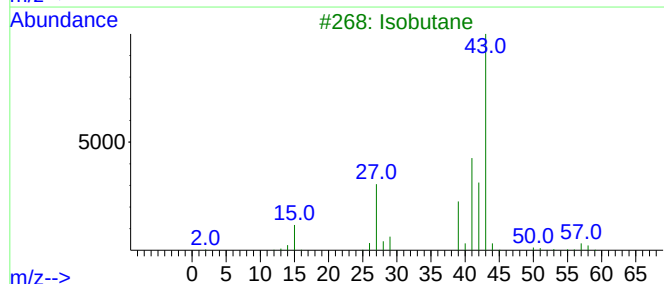
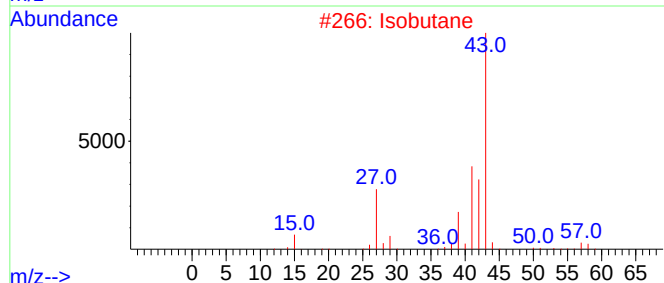
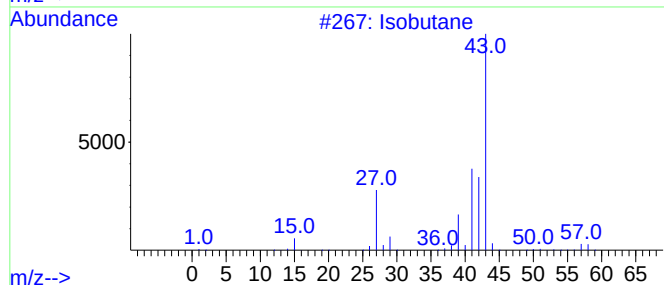
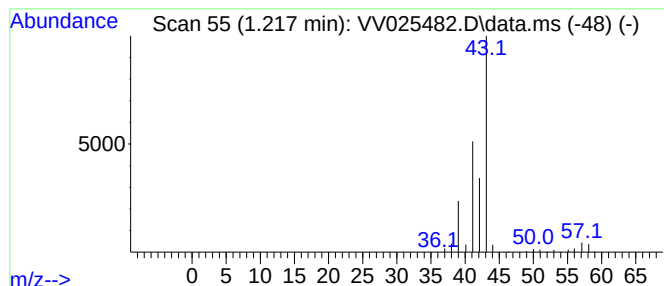
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.217	6.65 ug/L	487983	1,4-Difluorobenzene	5.619

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



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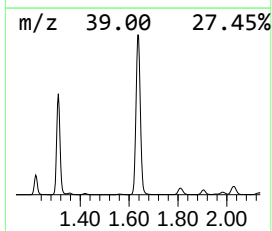
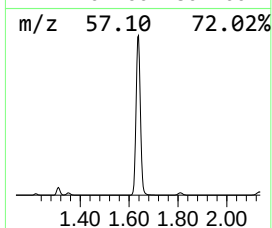
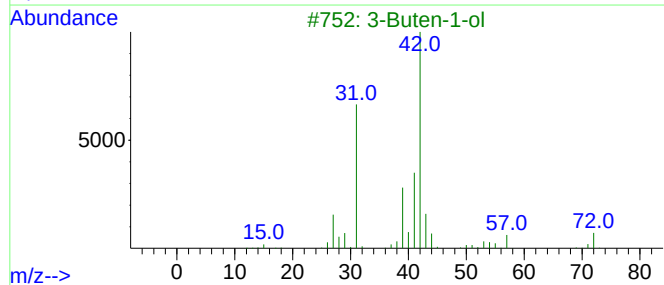
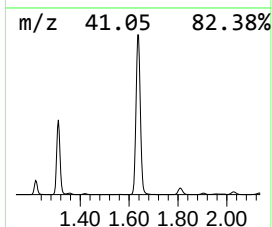
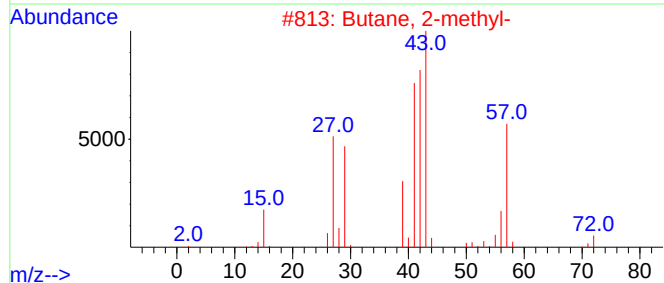
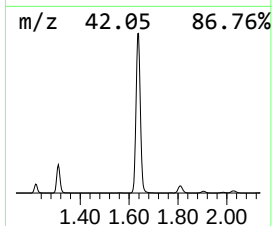
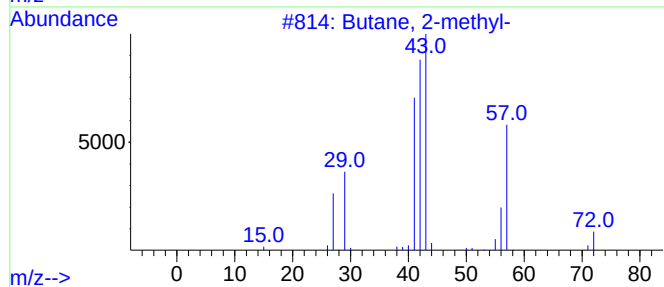
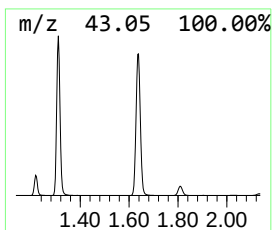
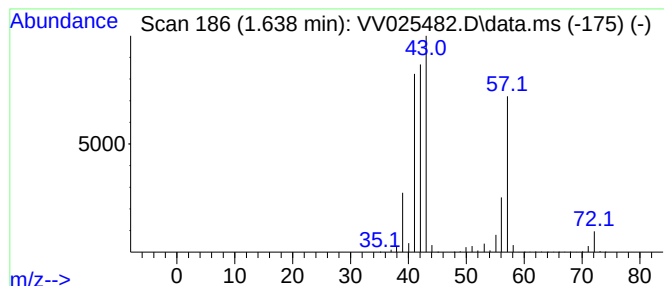
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR040822WMA.M
Quant Title : TRACE VOA SFAM1.0

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.
1.638	117.50 ug/L	8625540	1,4-Difluorobenzene	5.619

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2-methyl-	72	C5H12	000078-78-4	91
2		Butane, 2-methyl-	72	C5H12	000078-78-4	90
3		3-Buten-1-ol	72	C4H8O	000627-27-0	47
4		Pentane	72	C5H12	000109-66-0	43
5		1-Butene	56	C4H8	000106-98-9	10



