

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100520\
 Data File : VX018770.D
 Acq On : 05 Oct 2020 17:03
 Operator : JC/SP
 Sample : L4239-05DL2 400X
 Misc : 25.0mL/MSVOA X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BG3N5DL2

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_X\METHOD\SOMXTR100220WMA.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	1.392	46	50	56	rBV	71773	76488	8.34%	0.795%
2	1.697	95	100	107	rBV	58580	74801	8.16%	0.778%
3	2.343	200	206	214	rBV	133108	210665	22.97%	2.190%
4	2.873	278	293	302	rVB3	21366	63837	6.96%	0.664%
5	3.154	329	339	348	rBV	42357	102233	11.15%	1.063%
6	3.501	388	396	406	rBV	77470	175910	19.18%	1.829%
7	3.672	418	424	430	rVB5	4474	11522	1.26%	0.120%
8	4.373	530	539	551	rVB2	53566	155452	16.95%	1.616%
9	4.568	559	571	584	rBV2	101256	376873	41.10%	3.918%
10	5.141	654	665	678	rBV	86170	264958	28.89%	2.754%
11	5.464	705	718	727	rBV2	58782	188438	20.55%	1.959%
12	5.550	727	732	737	rVB6	4344	10669	1.16%	0.111%
13	6.044	800	813	829	rBV3	207092	611704	66.71%	6.359%
14	6.830	931	942	957	rBV	148541	353517	38.55%	3.675%
15	7.373	1020	1031	1044	rBV	127998	285740	31.16%	2.970%
16	8.372	1188	1195	1205	rBV	84476	138227	15.07%	1.437%
17	8.696	1241	1248	1254	rBV	335798	506225	55.21%	5.263%
18	8.763	1254	1259	1268	rVB	582627	916982	100.00%	9.533%
19	8.994	1292	1297	1305	rBV	62236	89423	9.75%	0.930%
20	9.427	1362	1368	1379	rBV	427336	631247	68.84%	6.562%
21	9.738	1411	1419	1426	rBV	23066	34465	3.76%	0.358%
22	10.098	1472	1478	1486	rVB	305998	419917	45.79%	4.365%
23	10.238	1496	1501	1505	rBV	60827	82872	9.04%	0.862%
24	10.342	1512	1518	1531	rVB	227474	302801	33.02%	3.148%
25	10.683	1568	1574	1579	rBV	115954	151097	16.48%	1.571%
26	11.000	1622	1626	1632	rVB	24547	32786	3.58%	0.341%
27	11.232	1658	1664	1673	rVB	142967	181096	19.75%	1.883%
28	11.347	1676	1683	1689	rBV	63903	85799	9.36%	0.892%
29	11.421	1689	1695	1702	rVV	292593	498420	54.35%	5.181%
30	11.488	1702	1706	1715	rVB	137871	177900	19.40%	1.849%
31	11.610	1720	1726	1729	rVV	267257	327273	35.69%	3.402%
32	11.652	1729	1733	1739	rVB	149129	184565	20.13%	1.919%
33	11.793	1750	1756	1763	rBV	498166	596925	65.10%	6.206%
34	11.866	1763	1768	1772	rBV4	7669	10698	1.17%	0.111%

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BG3N5DL2

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_X\METHOD\SOMXTR100220WMA.M
Title : TRACE VOA SOM01.0

35	12.061	1795	1800	1806	rBV	332184	416947	45.47%	4.334%
36	12.128	1806	1811	1819	rVB	134877	164129	17.90%	1.706%
37	12.280	1830	1836	1844	rBV	55363	82720	9.02%	0.860%
38	12.360	1844	1849	1857	rVV	392310	504205	54.99%	5.242%
39	12.573	1880	1884	1886	rBV2	7495	11500	1.25%	0.120%
40	12.652	1891	1897	1901	rBV	13049	16068	1.75%	0.167%
41	12.890	1931	1936	1943	rBV5	7416	10006	1.09%	0.104%
42	12.963	1943	1948	1951	rVV	8138	10163	1.11%	0.106%
43	13.000	1951	1954	1958	rVB	12670	14759	1.61%	0.153%
44	13.329	2003	2008	2013	rVB2	11830	18306	2.00%	0.190%
45	13.817	2082	2088	2094	rBV	29527	38955	4.25%	0.405%

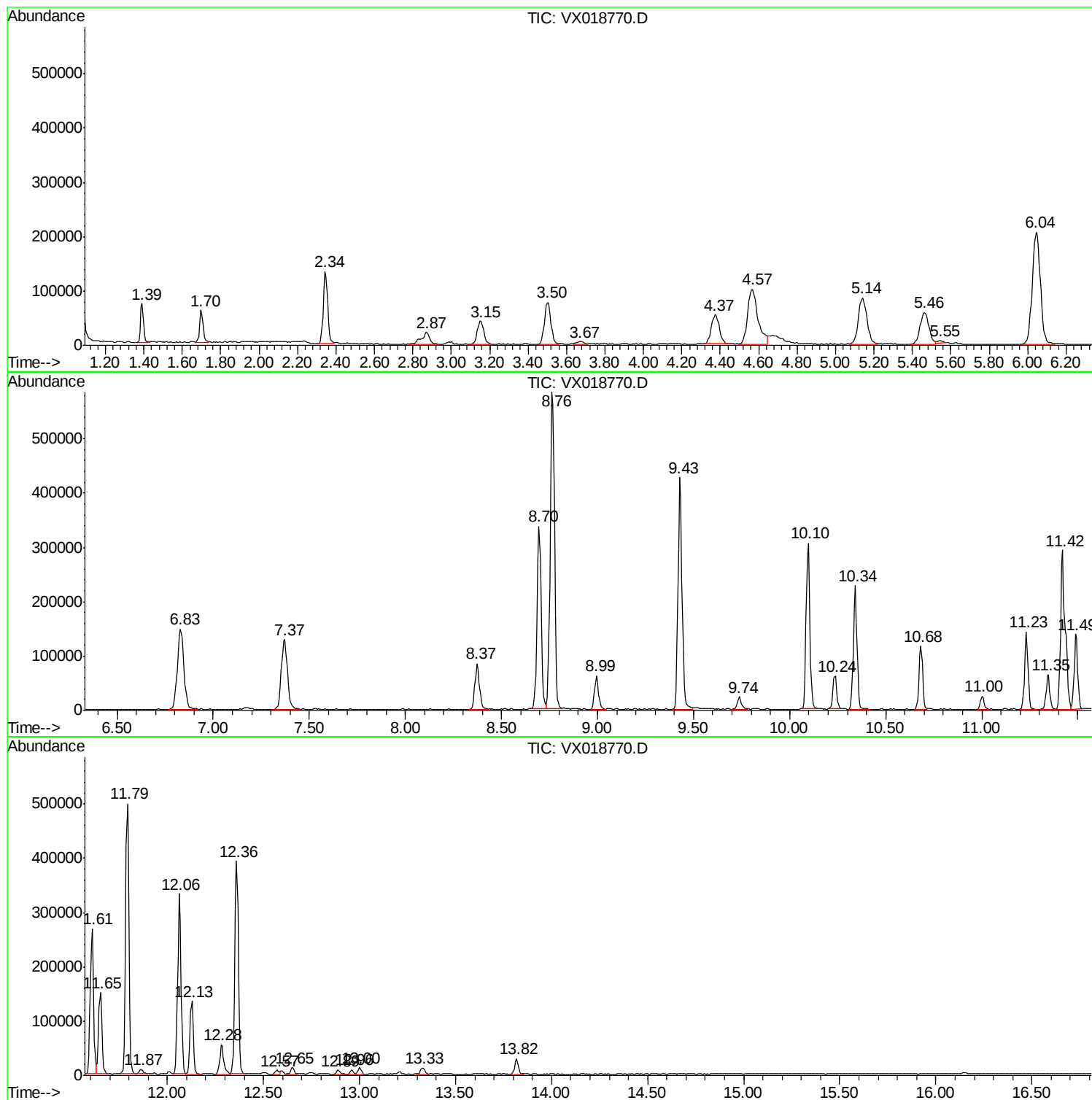
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

