

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100920\
 Data File : VU040580.D
 Acq On : 09 Oct 2020 13:48
 Operator : SY/MD
 Sample : L4240-11DL 40X
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BG3T0DL

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.1

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR100820WMA.M

Title : TRACE VOA SOM01.0

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.607	75	83	94	rVB	93930	103202	2.39%	0.594%
2	1.922	172	181	197	rVB	73674	96487	2.23%	0.555%
3	2.578	373	385	404	rBV	153594	253441	5.86%	1.458%
4	3.099	518	547	566	rBV	454263	1025114	23.71%	5.897%
5	3.379	612	634	661	rBV	780492	1734119	40.11%	9.976%
6	3.716	720	739	766	rBV	1963229	4323007	100.00%	24.868%
7	4.314	907	925	949	rBV2	49115	142561	3.30%	0.820%
8	4.452	950	968	980	rVV3	37966	95702	2.21%	0.551%
9	4.549	980	998	1017	rVV	694271	1649440	38.15%	9.489%
10	4.658	1017	1032	1086	rVB2	105469	428459	9.91%	2.465%
11	5.079	1147	1163	1180	rBV	127490	282056	6.52%	1.623%
12	5.404	1246	1264	1278	rBV3	68838	157488	3.64%	0.906%
13	5.739	1346	1368	1389	rBV2	260155	675584	15.63%	3.886%
14	6.260	1515	1530	1554	rBV	251170	476088	11.01%	2.739%
15	6.703	1654	1668	1683	rBV	163614	321594	7.44%	1.850%
16	7.571	1925	1938	1954	rBV	81706	145841	3.37%	0.839%
17	7.903	2026	2041	2058	rBV	299827	511667	11.84%	2.943%
18	8.185	2113	2129	2150	rBV	55316	96687	2.24%	0.556%
19	8.639	2257	2270	2303	rBV	484388	912823	21.12%	5.251%
20	9.420	2499	2513	2545	rVB	363822	609240	14.09%	3.505%
21	10.754	2917	2928	2947	rVB	133163	215507	4.99%	1.240%
22	10.902	2959	2974	2990	rBV	77508	115588	2.67%	0.665%
23	10.996	2990	3003	3009	rBV	249128	429699	9.94%	2.472%
24	11.086	3021	3031	3048	rVB	145245	225779	5.22%	1.299%
25	11.288	3081	3094	3105	rBV	74891	115083	2.66%	0.662%
26	11.462	3134	3148	3164	rBV	378528	581059	13.44%	3.343%
27	11.806	3241	3255	3270	rBV	399812	639829	14.80%	3.681%
28	11.889	3270	3281	3293	rVB	58996	94965	2.20%	0.546%
29	12.185	3361	3373	3390	rVB2	363376	617344	14.28%	3.551%
30	12.565	3480	3491	3501	rBV	50589	75642	1.75%	0.435%
31	12.963	3599	3615	3623	rBV2	29387	46576	1.08%	0.268%
32	13.018	3623	3632	3642	rVB	47042	70412	1.63%	0.405%
33	13.442	3755	3764	3780	rVB4	35948	64499	1.49%	0.371%
34	13.828	3872	3884	3901	rVB5	27279	50952	1.18%	0.293%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100920\
Data File : VU040580.D
Acq On : 09 Oct 2020 13:48
Operator : SY/MD
Sample : L4240-11DL 40X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3T0DL

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\SOMUTR100820WMA.M
Title : TRACE VOA SOM01.0

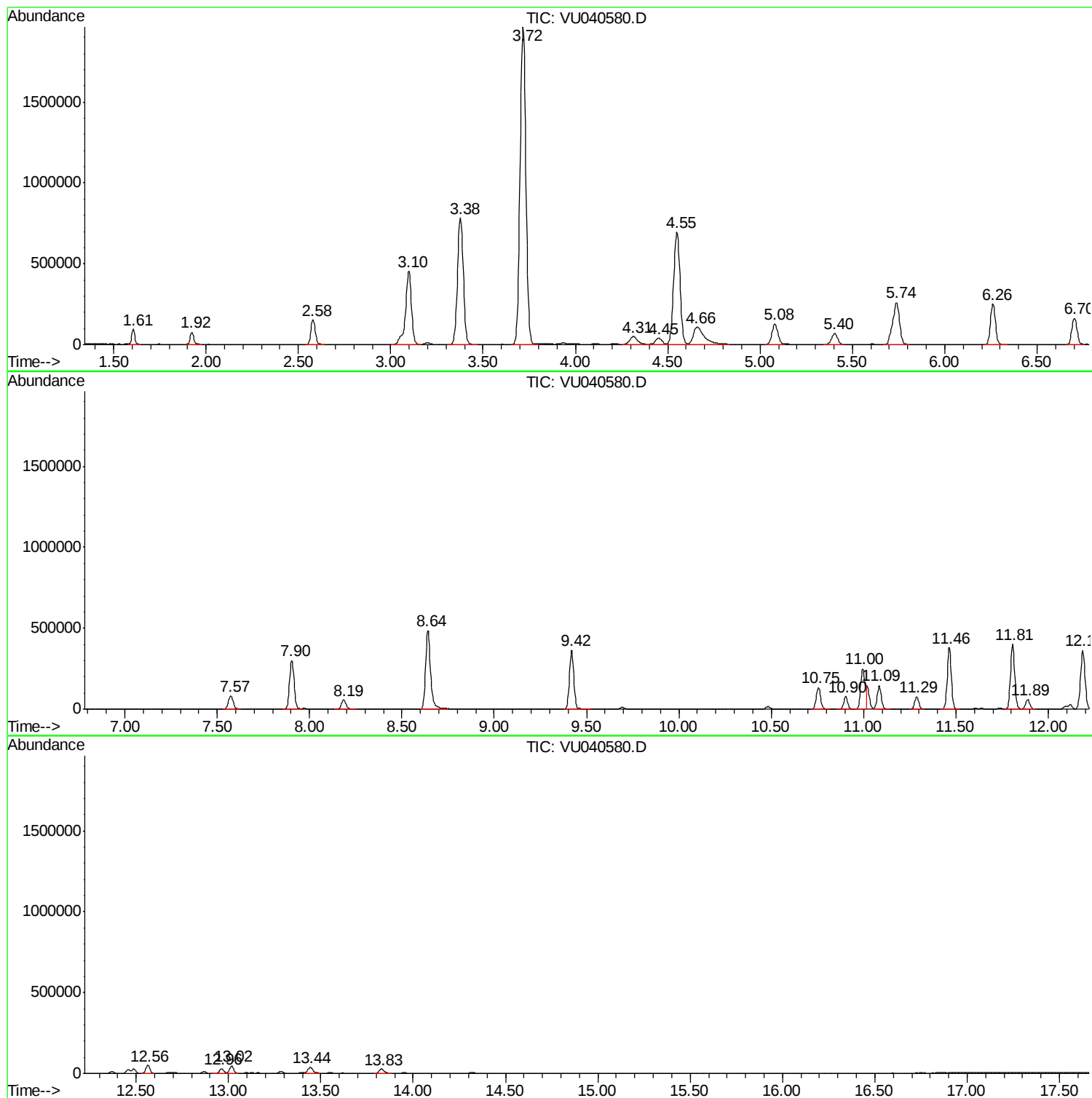
Sum of corrected areas: 17383534

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100920\
Data File : VU040580.D
Acq On : 09 Oct 2020 13:48
Operator : SY/MD
Sample : L4240-11DL 40X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3T0DL

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR100820WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100920\
Data File : VU040580.D
Acq On : 09 Oct 2020 13:48
Operator : SY/MD
Sample : L4240-11DL 40X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3T0DL

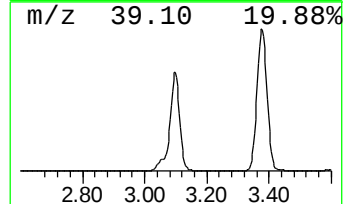
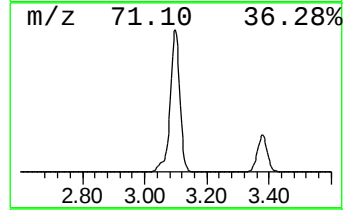
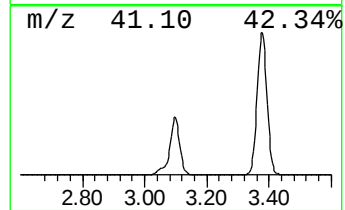
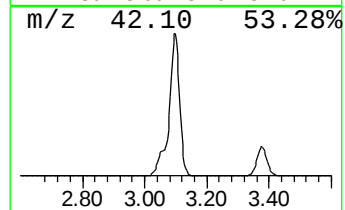
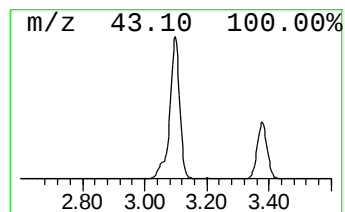
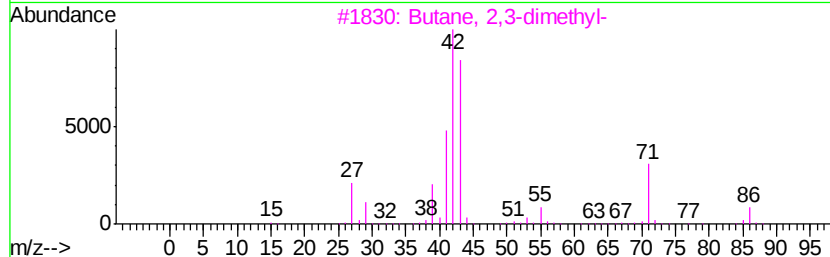
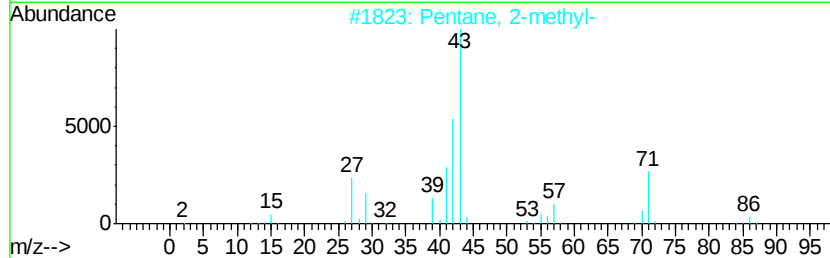
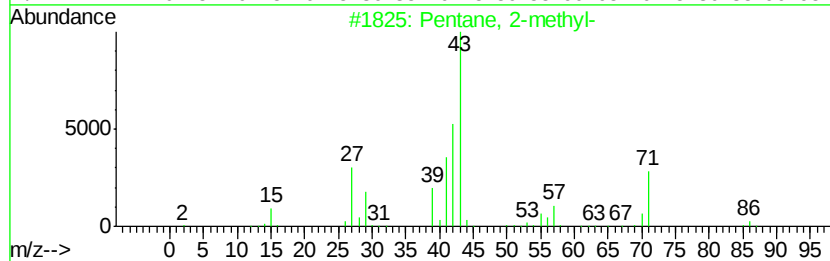
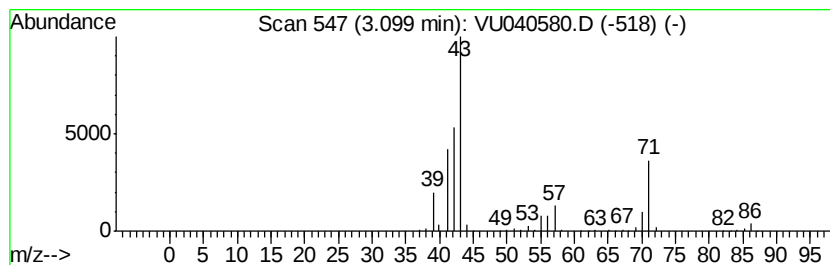
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR100820WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.10	10.77 ug/L	1025110	1,4-Difluorobenzene	6.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2-methyl-	86	C6H14	000107-83-5	91
2		Pentane, 2-methyl-	86	C6H14	000107-83-5	91
3		Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	53
4		Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	50
5		1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100920\
Data File : VU040580.D
Acq On : 09 Oct 2020 13:48
Operator : SY/MD
Sample : L4240-11DL 40X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3T0DL