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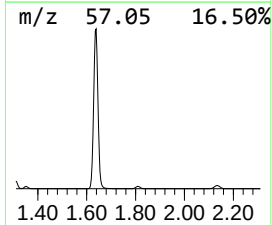
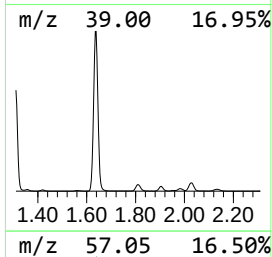
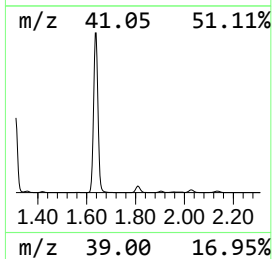
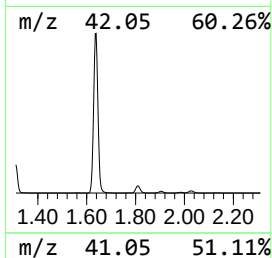
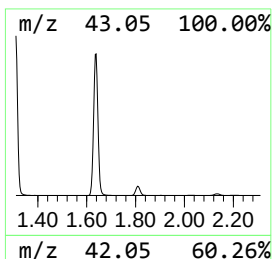
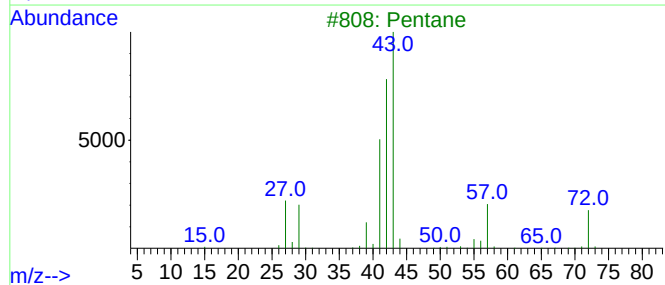
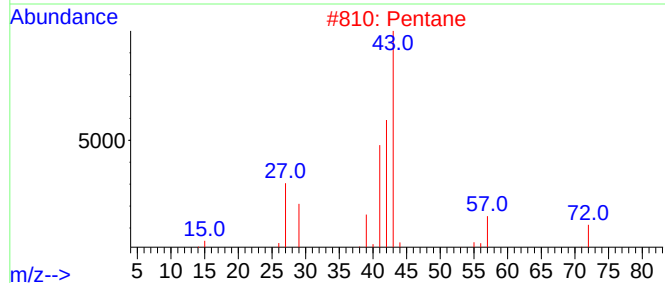
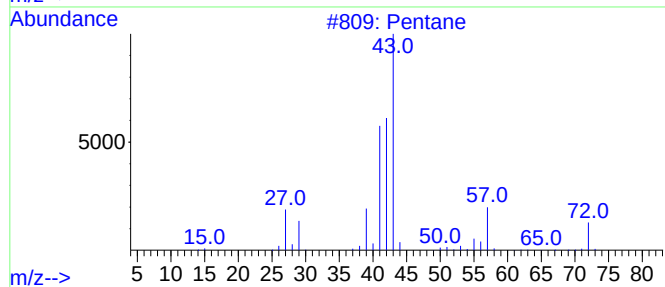
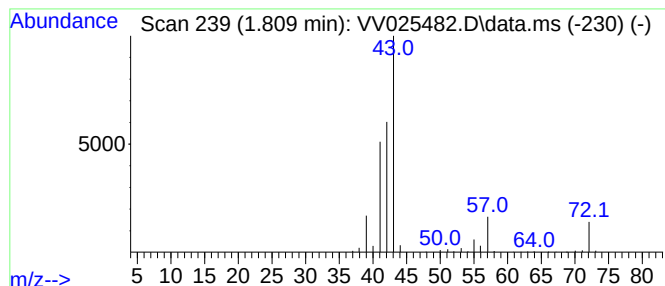
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR040822WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.809	5.36 ug/L	393829	1,4-Difluorobenzene	5.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane			72	C5H12	000109-66-0	90
2	Pentane			72	C5H12	000109-66-0	86
3	Pentane			72	C5H12	000109-66-0	86
4	Pentane			72	C5H12	000109-66-0	86
5	Pentane			72	C5H12	000109-66-0	80



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ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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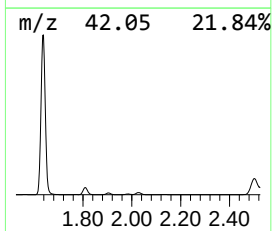
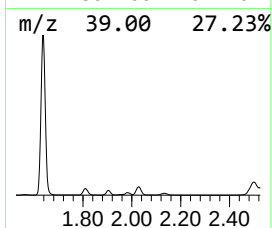
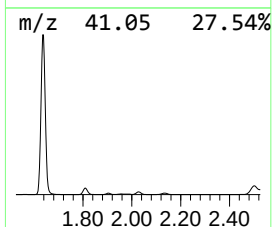
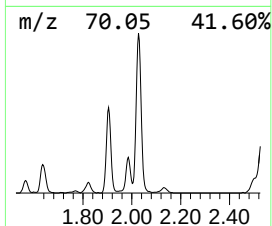
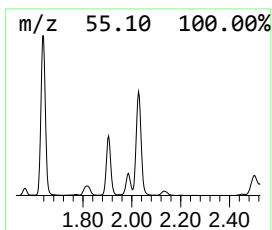
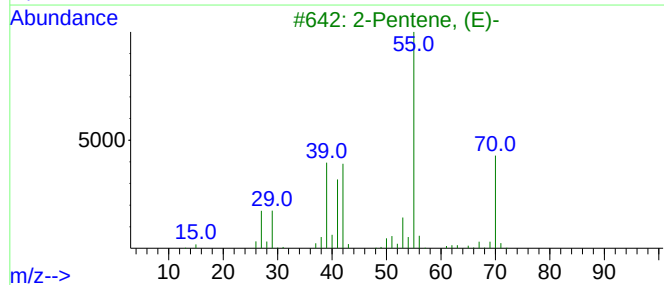
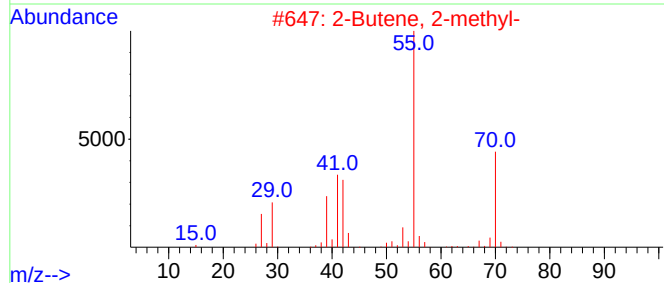
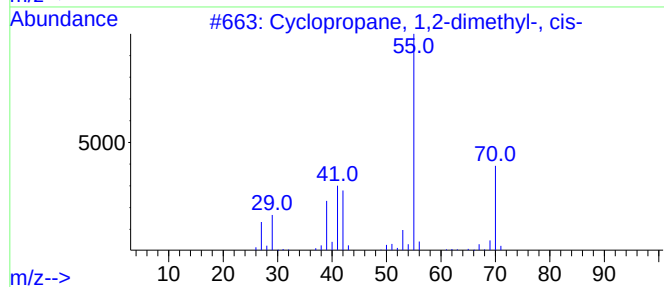
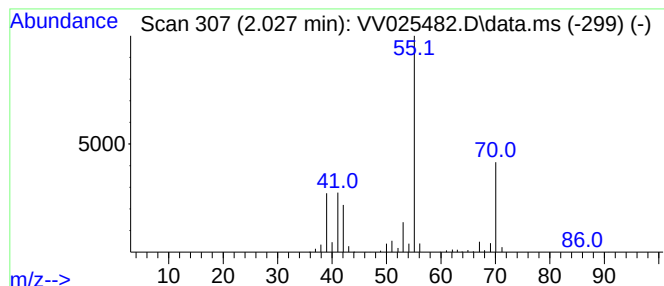
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 4 (DEL) Alkane: Cyclic2.027 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.027	4.52 ug/L	331713	1,4-Difluorobenzene	5.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	91
2			2-Butene, 2-methyl-	70	C5H10	000513-35-9	90
3			2-Pentene, (E)-	70	C5H10	000646-04-8	87
4			Cyclopropane, 1,2-dimethyl-, trans-	70	C5H10	002402-06-4	83
5			Cyclopropane, 1,1-dimethyl-	70	C5H10	001630-94-0	80