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



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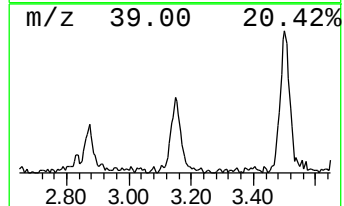
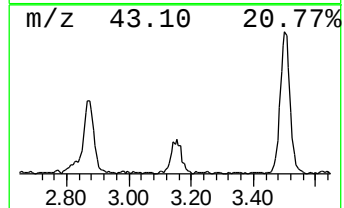
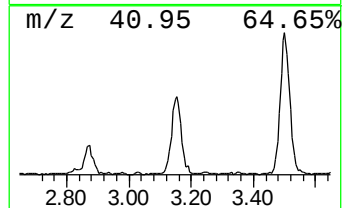
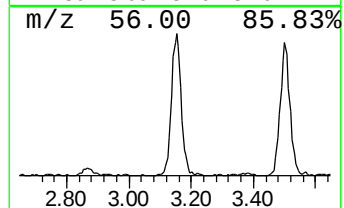
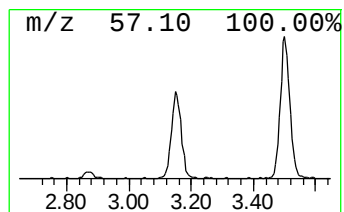
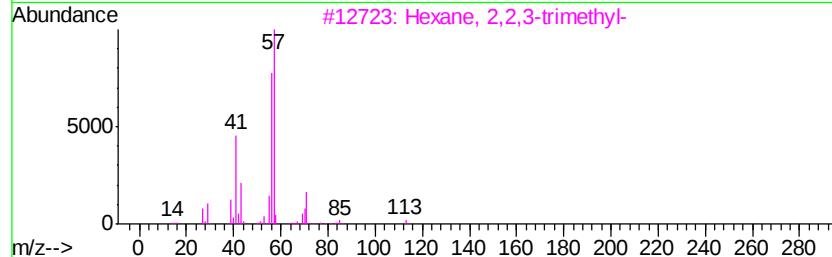
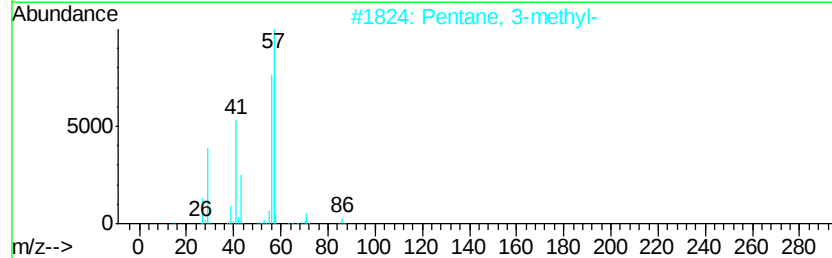
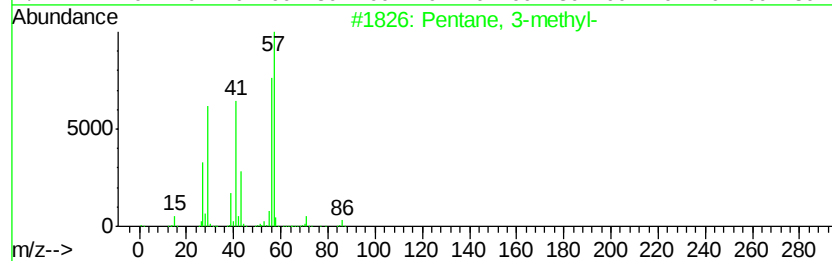
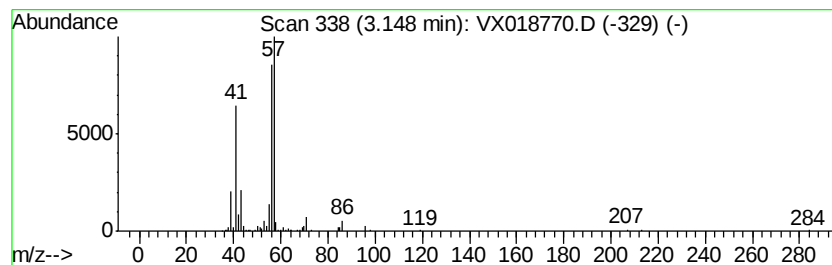
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.15	1.45 ug/L	102233	1,4-Difluorobenzene	6.83

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 3-methyl-	86	C6H14	000096-14-0	72
2			Pentane, 3-methyl-	86	C6H14	000096-14-0	72
3			Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	64
4			Decane, 2,2,3-trimethyl-	184	C13H28	062338-09-4	56
5			Pentane, 2,2,3-trimethyl-	114	C8H18	000564-02-3	50



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Instrument :
MSVOA_X
ClientSampleId :
BG3N5DL2

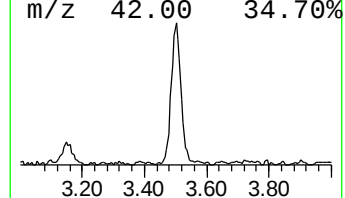
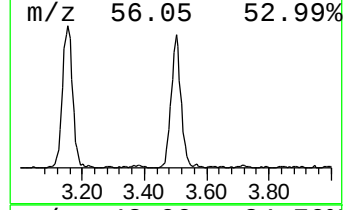
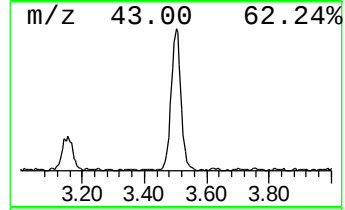
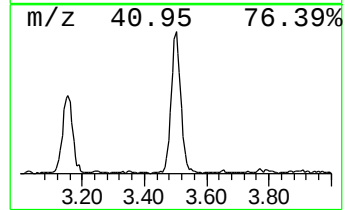
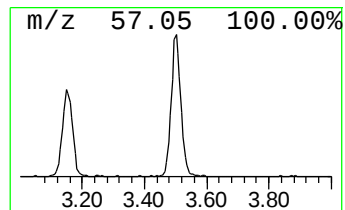
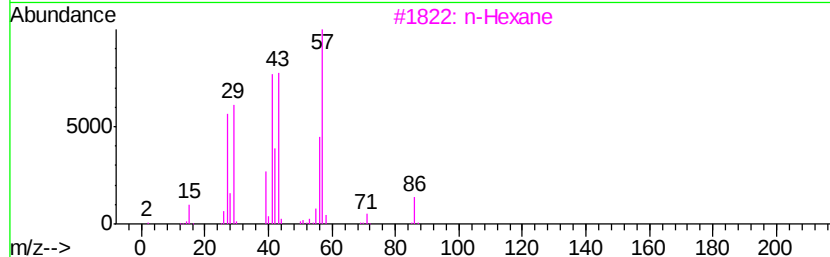
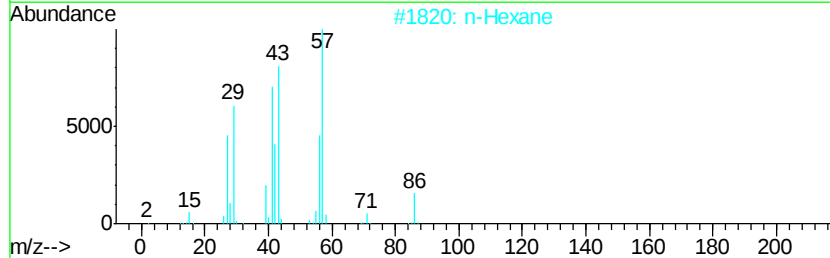
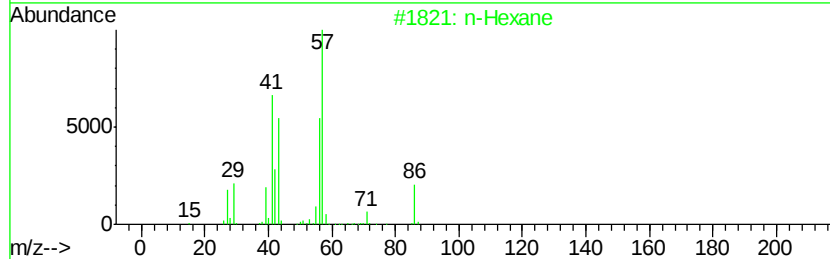
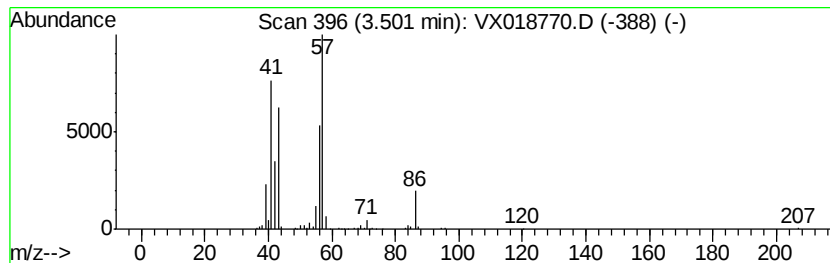
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.50	2.49 ug/L	175910	1,4-Difluorobenzene	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexane	86	C6H14	000110-54-3	90
2		n-Hexane	86	C6H14	000110-54-3	59
3		n-Hexane	86	C6H14	000110-54-3	59
4		Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	47
5		Hydroperoxide, hexyl	118	C6H14O2	004312-76-9	47