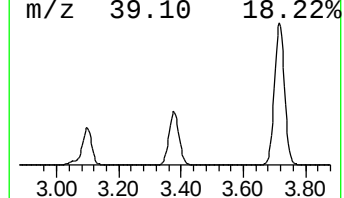
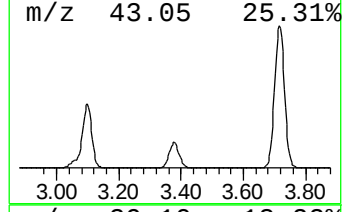
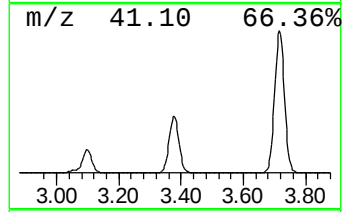
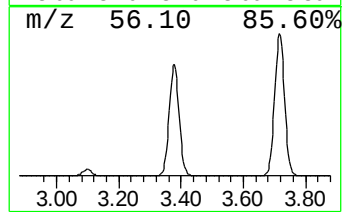
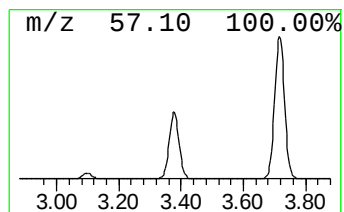
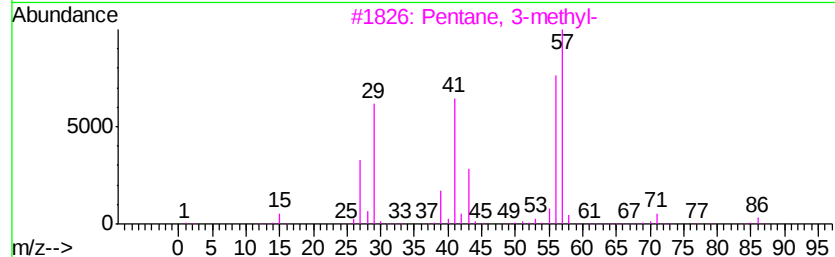
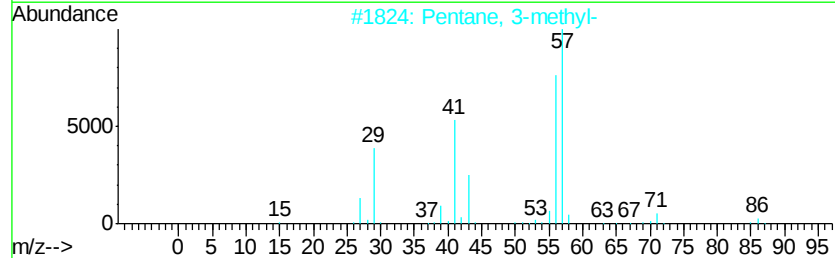
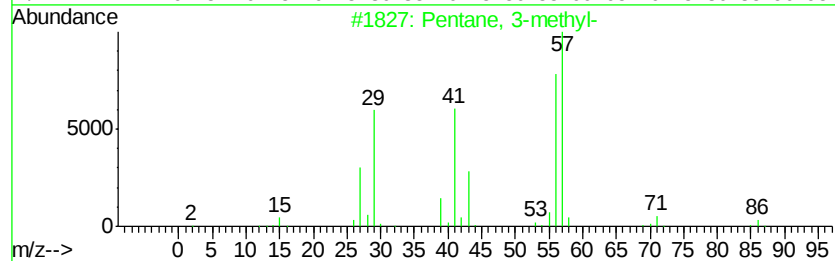
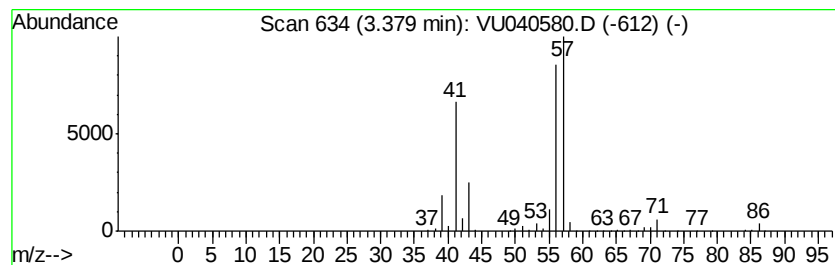
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR100820WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.38	18.21 ug/L	1734120	1,4-Difluorobenzene	6.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2		Pentane, 3-methyl-	86	C6H14	000096-14-0	91
3		Pentane, 3-methyl-	86	C6H14	000096-14-0	91
4		Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	78
5		Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	56



Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100920\
Data File : VU040580.D
Acq On : 09 Oct 2020 13:48
Operator : SY/MD
Sample : L4240-11DL 40X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3T0DL

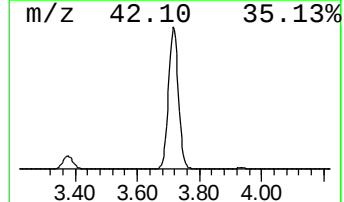
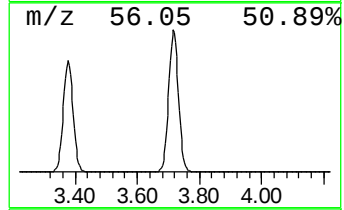
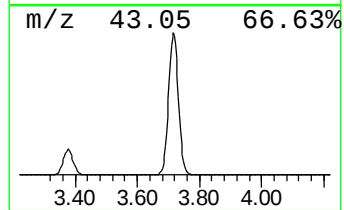
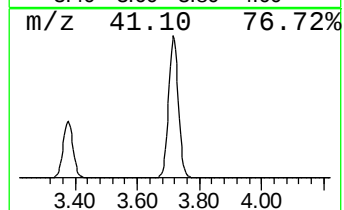
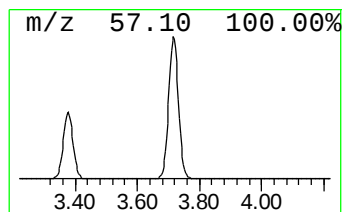
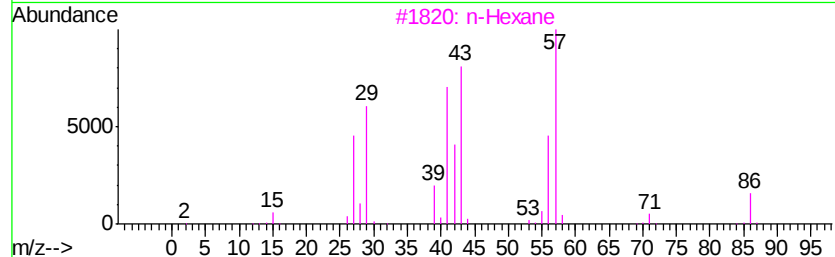
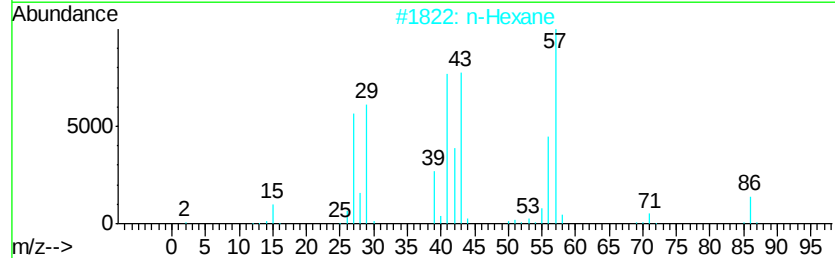
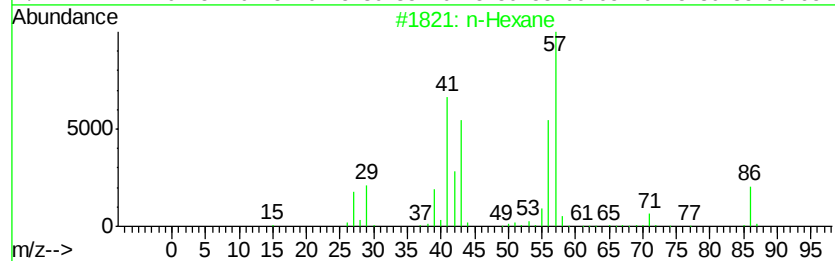
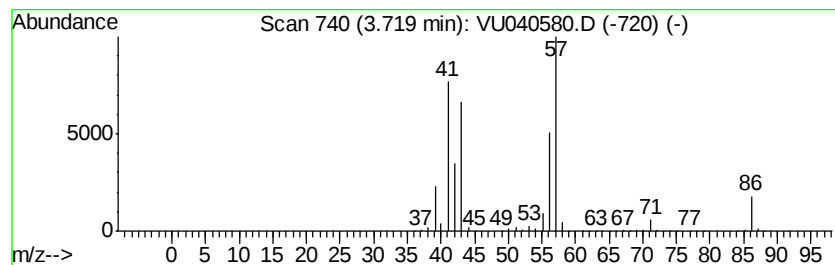
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR100820WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.72	45.40 ug/L	4323010	1,4-Difluorobenzene	6.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexane	86	C6H14	000110-54-3	91
2		n-Hexane	86	C6H14	000110-54-3	91
3		n-Hexane	86	C6H14	000110-54-3	91
4		Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	40
5		1-Pentene, 4-methyl-	84	C6H12	000691-37-2	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100920\
Data File : VU040580.D
Acq On : 09 Oct 2020 13:48
Operator : SY/MD
Sample : L4240-11DL 40X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3T0DL