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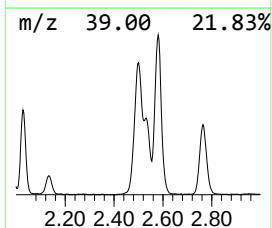
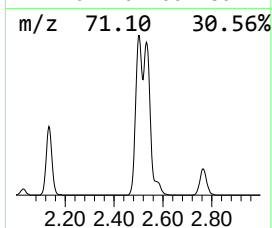
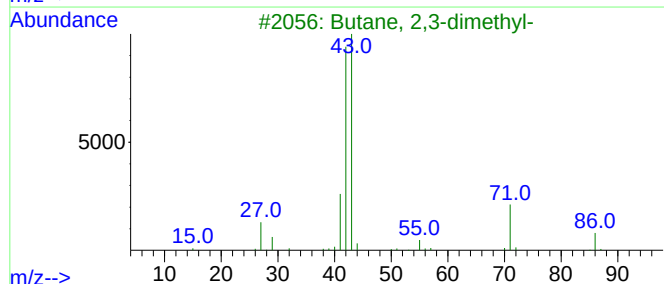
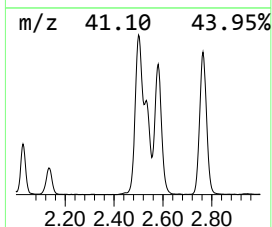
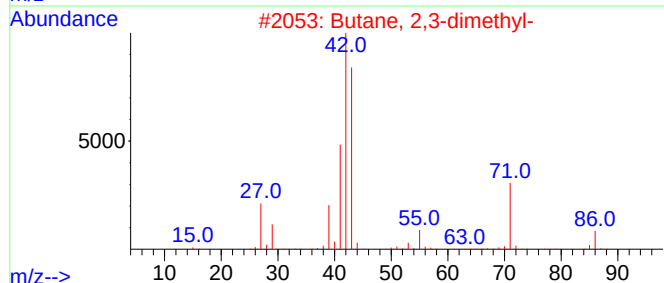
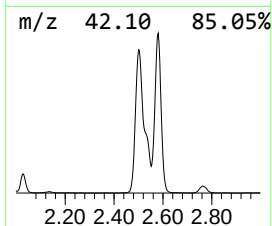
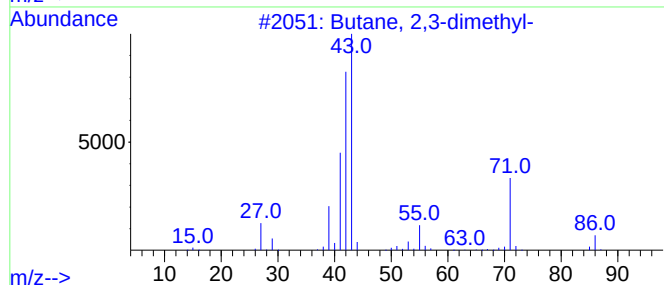
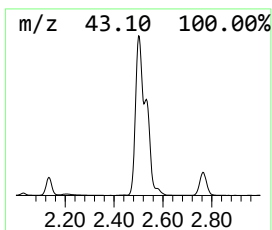
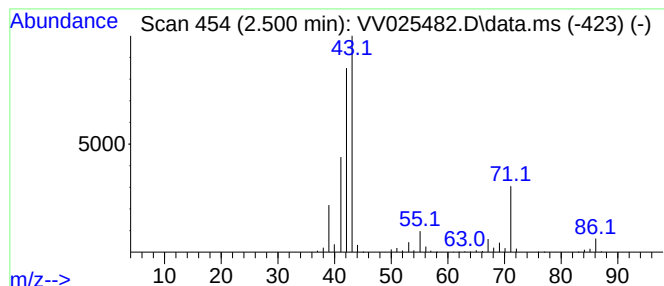
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.500	16.94 ug/L	1243640	1,4-Difluorobenzene	5.619

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	90
2	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	87
3	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	80
4	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	72
5	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	64