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100520\
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Sample : L4239-05DL2 400X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3N5DL2

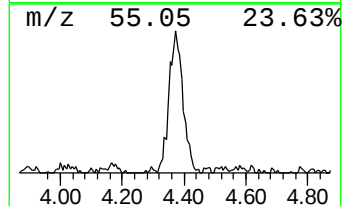
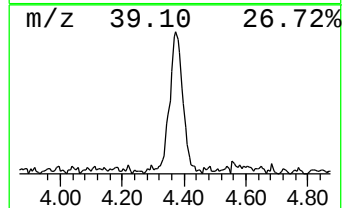
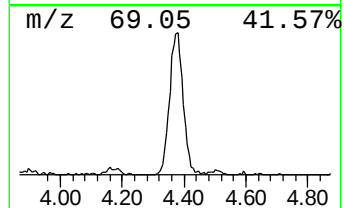
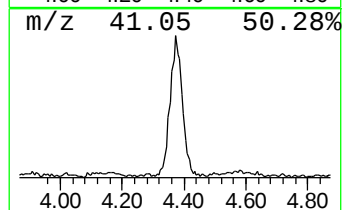
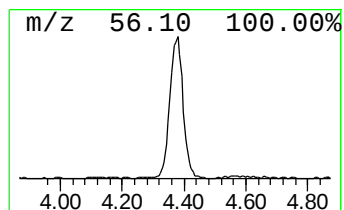
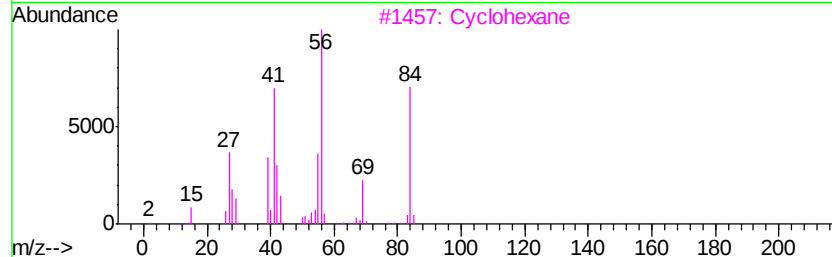
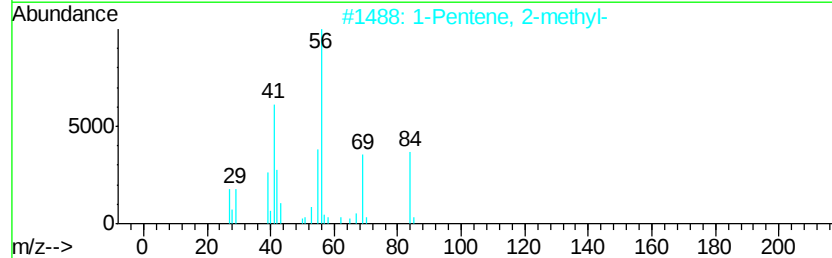
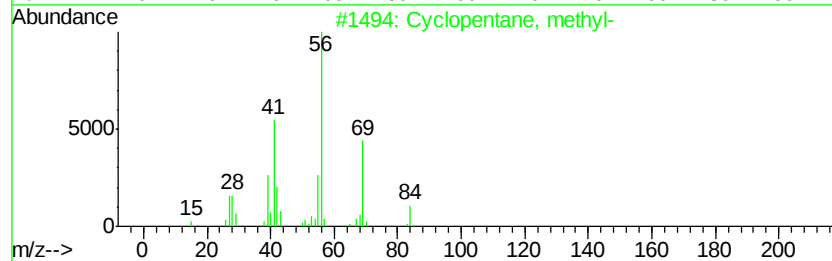
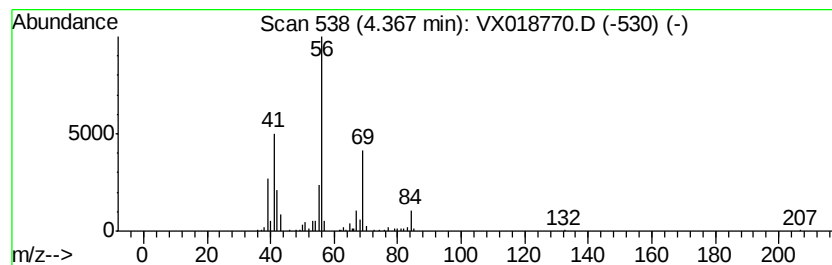
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 3 (DEL) Alkane: Cyclic4.37 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.37	2.20 ug/L	155452	1,4-Difluorobenzene	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	90
2		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	80
3		Cyclohexane	84	C6H12	000110-82-7	80
4		Cyclopentane, methyl-	84	C6H12	000096-37-7	80
5		Cyclohexane	84	C6H12	000110-82-7	72



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ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
BG3N5DL2