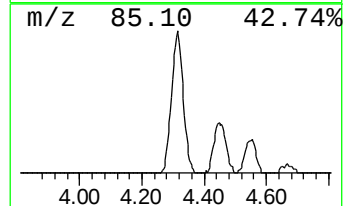
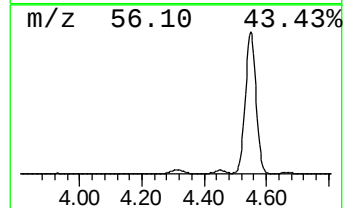
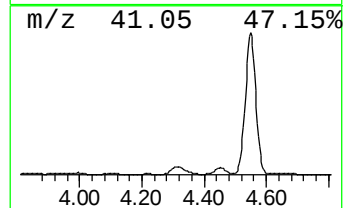
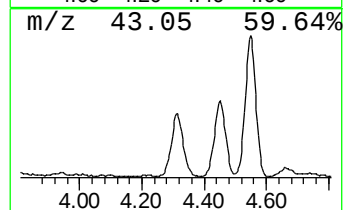
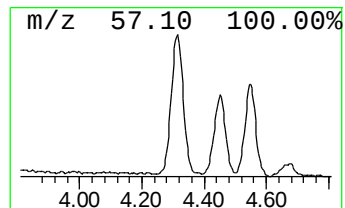
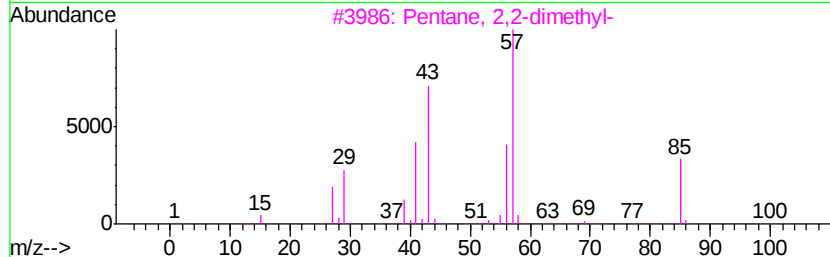
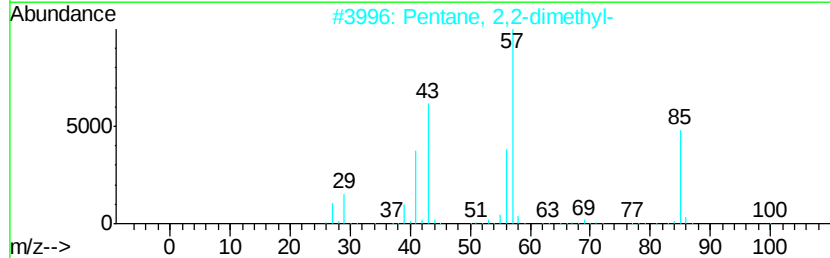
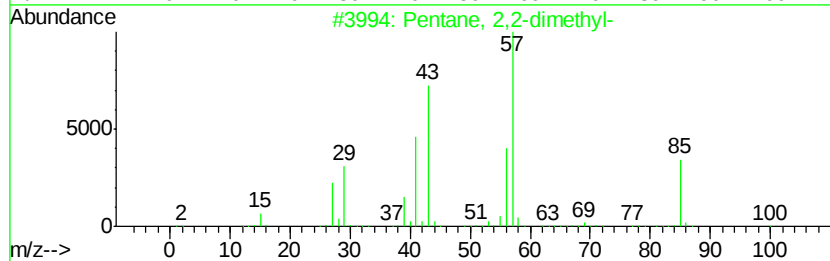
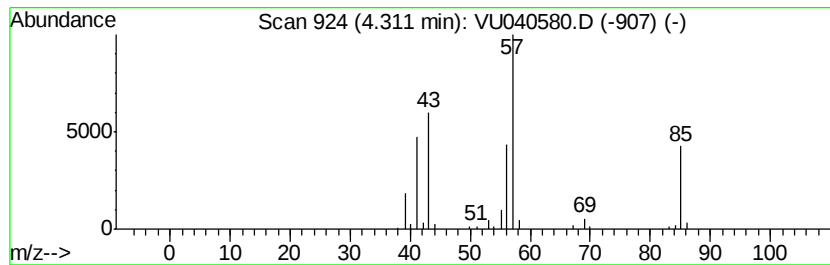
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR100820WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.31	1.50 ug/L	142561	1,4-Difluorobenzene	6.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	83
2		Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	83
3		Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	83
4		Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	78
5		Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	74



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100920\
Data File : VU040580.D
Acq On : 09 Oct 2020 13:48
Operator : SY/MD
Sample : L4240-11DL 40X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3T0DL

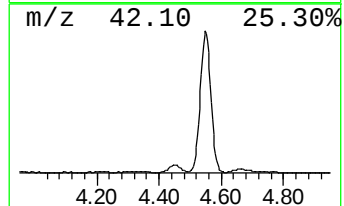
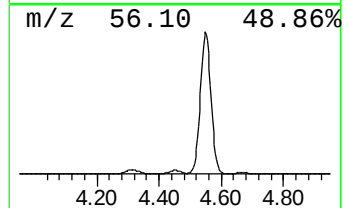
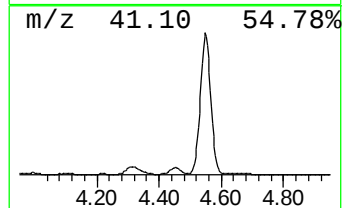
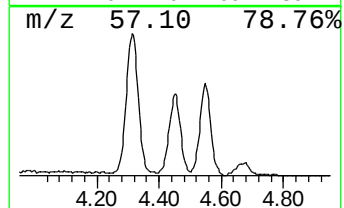
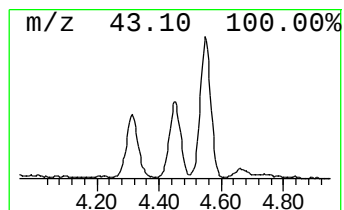
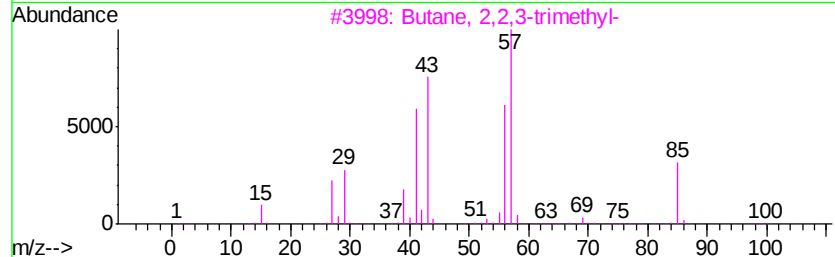
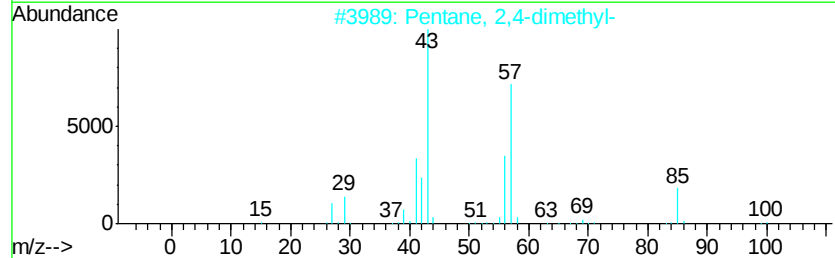
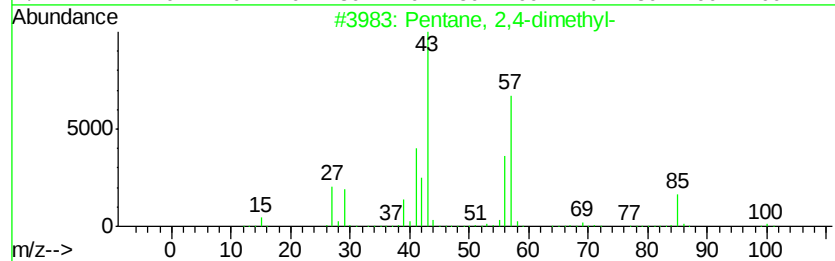
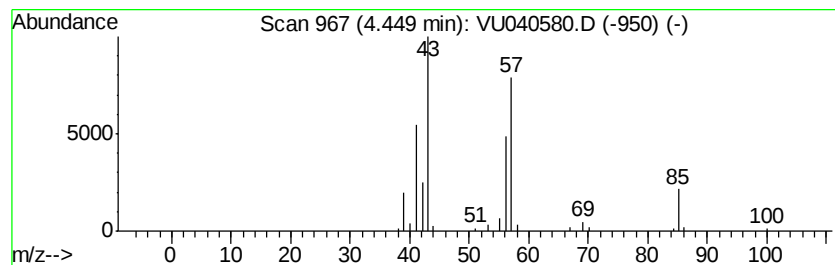
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR100820WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.45	1.01 ug/L	95702	1,4-Difluorobenzene	6.26