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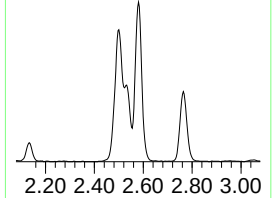
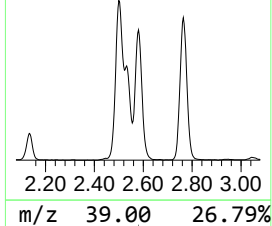
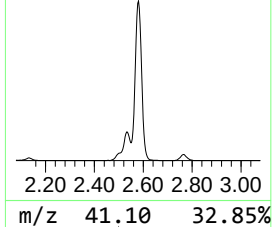
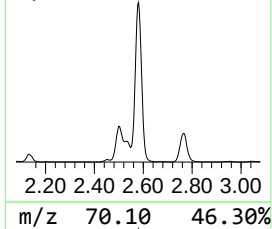
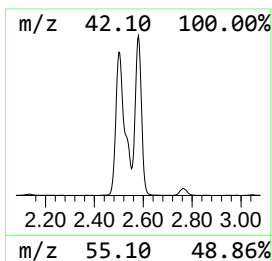
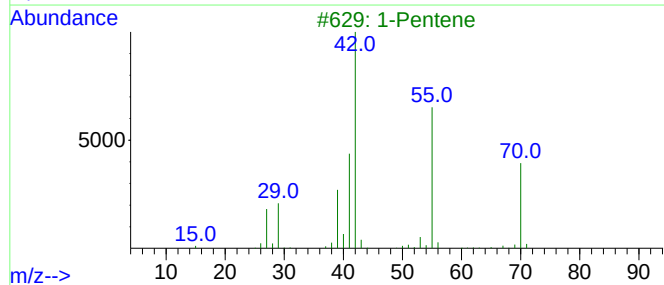
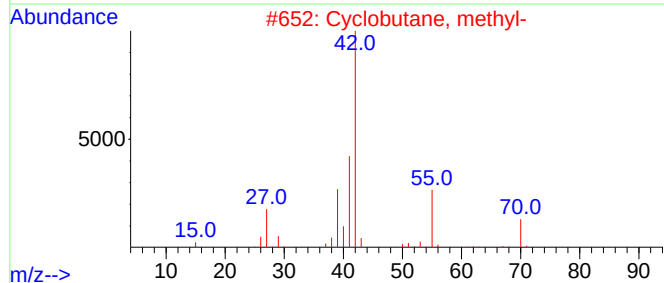
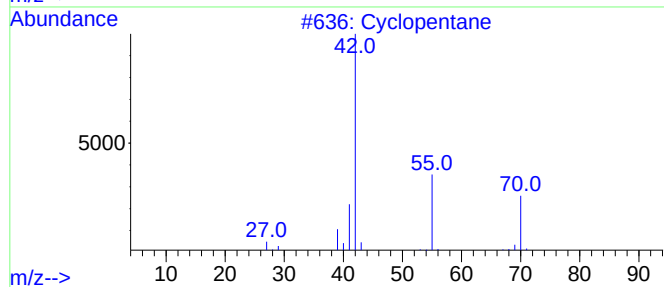
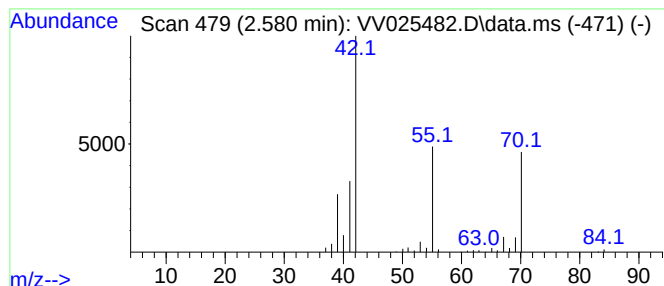
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 6 (DEL) Alkane: Cyclic2.580 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.580	12.48 ug/L	916011	1,4-Difluorobenzene	5.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane	70	C5H10	000287-92-3	80
2			Cyclobutane, methyl-	70	C5H10	000598-61-8	78
3			1-Pentene	70	C5H10	000109-67-1	78
4			1-Pentene	70	C5H10	000109-67-1	72
5			Cyclobutane, methyl-	70	C5H10	000598-61-8	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041222\
Data File : VV025482.D
Acq On : 12 Apr 2022 18:16
Operator : SY/MD
Sample : N2356-08
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBDS6

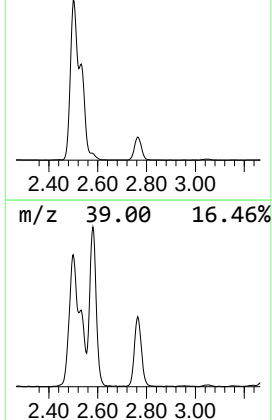
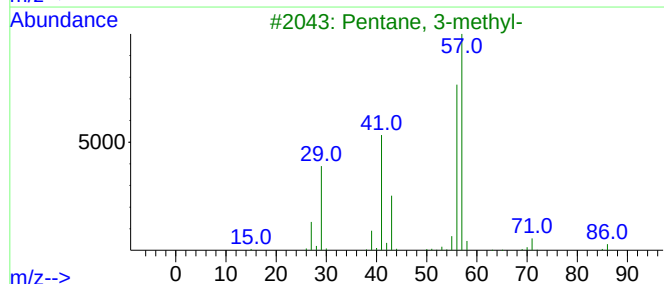
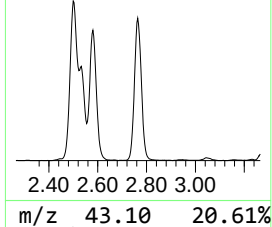
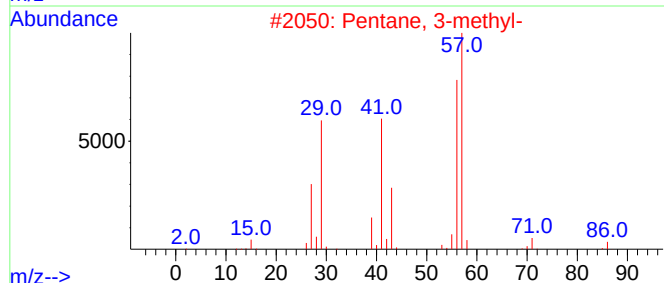
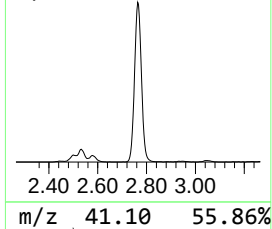
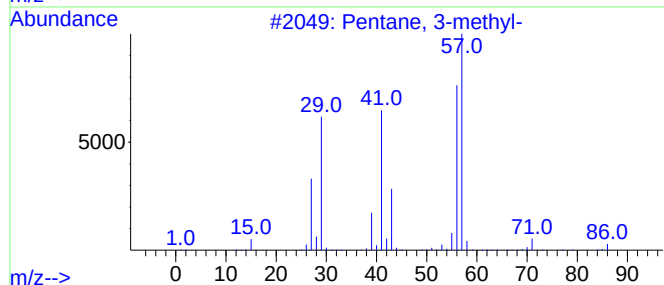
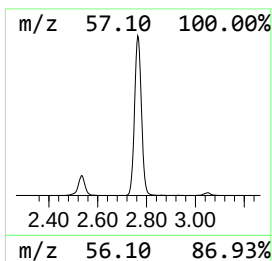
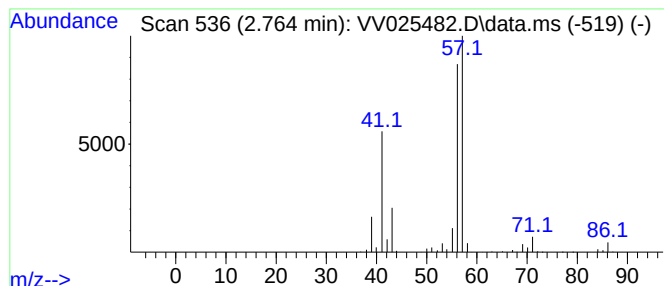
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.764	10.58 ug/L	776974	1,4-Difluorobenzene	5.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 3-methyl-	86	C6H14	000096-14-0	90
2			Pentane, 3-methyl-	86	C6H14	000096-14-0	90
3			Pentane, 3-methyl-	86	C6H14	000096-14-0	87
4			Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	78
5			Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041222\
Data File : VV025482.D
Acq On : 12 Apr 2022 18:16
Operator : SY/MD
Sample : N2356-08
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBDS6