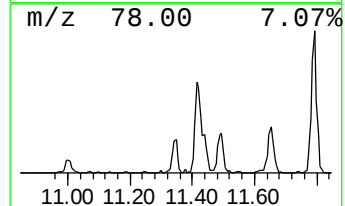
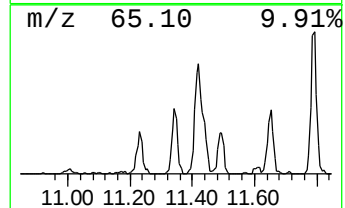
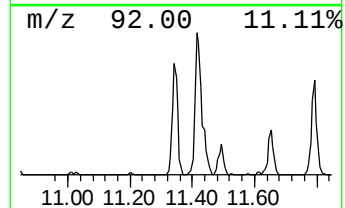
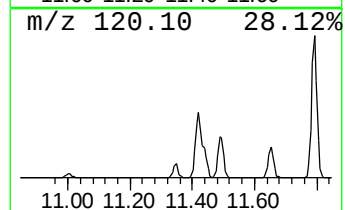
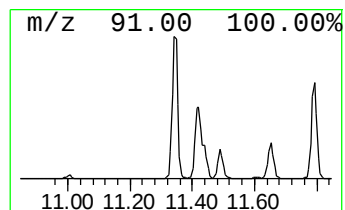
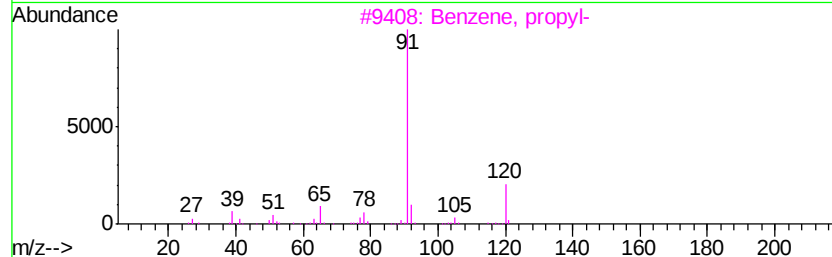
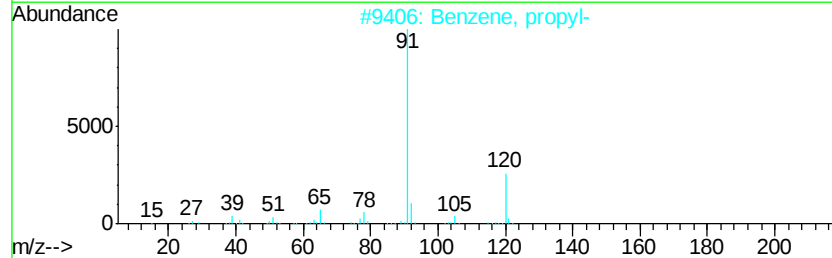
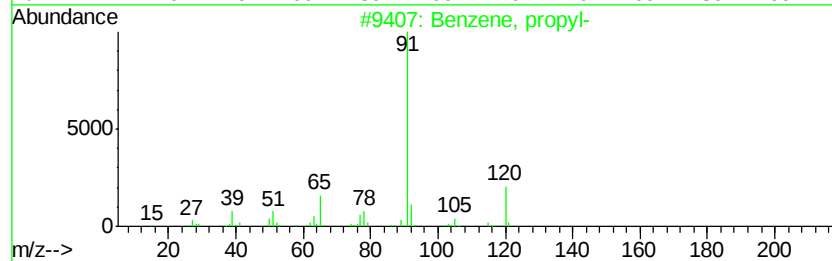
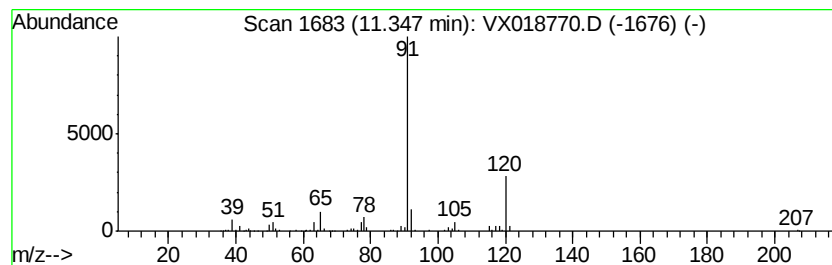
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, propyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.35	1.03 ug/L	85799	1,4-Dichlorobenzene-d4	12.06

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		N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	72
5		n-Butylbenzylamine	163	C11H17N	002403-22-7	72



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ALS Vial : 11 Sample Multiplier: 1

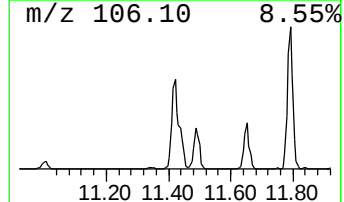
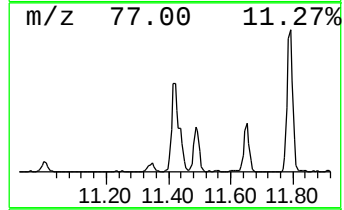
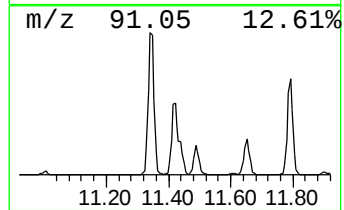
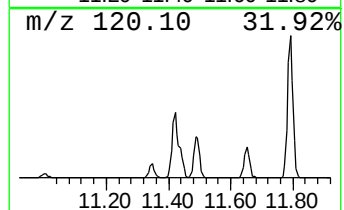
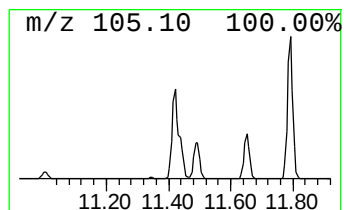
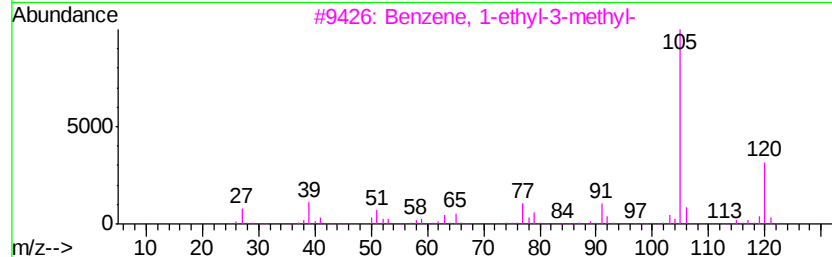
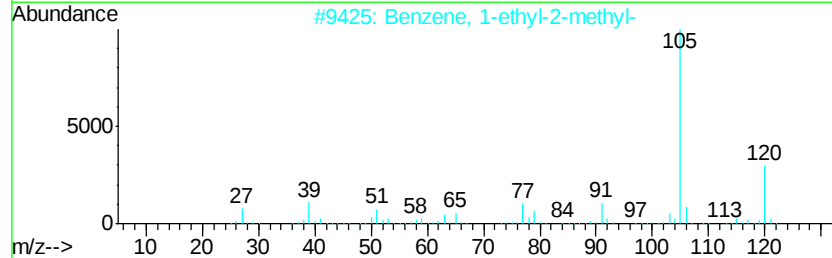
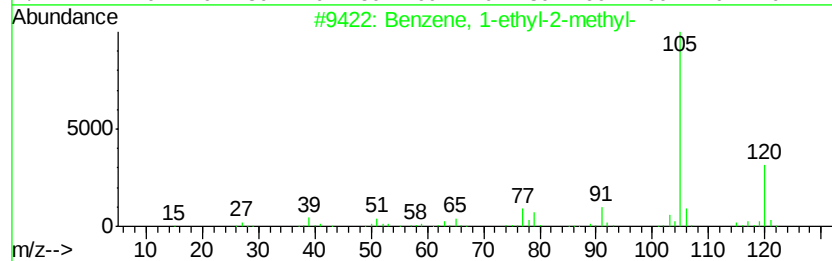
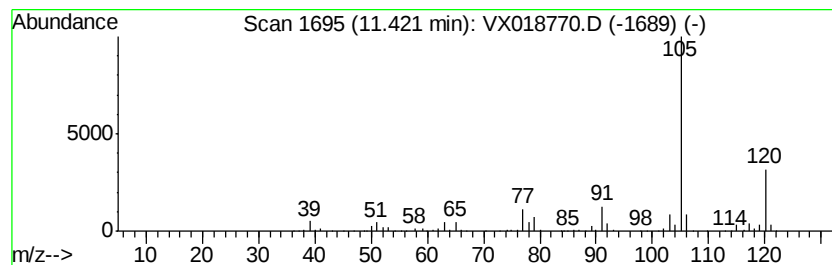
Instrument :
MSVOA_X
ClientSampled :
BG3N5DL2