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	91	
2	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	83	
3	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	59	
4	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	59	
5	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	50	



Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100920\
Data File : VU040580.D
Acq On : 09 Oct 2020 13:48
Operator : SY/MD
Sample : L4240-11DL 40X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3T0DL

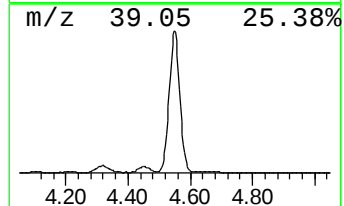
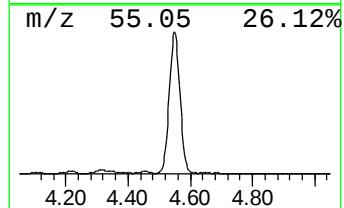
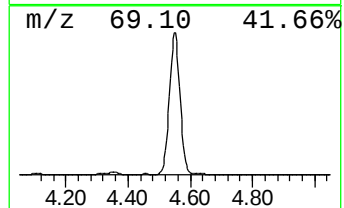
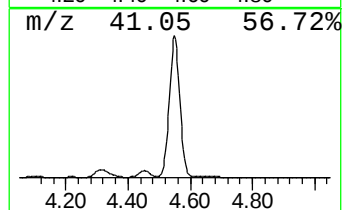
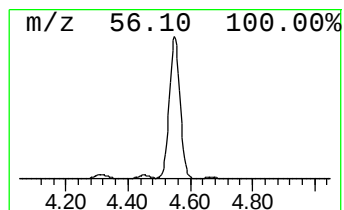
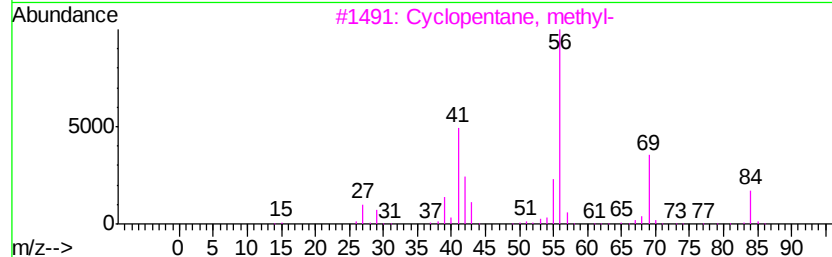
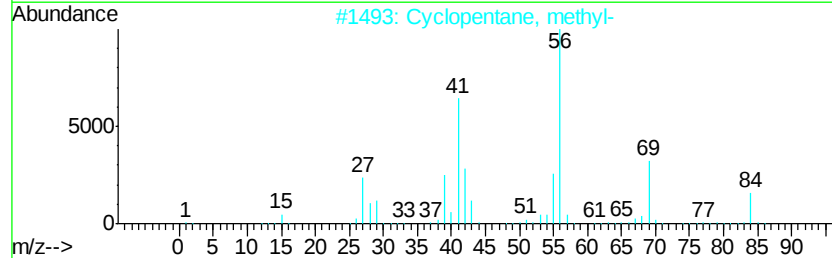
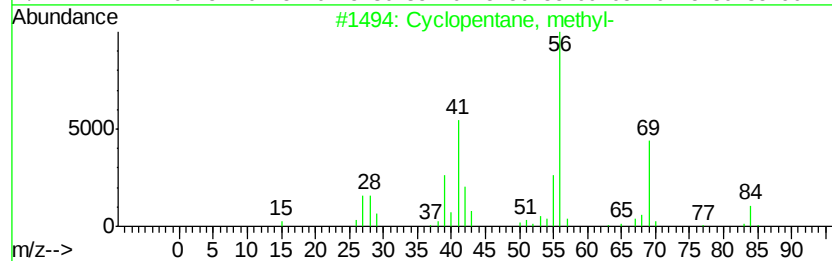
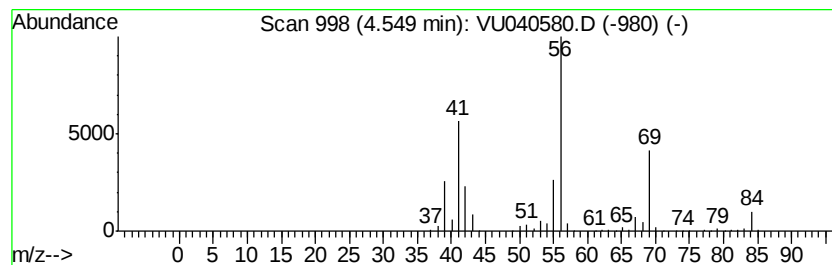
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR100820WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 (DEL) Alkane: Cyclic4.55 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.55	17.32 ug/L	1649440	1,4-Difluorobenzene	6.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2		Cyclopentane, methyl-	84	C6H12	000096-37-7	91
3		Cyclopentane, methyl-	84	C6H12	000096-37-7	78
4		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
5		Cyclobutane, ethyl-	84	C6H12	004806-61-5	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100920\
Data File : VU040580.D
Acq On : 09 Oct 2020 13:48
Operator : SY/MD
Sample : L4240-11DL 40X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 6 Sample Multiplier: 1

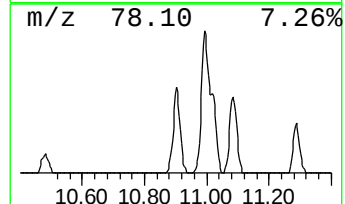
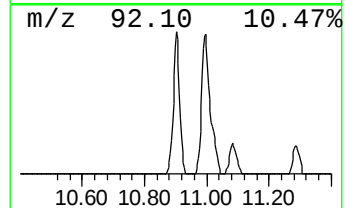
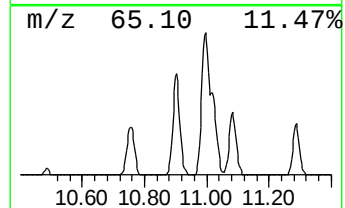
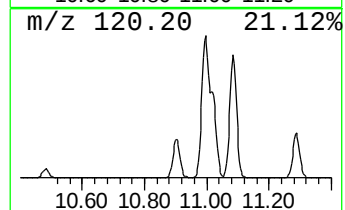
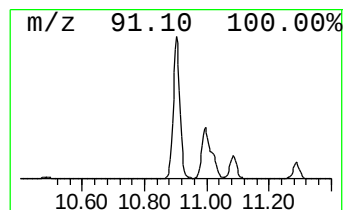
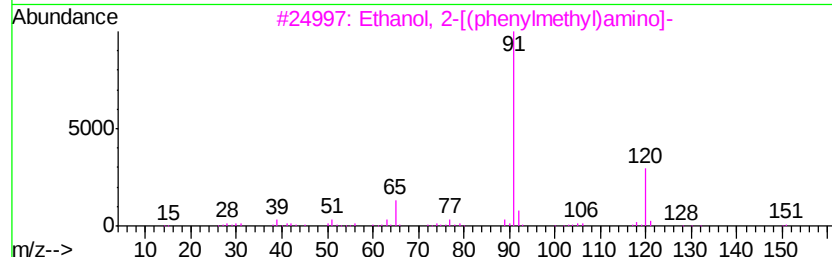
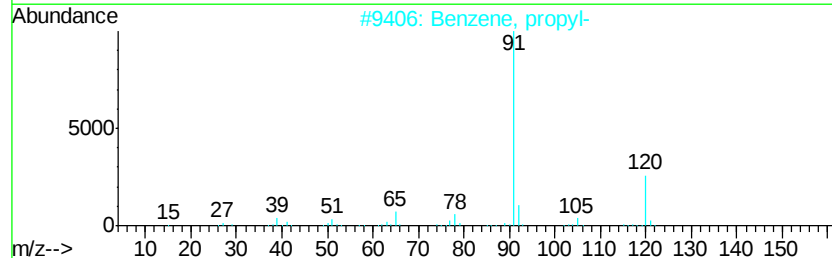
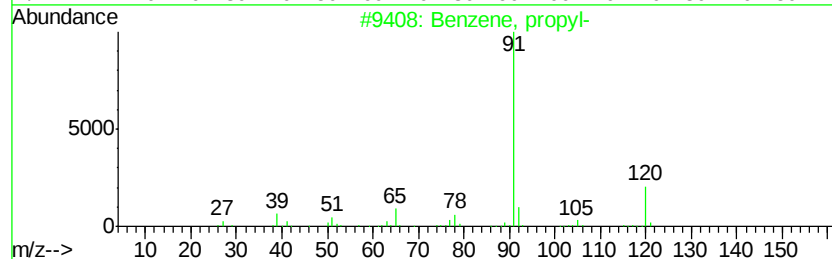
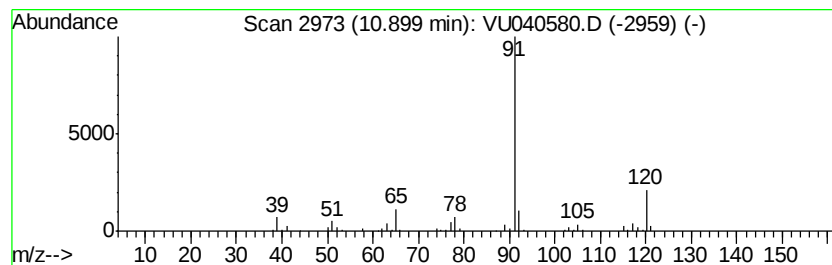
Instrument :
MSVOA_U
ClientSampleId :
BG3T0DL