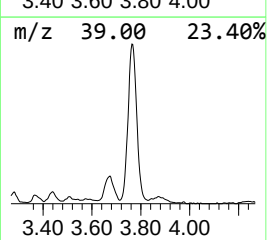
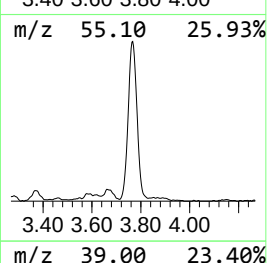
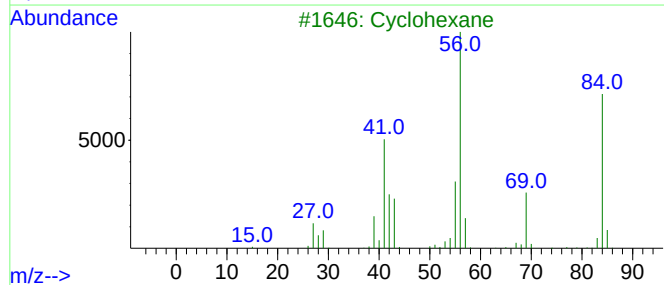
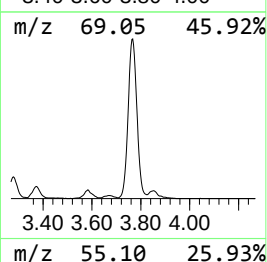
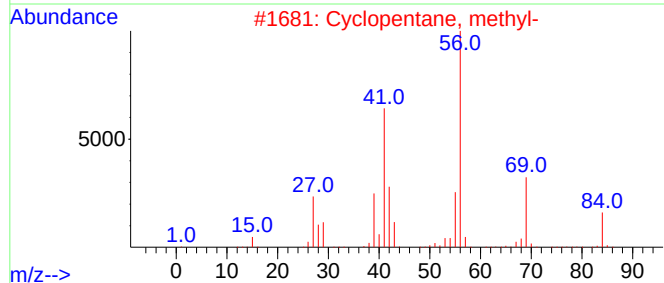
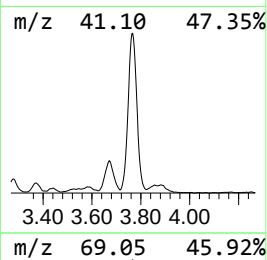
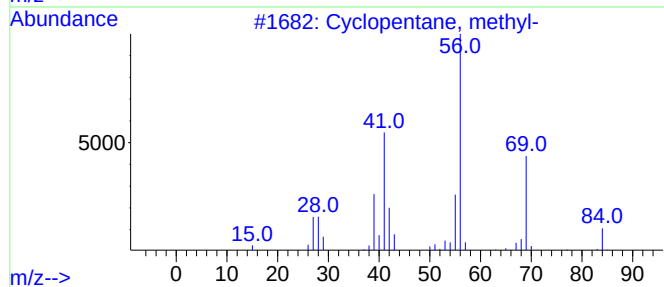
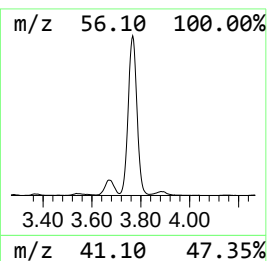
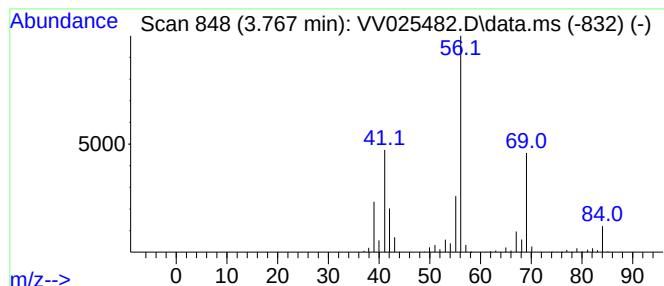
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 8 (DEL) Alkane: Cyclic3.767 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.767	12.75 ug/L	935932	1,4-Difluorobenzene	5.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, methyl-	84	C6H12	000096-37-7	90
2			Cyclopentane, methyl-	84	C6H12	000096-37-7	87
3			Cyclohexane	84	C6H12	000110-82-7	78
4			Cyclopentane, methyl-	84	C6H12	000096-37-7	72
5			Cyclopentane, methyl-	84	C6H12	000096-37-7	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041222\
Data File : VV025482.D
Acq On : 12 Apr 2022 18:16
Operator : SY/MD
Sample : N2356-08
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBDS6