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Quant Title : TRACE VOA SOM01.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.42	5.98 ug/L	498420	1,4-Dichlorobenzene-d4	12.06	
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-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:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100520\
Data File : VX018770.D
Acq On : 05 Oct 2020 17:03
Operator : JC/SP
Sample : L4239-05DL2 400X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 11 Sample Multiplier: 1

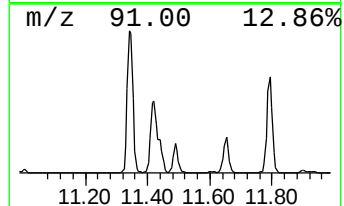
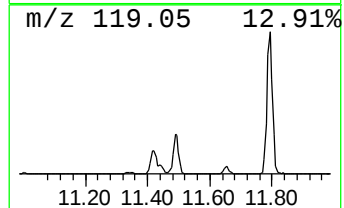
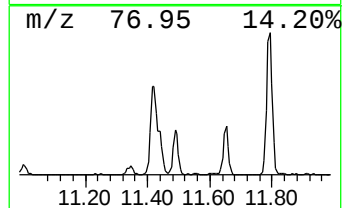
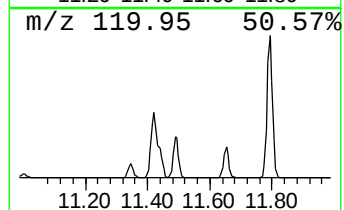
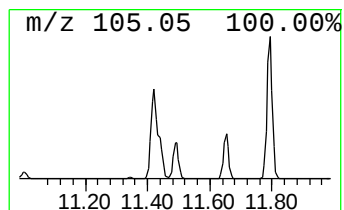
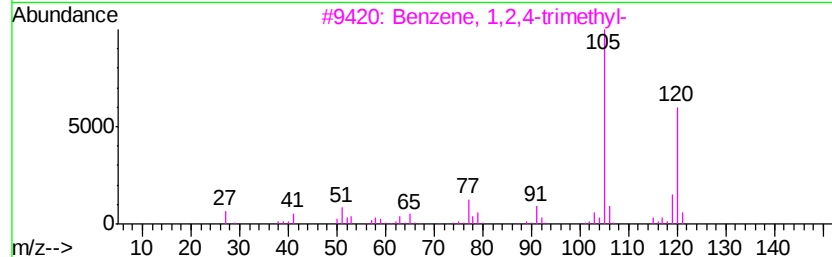
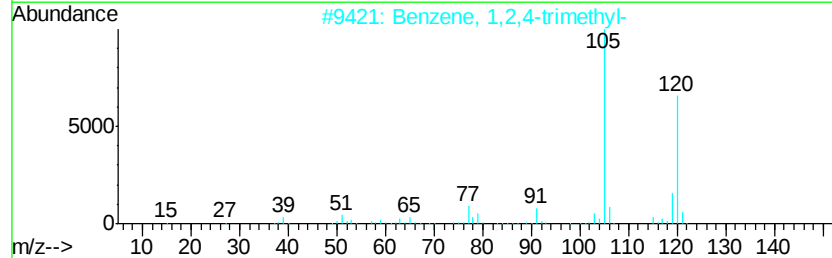
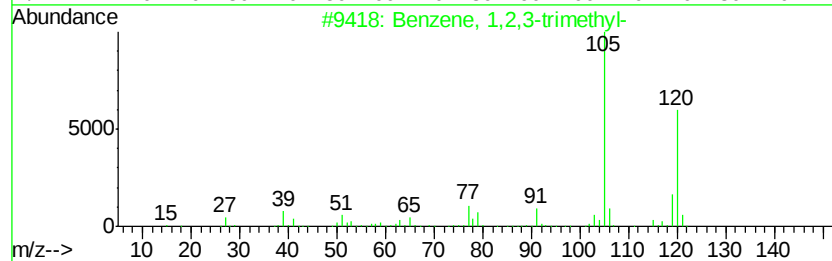
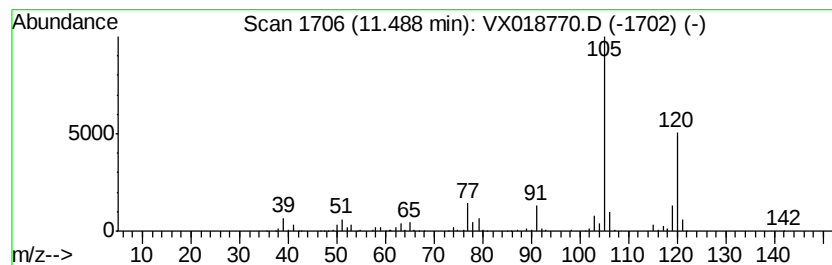
Instrument :
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ClientSampleId :
BG3N5DL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.49	2.13 ug/L	177900	1,4-Dichlorobenzene-d4	12.06	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
4	Mesitylene	120	C9H12	000108-67-8	94
5	Mesitylene	120	C9H12	000108-67-8	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100520\
Data File : VX018770.D
Acq On : 05 Oct 2020 17:03
Operator : JC/SP
Sample : L4239-05DL2 400X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BG3N5DL2