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR100820WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzene, propyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.90	0.90 ug/L	115588	1,4-Dichlorobenzene-d4	11.81
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, propyl-	120 C9H12	000103-65-1 87
2		Benzene, propyl-	120 C9H12	000103-65-1 83
3		Ethanol, 2-[(phenylmethyl)amino]-	151 C9H13NO	000104-63-2 72
4		1,2-Ethanediamine, N-(phenylmeth...	150 C9H14N2	004152-09-4 72
5		1,2-Ethanediamine, N,N'-bis(phen...	240 C16H20N2	000140-28-3 72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100920\
Data File : VU040580.D
Acq On : 09 Oct 2020 13:48
Operator : SY/MD
Sample : L4240-11DL 40X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3T0DL

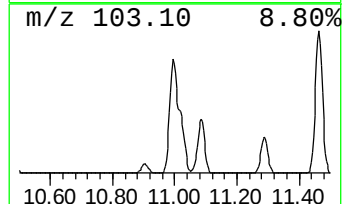
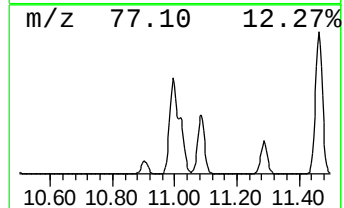
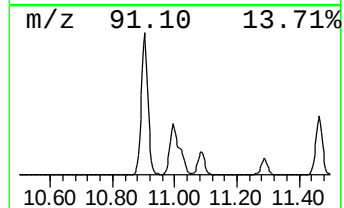
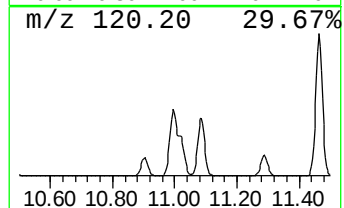
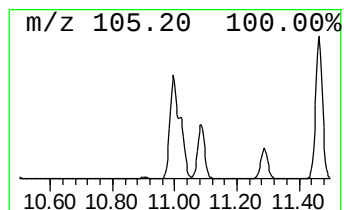
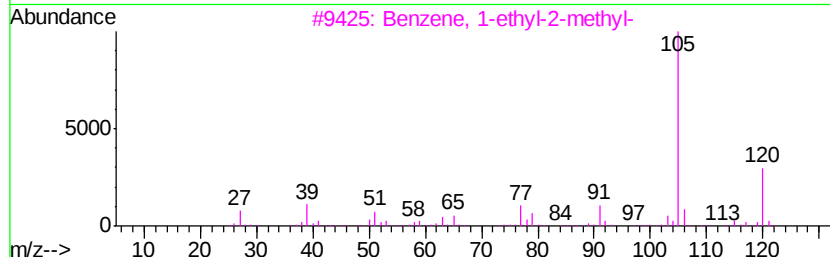
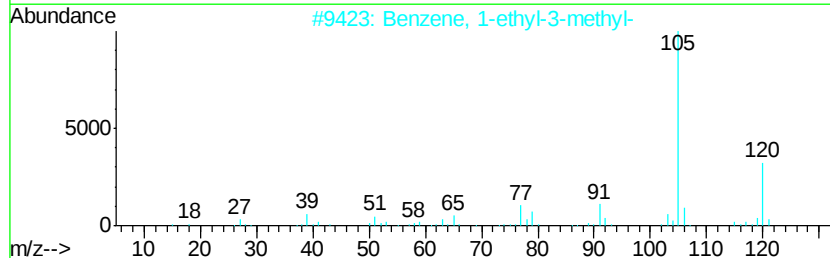
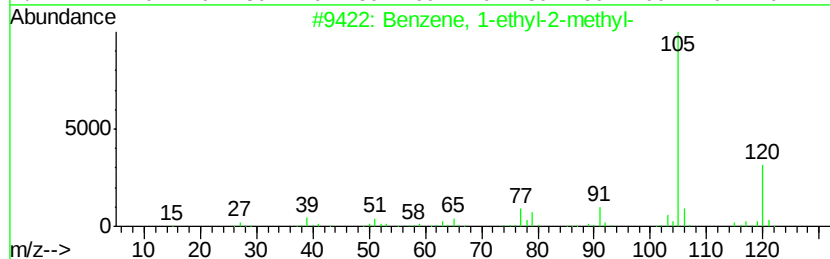
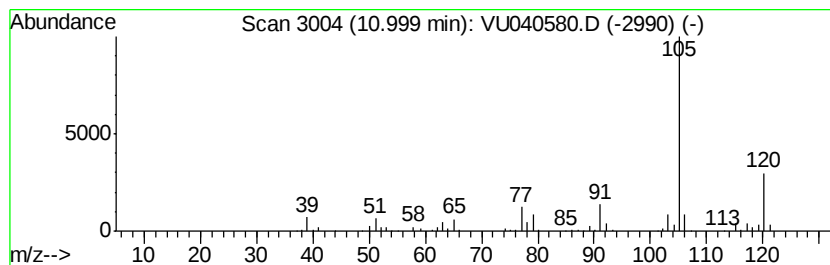
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR100820WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Benzene, 1-ethyl-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.00	3.36 ug/L	429699	1,4-Dichlorobenzene-d4	11.81

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12		000611-14-3	95
2	Benzene, 1-ethyl-3-methyl-	120	C9H12		000620-14-4	95
3	Benzene, 1-ethyl-2-methyl-	120	C9H12		000611-14-3	94
4	Benzene, 1-ethyl-3-methyl-	120	C9H12		000620-14-4	91
5	Benzene, 1-ethyl-3-methyl-	120	C9H12		000620-14-4	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100920\
Data File : VU040580.D
Acq On : 09 Oct 2020 13:48
Operator : SY/MD
Sample : L4240-11DL 40X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3T0DL

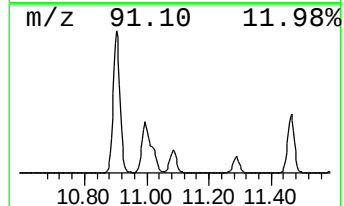
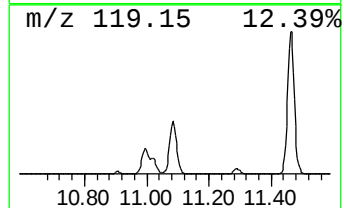
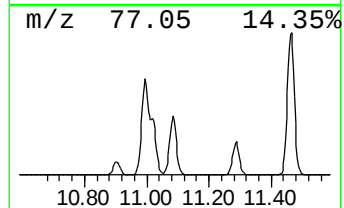
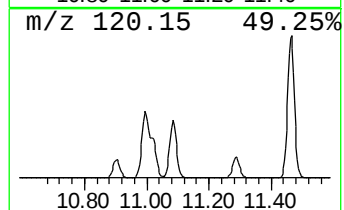
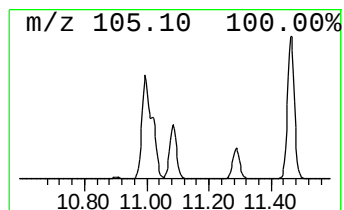
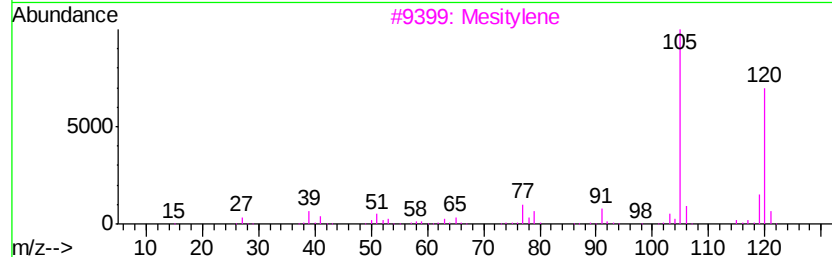
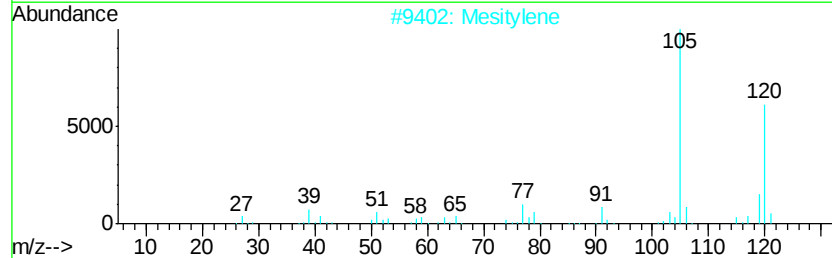
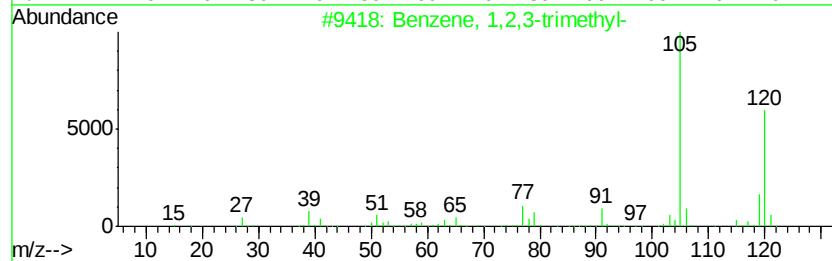
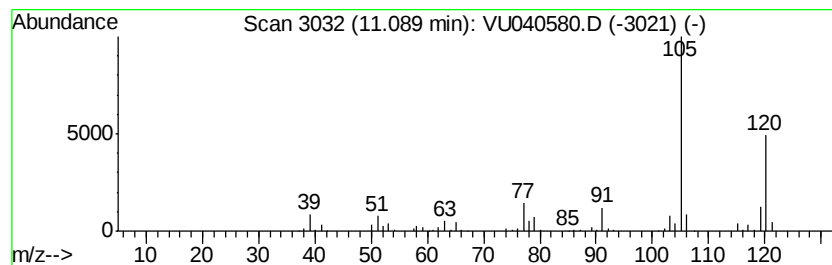
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR100820WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Benzene, 1,2,3-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.09	1.76 ug/L	225779	1,4-Dichlorobenzene-d4	11.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2		Mesitylene	120	C9H12	000108-67-8	97
3		Mesitylene	120	C9H12	000108-67-8	95
4		Mesitylene	120	C9H12	000108-67-8	95
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95



Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100920\
Data File : VU040580.D
Acq On : 09 Oct 2020 13:48
Operator : SY/MD
Sample : L4240-11DL 40X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3T0DL

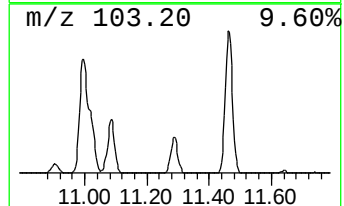
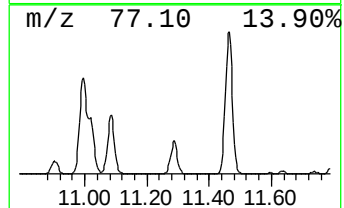
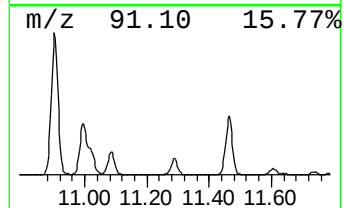
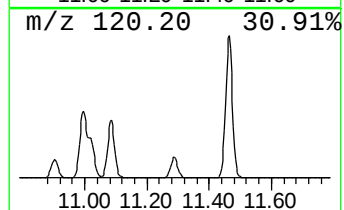
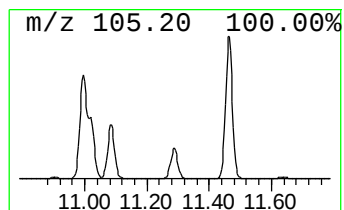
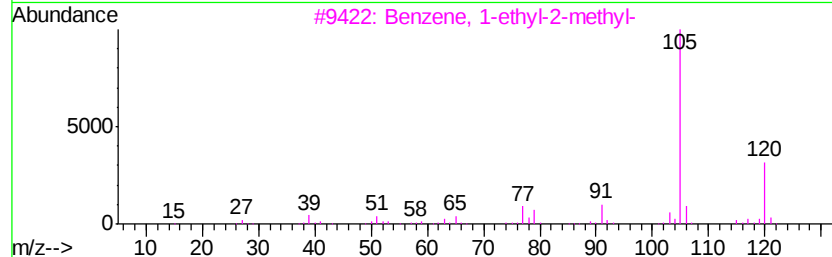
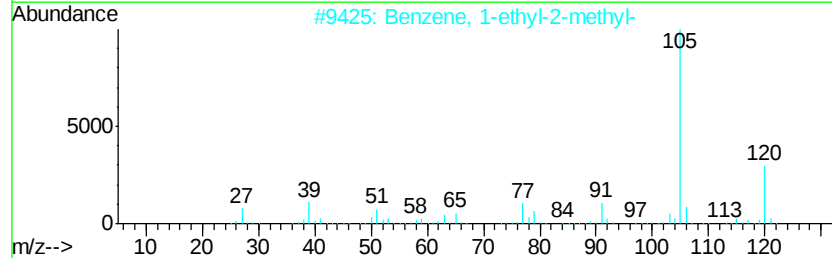
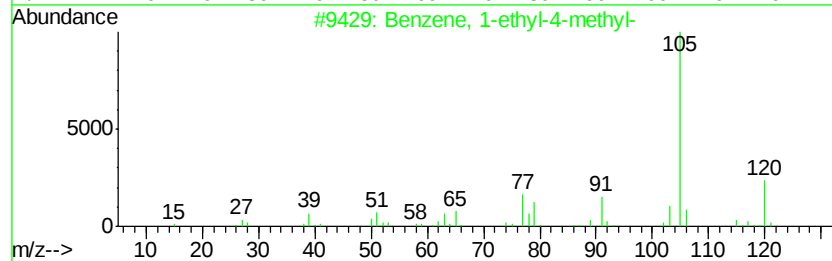
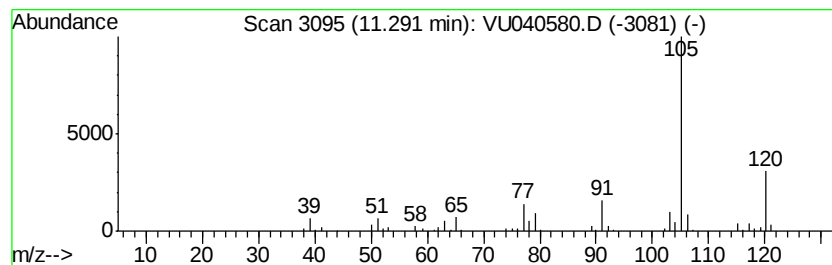
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR100820WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Benzene, 1-ethyl-4-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.29	0.90 ug/L	115083	1,4-Dichlorobenzene-d4	11.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
3		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5		Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100920\
Data File : VU040580.D
Acq On : 09 Oct 2020 13:48
Operator : SY/MD
Sample : L4240-11DL 40X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3T0DL

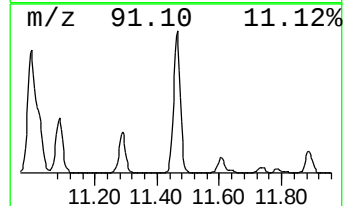
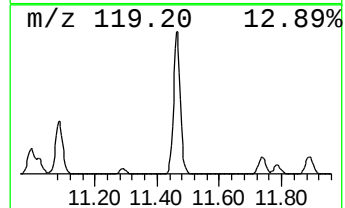
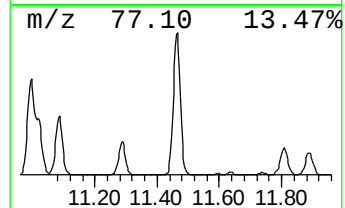
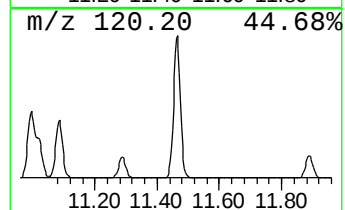
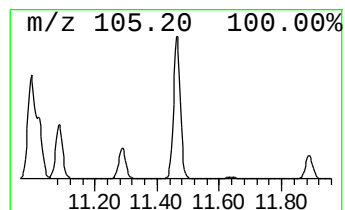
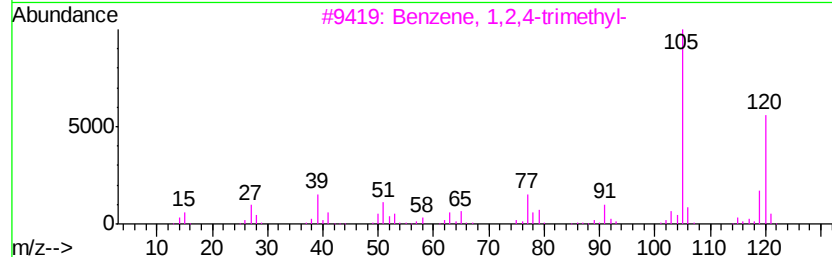
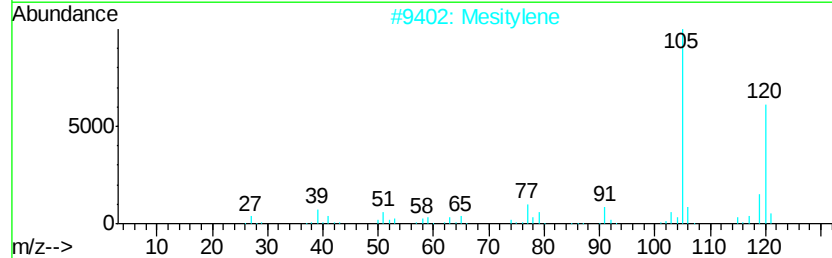
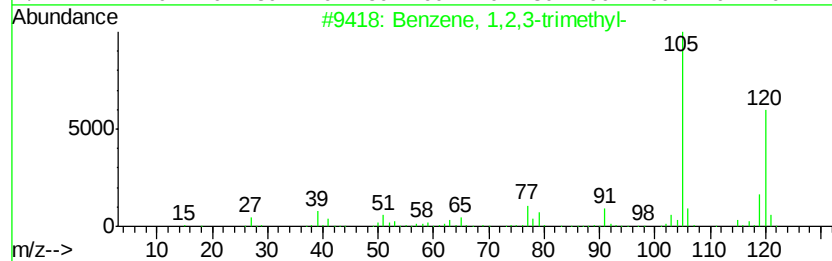
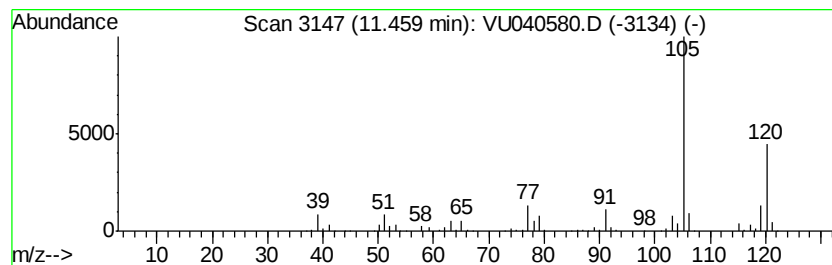
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR100820WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Mesitylene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.46	4.54 ug/L	581059	1,4-Dichlorobenzene-d4	11.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2		Mesitylene	120	C9H12	000108-67-8	97
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
4		Mesitylene	120	C9H12	000108-67-8	94
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100920\
Data File : VU040580.D
Acq On : 09 Oct 2020 13:48
Operator : SY/MD
Sample : L4240-11DL 40X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3T0DL