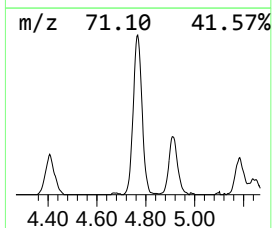
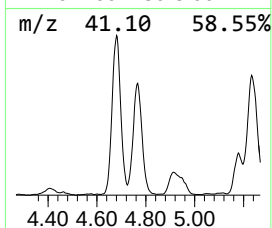
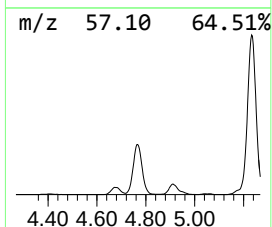
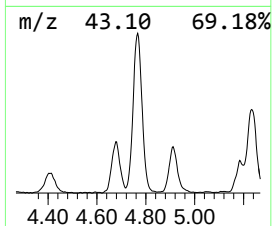
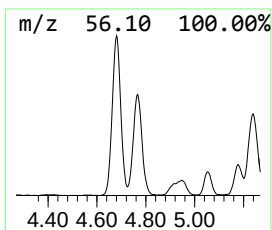
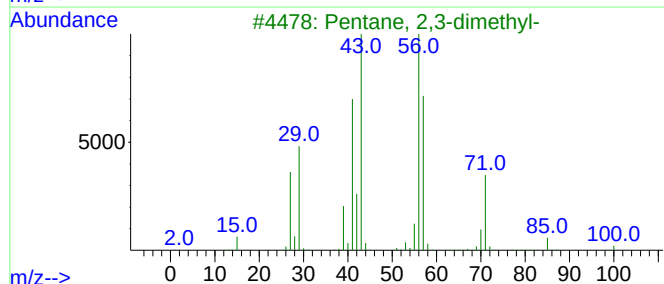
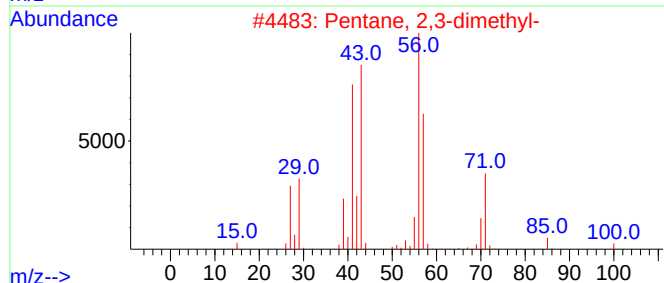
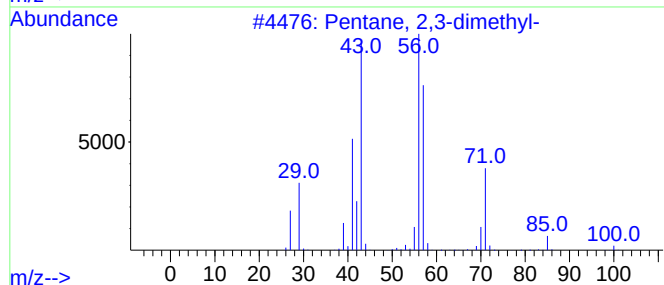
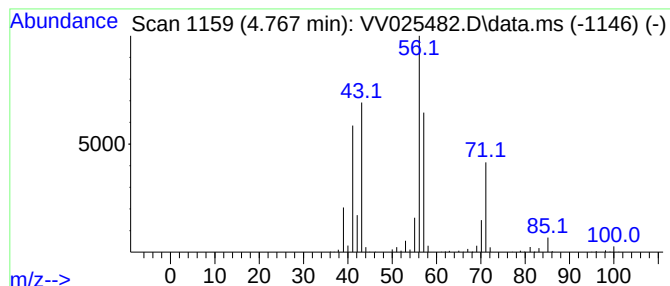
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Quant Title : TRACE VOA SFAM1.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.767	8.03 ug/L	589689	1,4-Difluorobenzene	5.619

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	87
2		Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	87
3		Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	83
4		Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	80
5		Heptane	100	C7H16	000142-82-5	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041222\
Data File : VV025482.D
Acq On : 12 Apr 2022 18:16
Operator : SY/MD
Sample : N2356-08
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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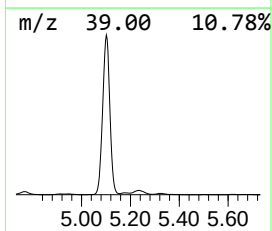
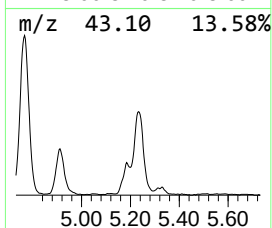
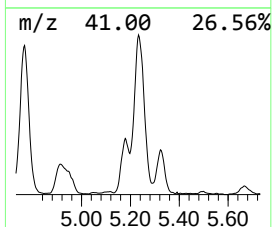
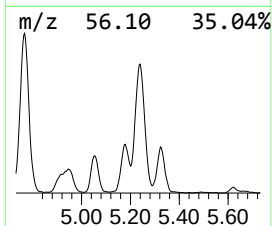
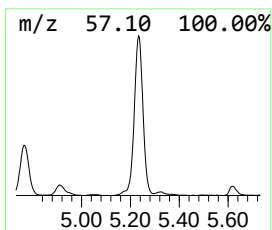
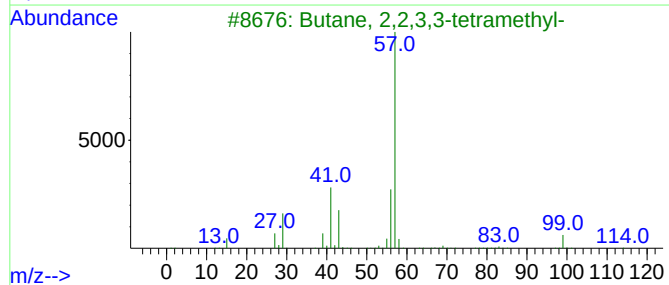
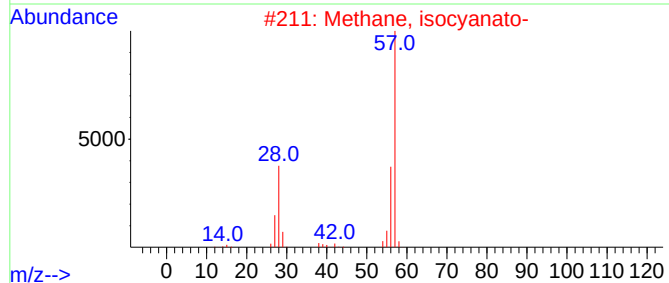
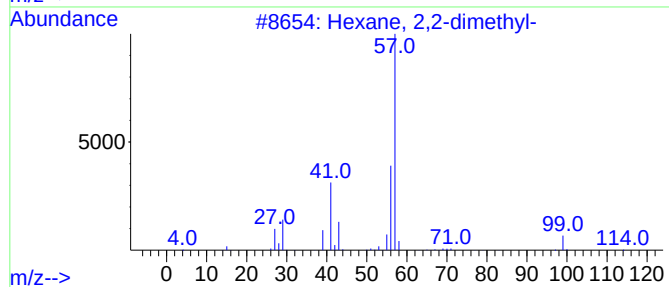
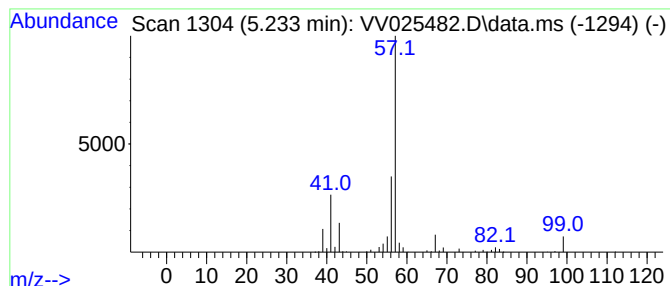
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.233	10.51 ug/L	771238	1,4-Difluorobenzene	5.619

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	59
2		Methane, isocyanato-	57	C2H3NO	000624-83-9	59
3		Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	59
4		Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	53
5		Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041222\
Data File : VV025482.D
Acq On : 12 Apr 2022 18:16
Operator : SY/MD
Sample : N2356-08
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBDS6