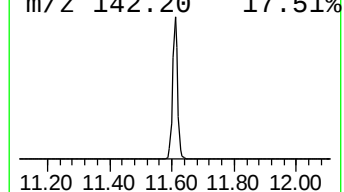
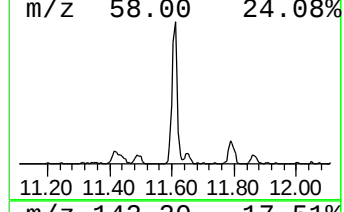
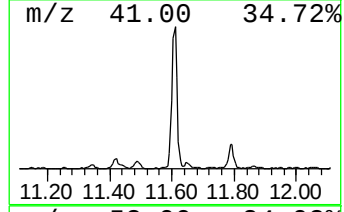
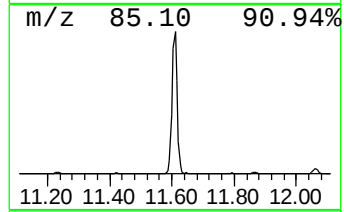
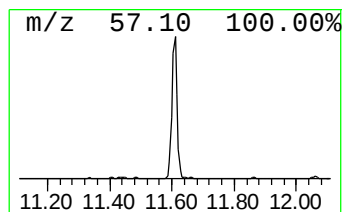
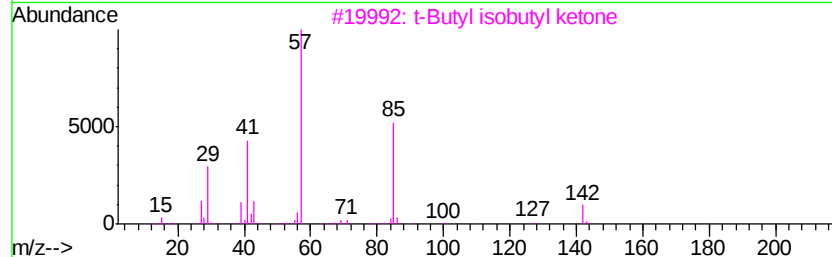
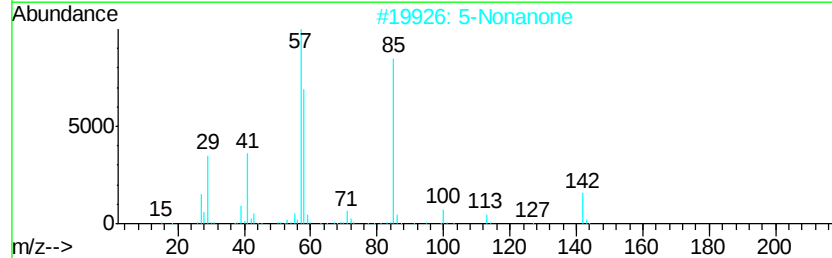
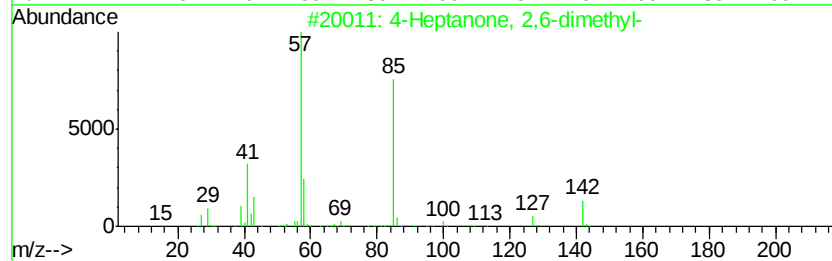
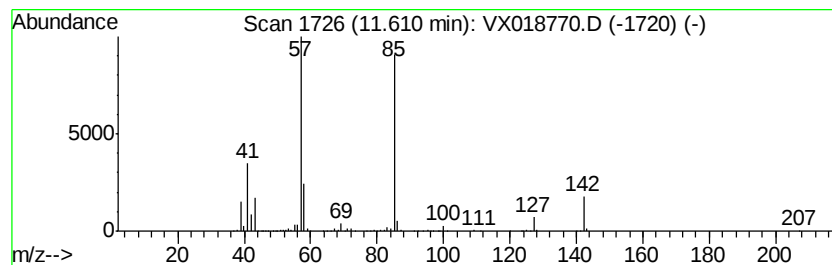
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 8 4-Heptanone, 2,6-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.61	3.92 ug/L	327273	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Heptanone, 2,6-dimethyl-	142	C9H18O	000108-83-8	91
2		5-Nonanone	142	C9H18O	000502-56-7	59
3		t-Butyl isobutyl ketone	142	C9H18O	014705-50-1	59
4		4-Heptanone, 2,6-dimethyl-	142	C9H18O	000108-83-8	50
5		3-Buten-2-ol, 2,3-dimethyl-	100	C6H12O	010473-13-9	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100520\
Data File : VX018770.D
Acq On : 05 Oct 2020 17:03
Operator : JC/SP
Sample : L4239-05DL2 400X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
BG3N5DL2

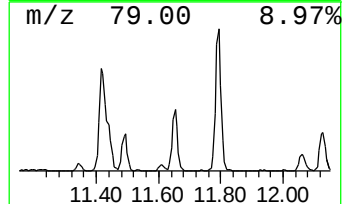
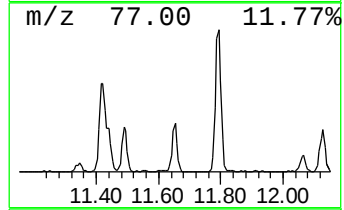
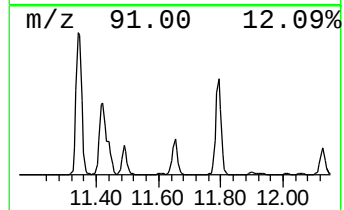
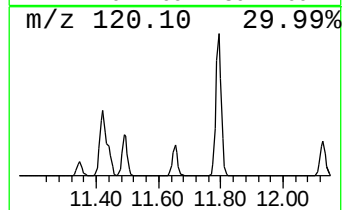
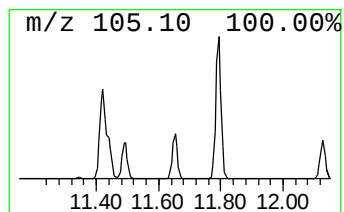
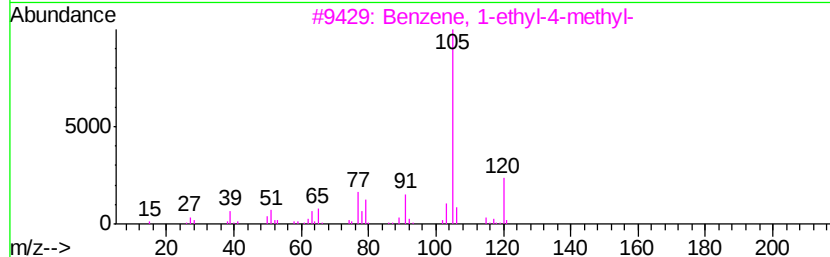
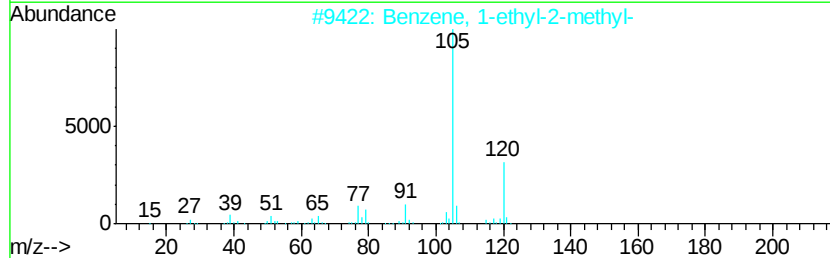
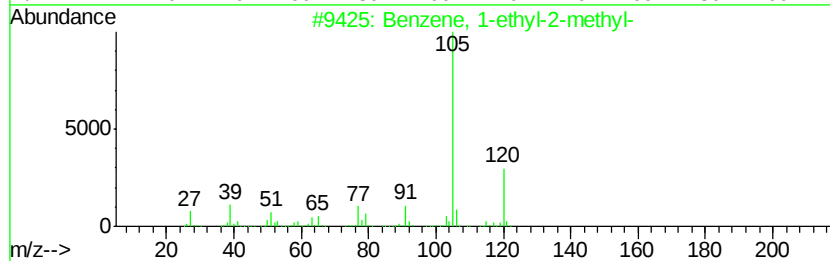
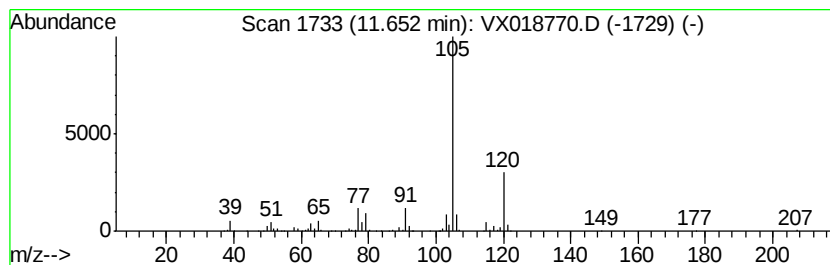
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 9 Benzene, 1-ethyl-4-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.65	2.21 ug/L	184565	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100520\
Data File : VX018770.D
Acq On : 05 Oct 2020 17:03
Operator : JC/SP
Sample : L4239-05DL2 400X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
BG3N5DL2