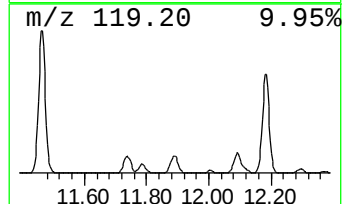
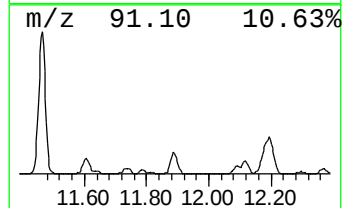
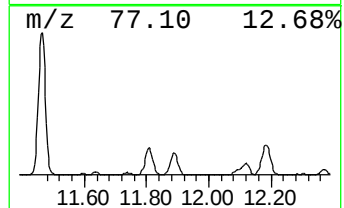
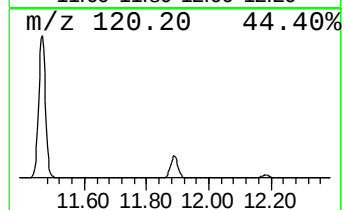
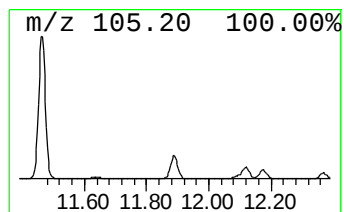
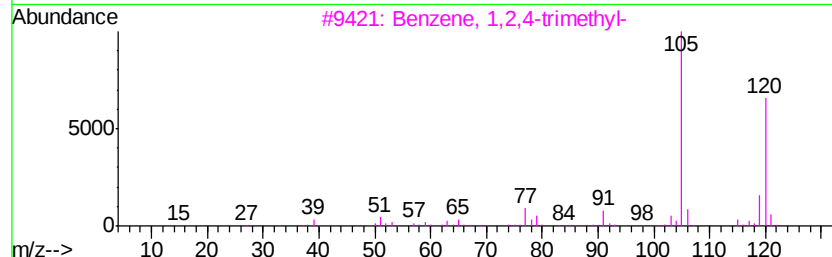
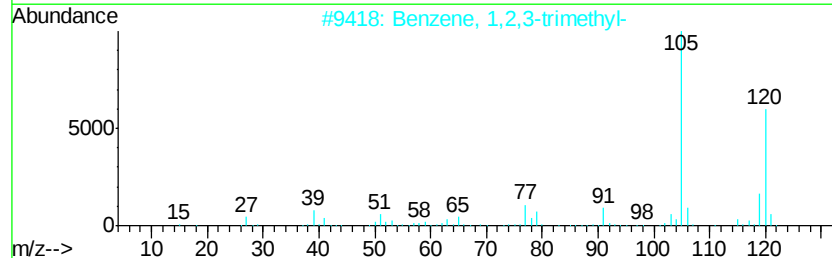
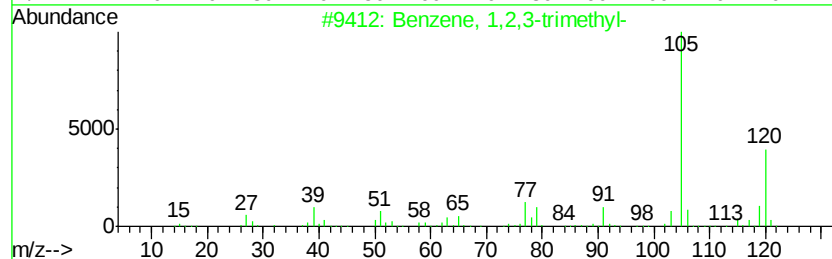
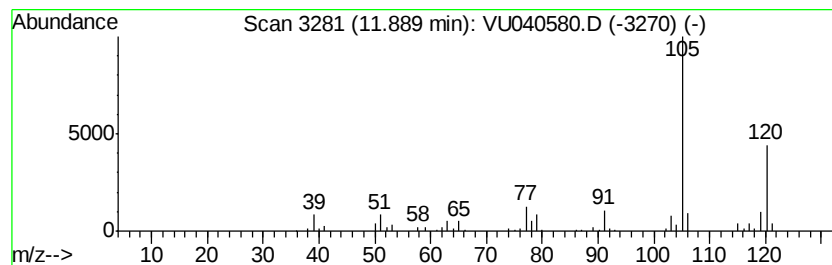
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR100820WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Benzene, 1,2,4-trimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.89	0.74 ug/L	94965	1,4-Dichlorobenzene-d4	11.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100920\
Data File : VU040580.D
Acq On : 09 Oct 2020 13:48
Operator : SY/MD
Sample : L4240-11DL 40X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3T0DL

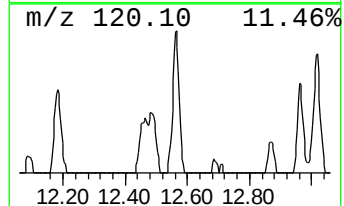
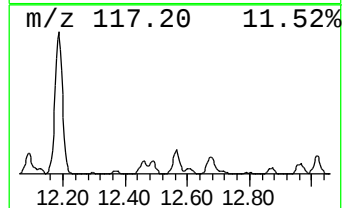
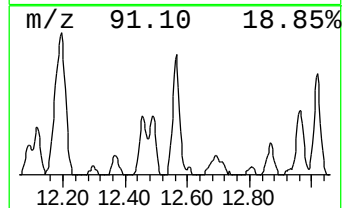
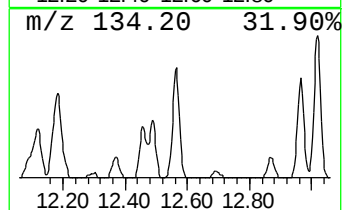
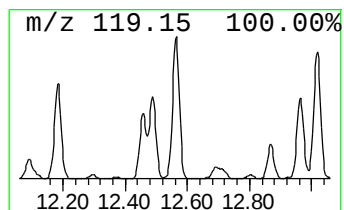
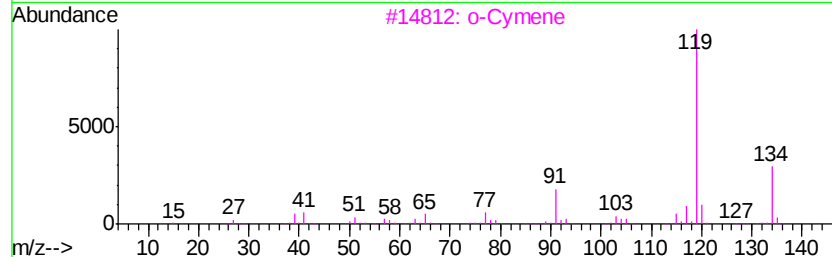
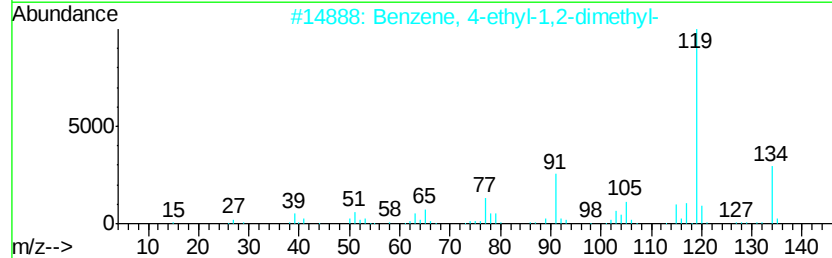
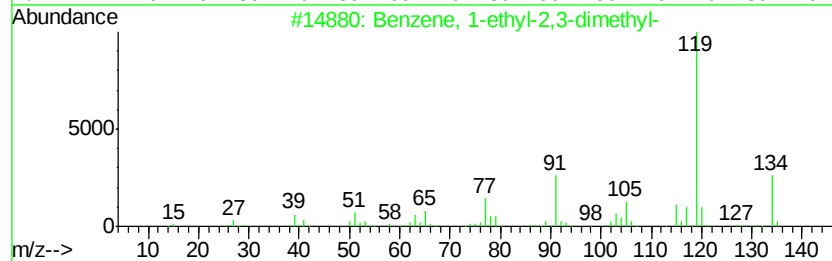
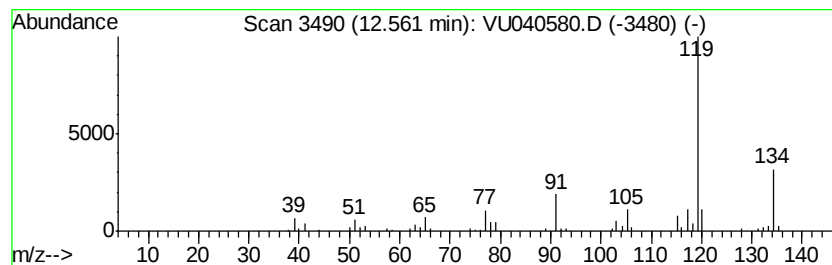
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR100820WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.56	0.59 ug/L	75642	1,4-Dichlorobenzene-d4	11.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3		o-Cymene	134	C10H14	000527-84-4	94
4		p-Cymene	134	C10H14	000099-87-6	93
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100920\
Data File : VU040580.D
Acq On : 09 Oct 2020 13:48
Operator : SY/MD
Sample : L4240-11DL 40X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3T0DL