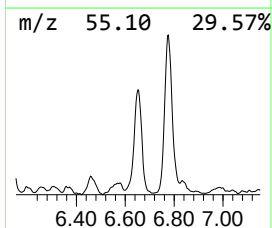
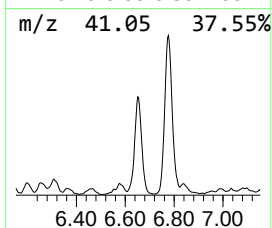
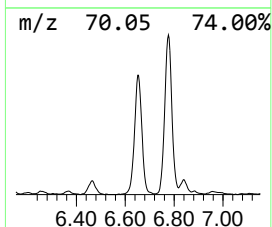
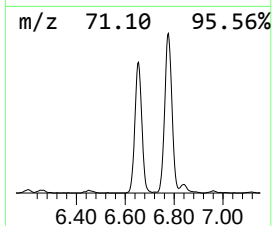
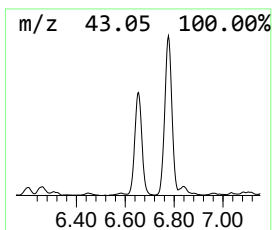
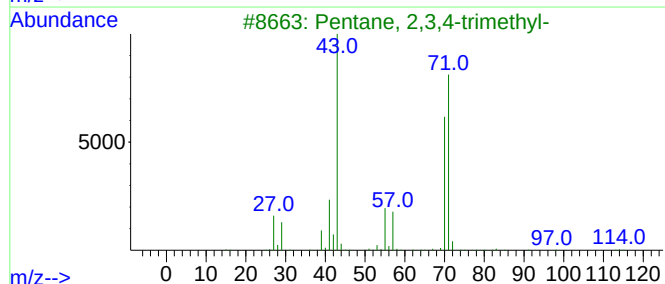
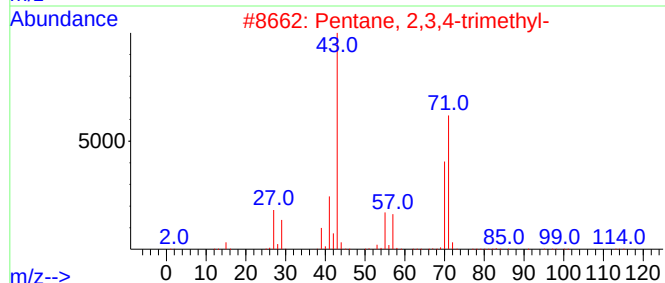
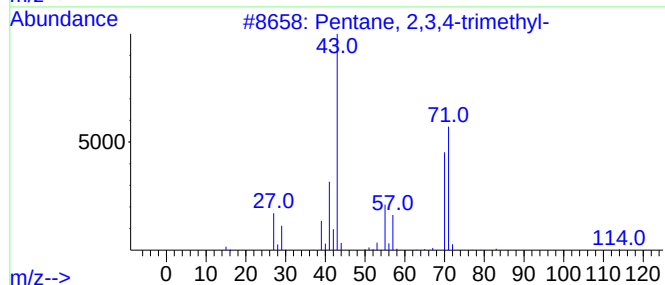
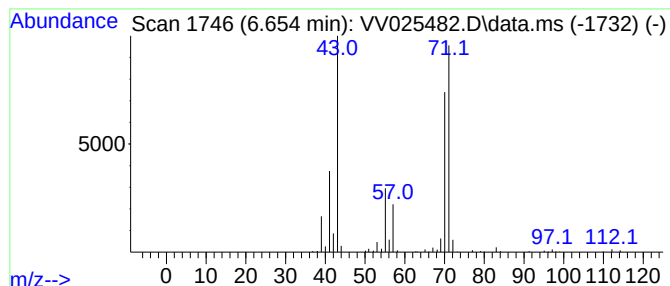
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.654	5.01 ug/L	367497	1,4-Difluorobenzene	5.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	80
2			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	80
3			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	80
4			Pentane, 3-ethyl-	100	C7H16	000617-78-7	78
5			3,3-Dimethyl-2-pentanol	116	C7H16O	019781-24-9	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041222\
Data File : VV025482.D
Acq On : 12 Apr 2022 18:16
Operator : SY/MD
Sample : N2356-08
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBDS6

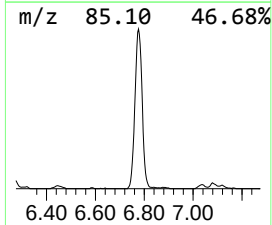
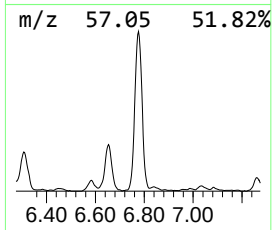
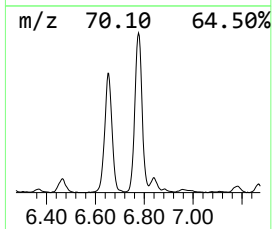
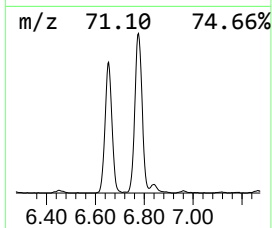
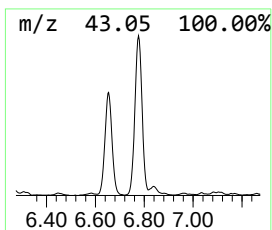
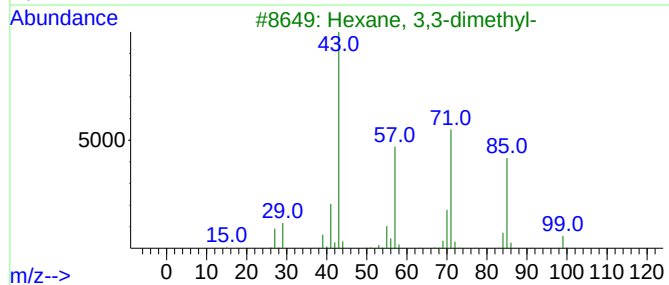
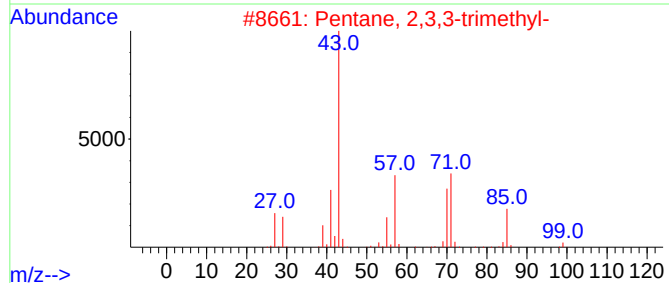
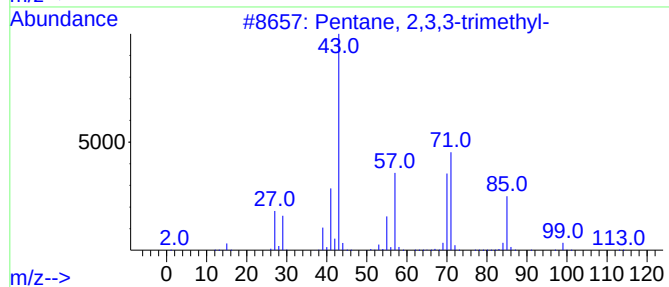
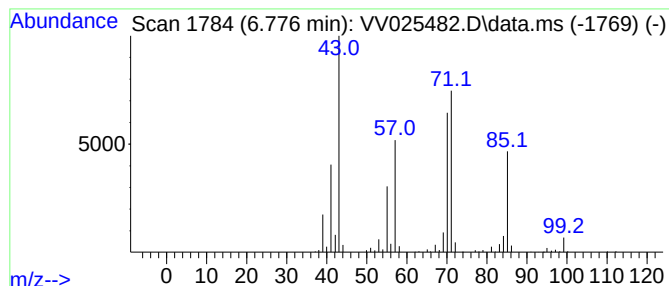
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.776	10.29 ug/L	755702	1,4-Difluorobenzene	5.619