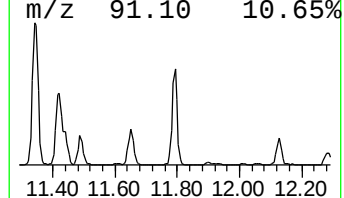
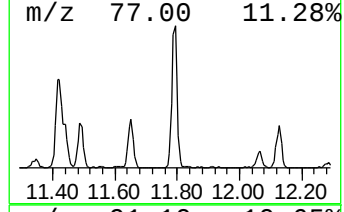
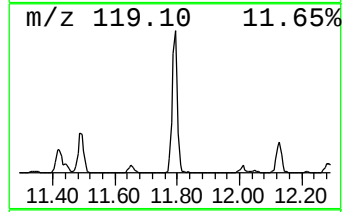
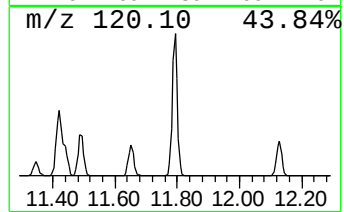
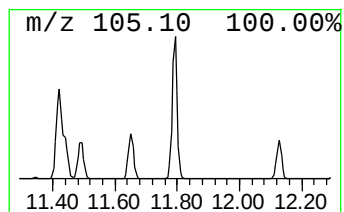
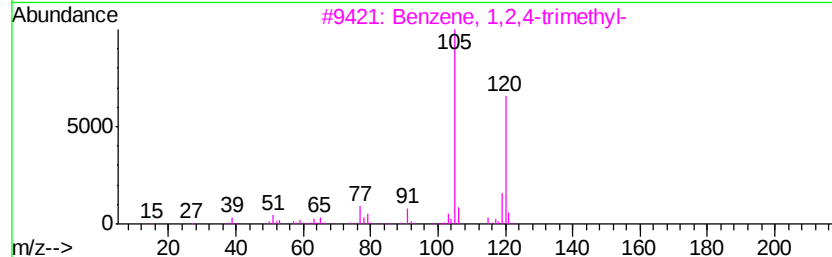
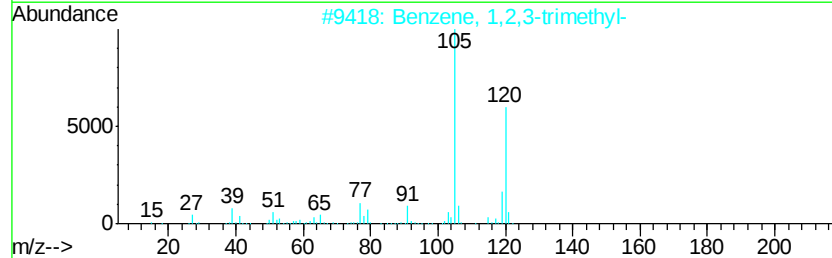
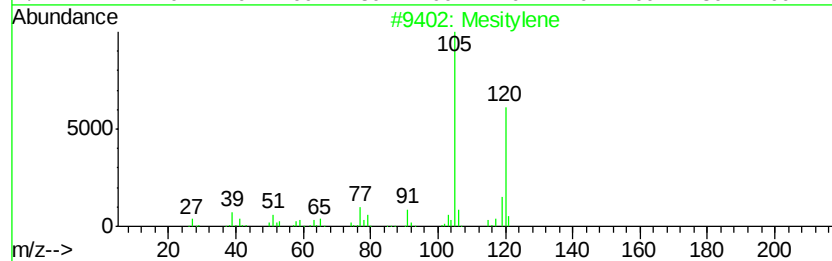
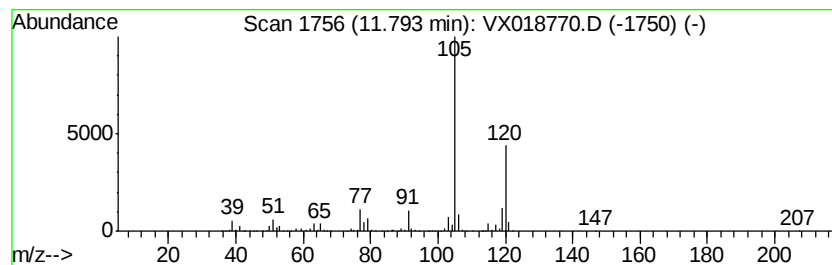
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 10 Mesitylene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.79	7.16 ug/L	596925	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Mesitylene	120	C9H12	000108-67-8	97
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	93
5		Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100520\
Data File : VX018770.D
Acq On : 05 Oct 2020 17:03
Operator : JC/SP
Sample : L4239-05DL2 400X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
BG3N5DL2

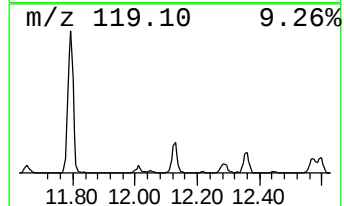
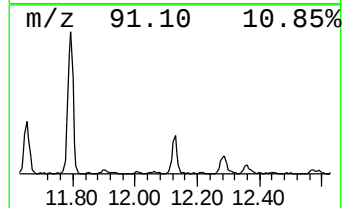
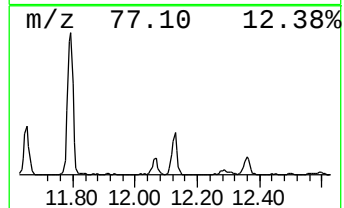
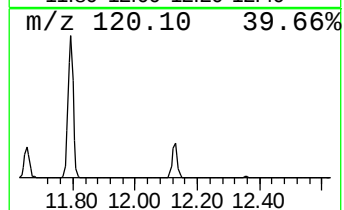
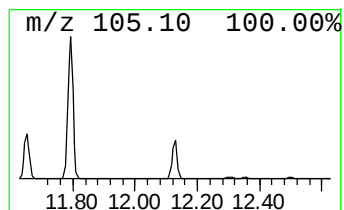
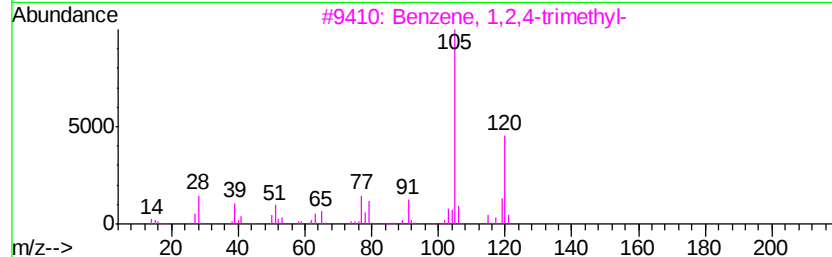
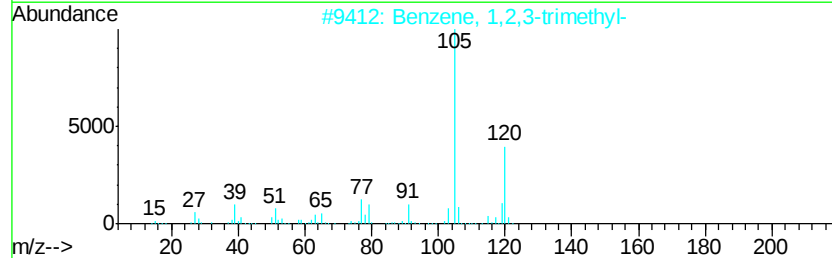
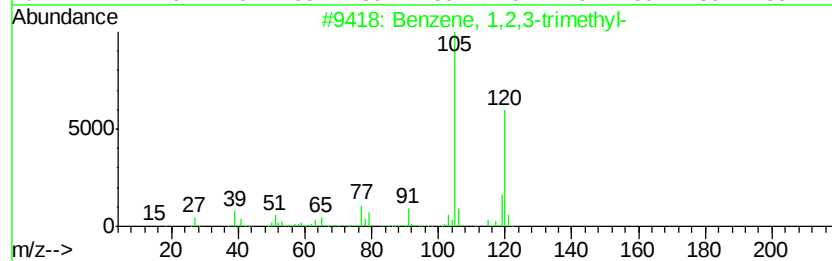
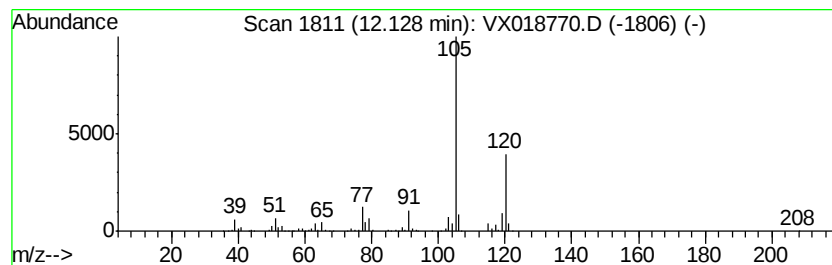
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 11 Benzene, 1,2,4-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.13	1.97 ug/L	164129	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4		Mesitylene	120	C9H12	000108-67-8	91
5		Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX100520\
Data File : VX018770.D
Acq On : 05 Oct 2020 17:03
Operator : JC/SP
Sample : L4239-05DL2 400X
Misc : 25.0mL/MSVOA X/WATER
ALS Vial : 11 Sample Multiplier: 1