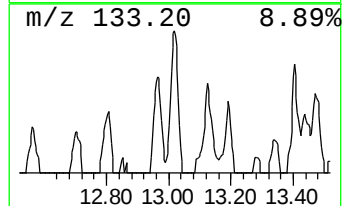
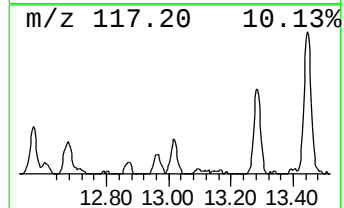
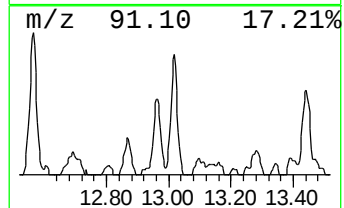
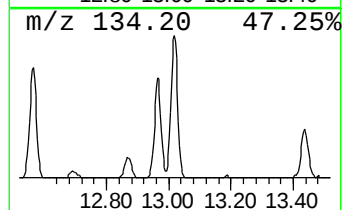
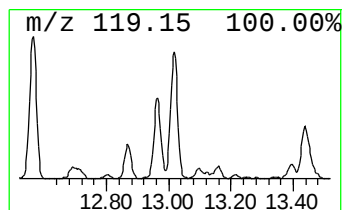
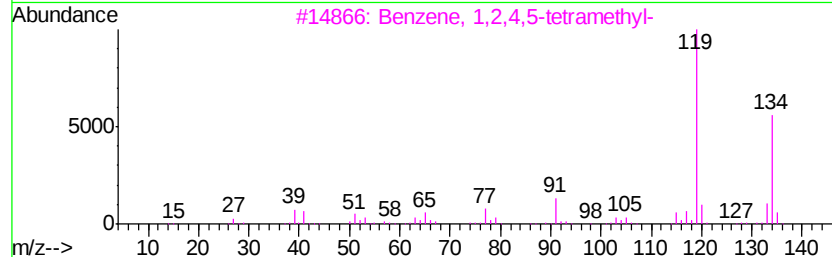
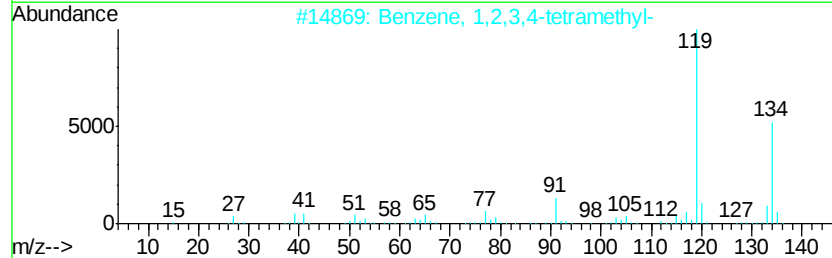
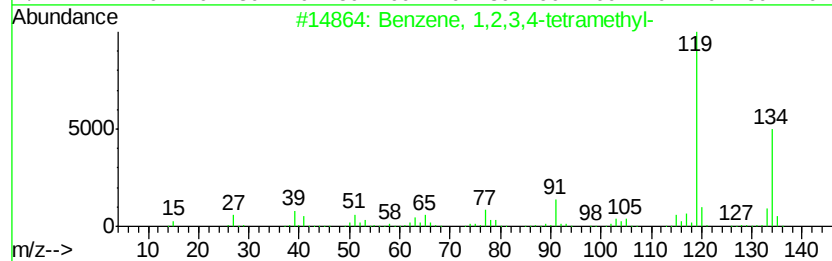
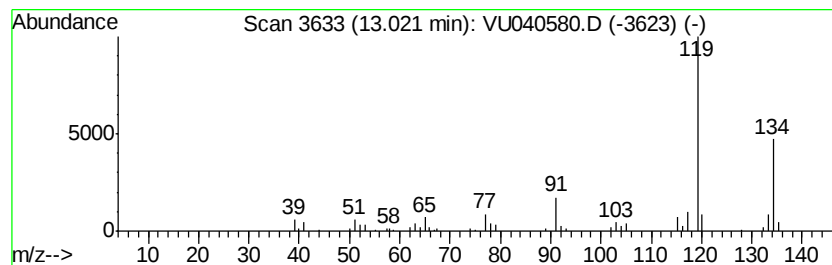
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR100820WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	0.55 ug/L	70412	1,4-Dichlorobenzene-d4	11.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	96
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
3		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
4		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
5		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100920\
Data File : VU040580.D
Acq On : 09 Oct 2020 13:48
Operator : SY/MD
Sample : L4240-11DL 40X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BG3T0DL

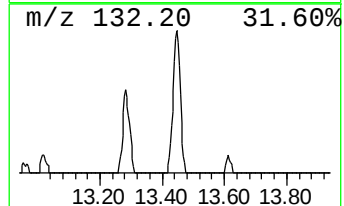
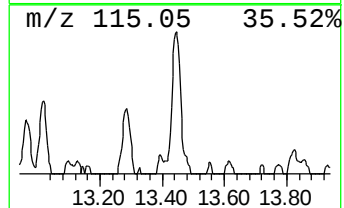
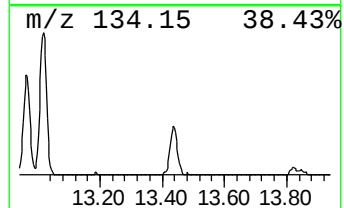
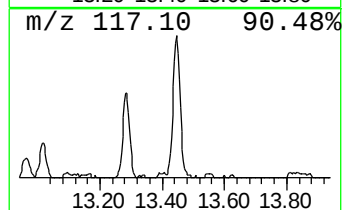
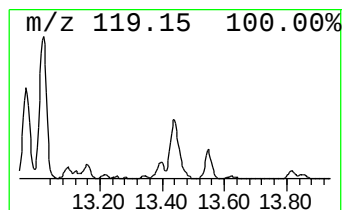
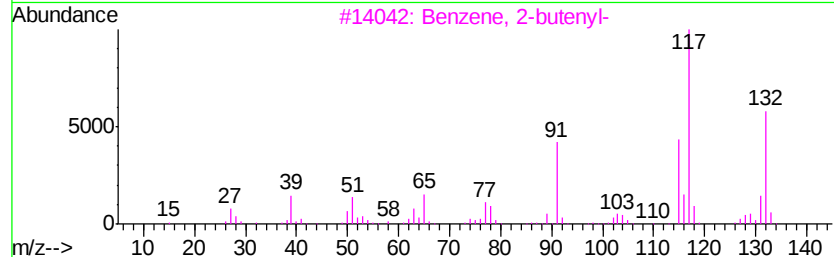
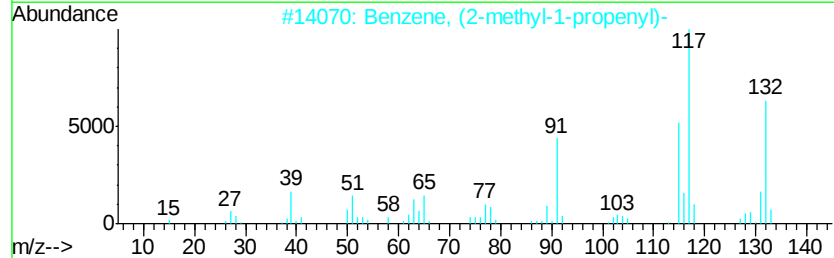
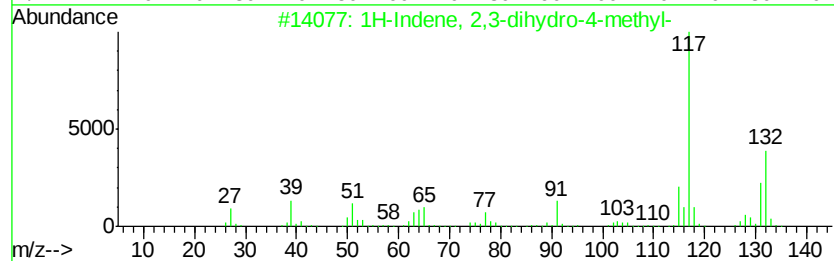
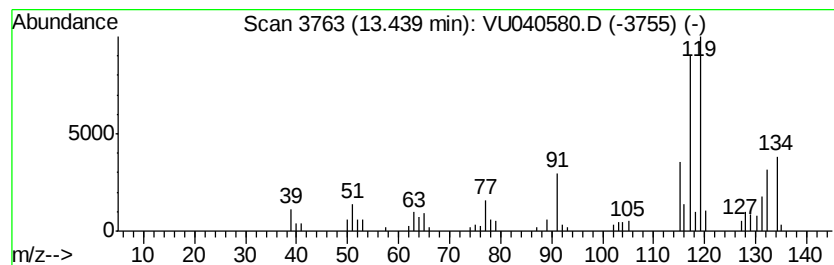
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR100820WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 unknown-01 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.44	0.50 ug/L	64499	1,4-Dichlorobenzene-d4	11.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	49
2			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	46
3			Benzene, 2-butenyl-	132	C10H12	001560-06-1	46
4			Benzene, 2-butenyl-	132	C10H12	001560-06-1	46
5			o-Cymene	134	C10H14	000527-84-4	46



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU100920\
 Data File : VU040580.D
 Acq On : 09 Oct 2020 13:48
 Operator : SY/MD
 Sample : L4240-11DL 40X
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BG3T0DL

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR100820WMA.M
 Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: Str...	3.10	10.8	ug/L	1025110	1	6.26	476088	5.0
(DEL) Alkane: Str...	3.38	18.2	ug/L	1734120	1	6.26	476088	5.0
(DEL) Alkane: Str...	3.72	45.4	ug/L	4323010	1	6.26	476088	5.0
(DEL) Alkane: Str...	4.31	1.5	ug/L	142561	1	6.26	476088	5.0
(DEL) Alkane: Str...	4.45	1.0	ug/L	95702	1	6.26	476088	5.0
(DEL) Alkane: Cyc...	4.55	17.3	ug/L	1649440	1	6.26	476088	5.0
Benzene, propyl-	10.90	0.9	ug/L	115588	3	11.81	639829	5.0
Benzene, 1-ethyl-...	11.00	3.4	ug/L	429699	3	11.81	639829	5.0
Benzene, 1,2,3-tr...	11.09	1.8	ug/L	225779	3	11.81	639829	5.0
Benzene, 1-ethyl-...	11.29	0.9	ug/L	115083	3	11.81	639829	5.0
Mesitylene	11.46	4.5	ug/L	581059	3	11.81	639829	5.0
Benzene, 1,2,4-tr...	11.89	0.7	ug/L	94965	3	11.81	639829	5.0
Benzene, 1-ethyl-...	12.56	0.6	ug/L	75642	3	11.81	639829	5.0
Benzene, 1,2,3,4-...	13.02	0.6	ug/L	70412	3	11.81	639829	5.0
unknown-01	13.44	0.5	ug/L	64499	3	11.81	639829	5.0