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	90
2			Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	83
3			Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	64
4			Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	59
5			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041222\
Data File : VV025482.D
Acq On : 12 Apr 2022 18:16
Operator : SY/MD
Sample : N2356-08
Misc : 25mL/MSVOA_V/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBDS6

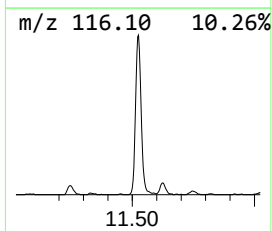
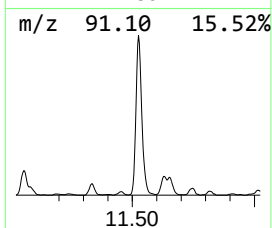
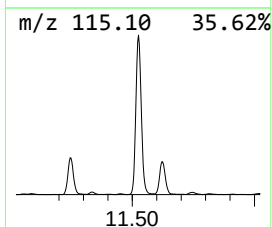
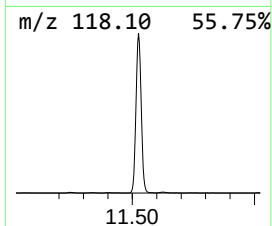
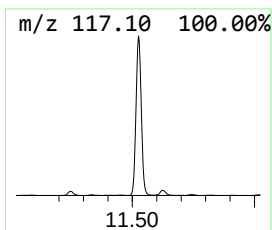
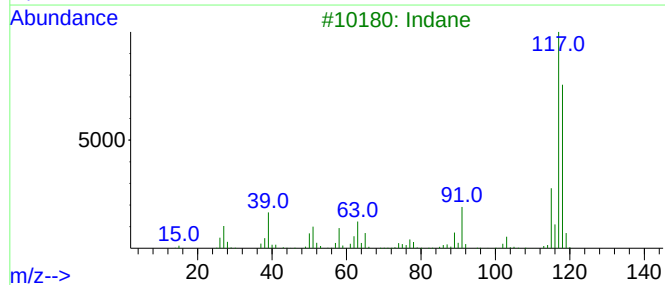
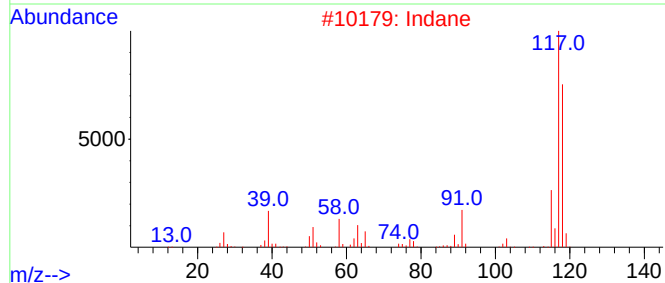
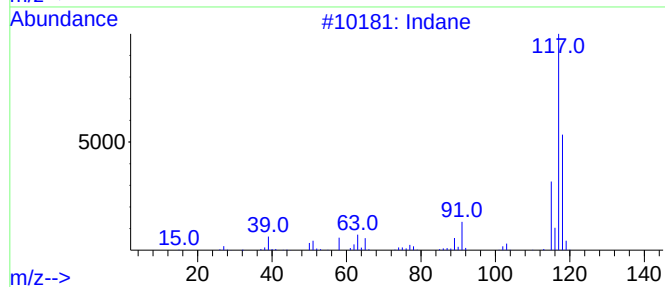
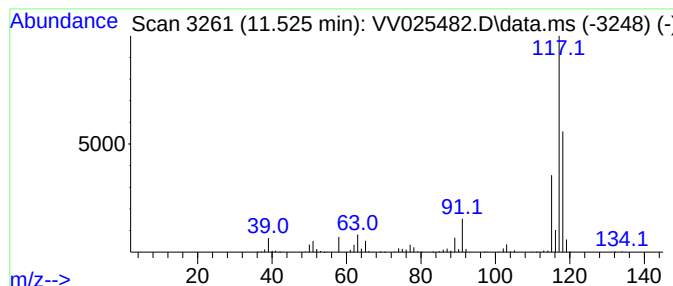
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 13 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.525	19.19 ug/L	1611340	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	95
2			Indane	118	C9H10	000496-11-7	91
3			Indane	118	C9H10	000496-11-7	90
4			Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	83
5			Benzene, cyclopropyl-	118	C9H10	000873-49-4	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV041222\
 Data File : VV025482.D
 Acq On : 12 Apr 2022 18:16
 Operator : SY/MD
 Sample : N2356-08
 Misc : 25mL/MSVOA_V/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 GBDS6

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR040822WMA.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...	1.217	6.7	ug/L	487983	1	5.619	367043	5.0
(DEL) Alkane: S...	1.638	117.5	ug/L	8625540	1	5.619	367043	5.0
(DEL) Alkane: S...	1.809	5.4	ug/L	393829	1	5.619	367043	5.0
(DEL) Alkane: C...	2.027	4.5	ug/L	331713	1	5.619	367043	5.0
(DEL) Alkane: S...	2.500	16.9	ug/L	1243640	1	5.619	367043	5.0
(DEL) Alkane: C...	2.580	12.5	ug/L	916011	1	5.619	367043	5.0
(DEL) Alkane: S...	2.764	10.6	ug/L	776974	1	5.619	367043	5.0
(DEL) Alkane: C...	3.767	12.8	ug/L	935932	1	5.619	367043	5.0
(DEL) Alkane: S...	4.767	8.0	ug/L	589689	1	5.619	367043	5.0
(DEL) Alkane: S...	5.233	10.5	ug/L	771238	1	5.619	367043	5.0
(DEL) Alkane: S...	6.654	5.0	ug/L	367497	1	5.619	367043	5.0
(DEL) Alkane: S...	6.776	10.3	ug/L	755702	1	5.619	367043	5.0
Indane	11.525	19.2	ug/L	1611340	3	11.249	419898	5.0