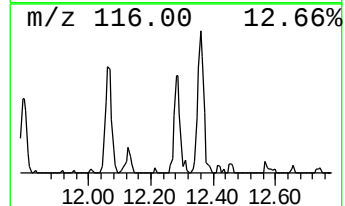
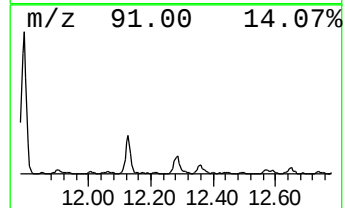
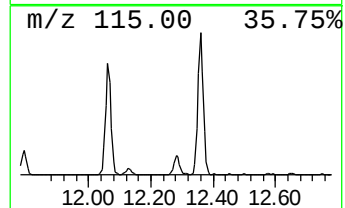
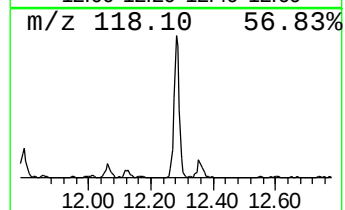
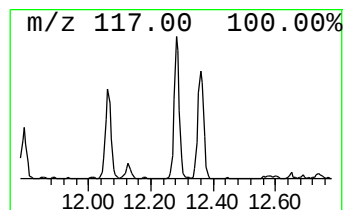
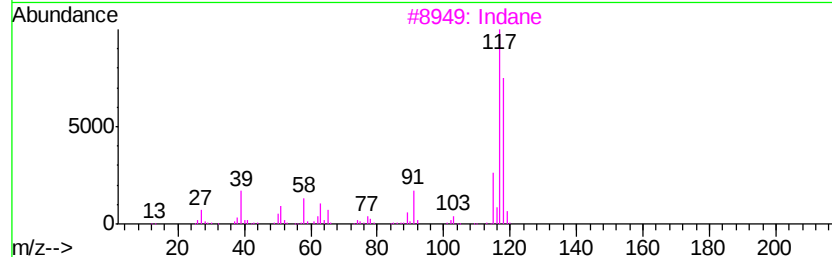
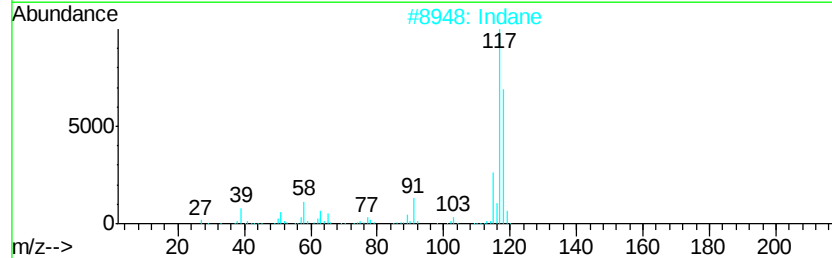
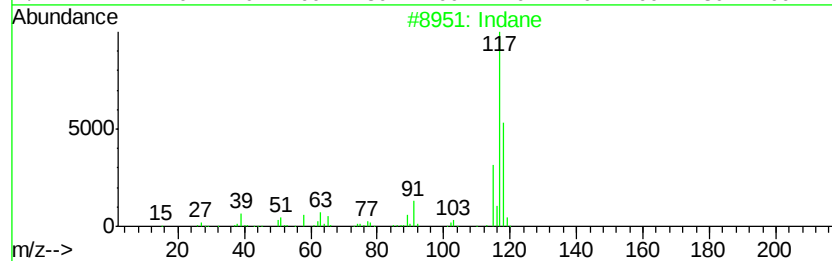
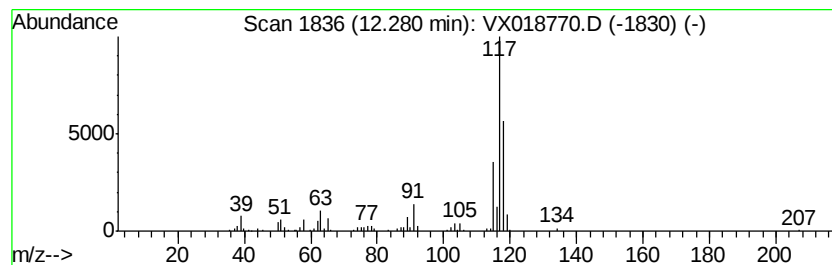
Instrument :
MSVOA_X
ClientSampled :
BG3N5DL2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 12 Indane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.28	0.99 ug/L	82720	1,4-Dichlorobenzene-d4	12.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Indane		118 C9H10	000496-11-7 94
2	Indane		118 C9H10	000496-11-7 94
3	Indane		118 C9H10	000496-11-7 91
4	Benzene, 1-ethenyl-2-methyl-		118 C9H10	000611-15-4 81
5	Benzene, 1-ethenyl-4-methyl-		118 C9H10	000622-97-9 74



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX100520\
 Data File : VX018770.D
 Acq On : 05 Oct 2020 17:03
 Operator : JC/SP
 Sample : L4239-05DL2 400X
 Misc : 25.0mL/MSVOA_X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BG3N5DL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\SOMXTR100220WMA.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...	3.15	1.5	ug/L	102233	1	6.83	353517	5.0
(DEL) Alkane: Str...	3.50	2.5	ug/L	175910	1	6.83	353517	5.0
(DEL) Alkane: Cyc...	4.37	2.2	ug/L	155452	1	6.83	353517	5.0
Benzene, propyl-	11.35	1.0	ug/L	85799	3	12.06	416947	5.0
Benzene, 1-ethyl-...	11.42	6.0	ug/L	498420	3	12.06	416947	5.0
Benzene, 1,2,3-tr...	11.49	2.1	ug/L	177900	3	12.06	416947	5.0
4-Heptanone, 2,6-...	11.61	3.9	ug/L	327273	3	12.06	416947	5.0
Benzene, 1-ethyl-...	11.65	2.2	ug/L	184565	3	12.06	416947	5.0
Mesitylene	11.79	7.2	ug/L	596925	3	12.06	416947	5.0
Benzene, 1,2,4-tr...	12.13	2.0	ug/L	164129	3	12.06	416947	5.0
Indane	12.28	1.0	ug/L	82720	3	12.06	416947	5.0