

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU052120\
 Data File : VU038274.D
 Acq On : 21 May 2020 19:21
 Operator : JC/MD
 Sample : L2724-12DL 40X
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BE4T8DL

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\SOMUTR050720WMA.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	1.610	76	84	93	rBV	176760	209967	2.07%	0.486%
2	1.929	176	183	197	rVB	186284	246663	2.43%	0.571%
3	2.398	320	329	351	rVB	243133	382681	3.77%	0.885%
4	2.591	377	389	404	rBV	302318	492232	4.85%	1.139%
5	4.035	817	838	862	rBV2	136530	355171	3.50%	0.822%
6	4.665	1017	1034	1082	rBV	384961	978768	9.65%	2.264%
7	5.102	1153	1170	1195	rBV2	372479	840105	8.29%	1.944%
8	5.761	1351	1375	1379	rBV2	627479	1502167	14.82%	3.475%
9	5.810	1379	1390	1419	rVB	4978668	10139203	100.00%	23.457%
10	6.279	1522	1536	1565	rVB	623329	1192870	11.76%	2.760%
11	6.723	1656	1674	1688	rBV	403306	790240	7.79%	1.828%
12	7.591	1928	1944	1957	rBV2	197922	342364	3.38%	0.792%
13	7.922	2026	2047	2061	rBV	754208	1338630	13.20%	3.097%
14	8.202	2121	2134	2148	rBV	124215	207778	2.05%	0.481%
15	8.655	2258	2275	2313	rBV	1274574	2336944	23.05%	5.406%
16	9.436	2503	2518	2524	rBV	887465	1619532	15.97%	3.747%
17	9.588	2552	2565	2589	rBV	963578	1585056	15.63%	3.667%
18	9.707	2589	2602	2619	rBV	1180154	1929732	19.03%	4.464%
19	10.118	2717	2730	2749	rVB2	256805	452354	4.46%	1.047%
20	10.497	2831	2848	2866	rBV	420526	695587	6.86%	1.609%
21	10.771	2913	2933	2945	rBV	346764	563806	5.56%	1.304%
22	10.915	2961	2978	2999	rBV2	192278	366316	3.61%	0.847%
23	11.099	3021	3035	3051	rBV	94773	149037	1.47%	0.345%
24	11.305	3088	3099	3116	rVB	70853	129616	1.28%	0.300%
25	11.478	3141	3153	3168	rVB	1075556	1639377	16.17%	3.793%
26	11.822	3247	3260	3275	rBV	987767	1707789	16.84%	3.951%
27	11.906	3275	3286	3297	rVB	460876	728695	7.19%	1.686%
28	12.105	3334	3348	3365	rBV	2202792	3450532	34.03%	7.983%
29	12.205	3366	3379	3400	rVB	845450	1463239	14.43%	3.385%
30	12.584	3485	3497	3506	rBV	107556	165098	1.63%	0.382%
31	12.989	3603	3623	3631	rBV2	61017	112130	1.11%	0.259%
32	13.044	3631	3640	3650	rVB	105098	166909	1.65%	0.386%
33	13.314	3713	3724	3737	rVB	132866	206943	2.04%	0.479%
34	13.475	3764	3774	3786	rVV3	191406	315835	3.11%	0.731%

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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\SOMUTR050720WMA.M
Title : TRACE VOA SOM01.0

35	13.542	3786	3795	3806	rVB2	79316	118915	1.17%	0.275%
36	13.652	3816	3829	3840	rBV2	180758	310461	3.06%	0.718%
37	14.111	3958	3972	3995	rVB	2036685	3290840	32.46%	7.613%
38	14.234	3999	4010	4022	rVB	75685	133084	1.31%	0.308%
39	15.263	4319	4330	4347	rVB	145360	238341	2.35%	0.551%
40	15.449	4376	4388	4403	rBV	189752	330399	3.26%	0.764%

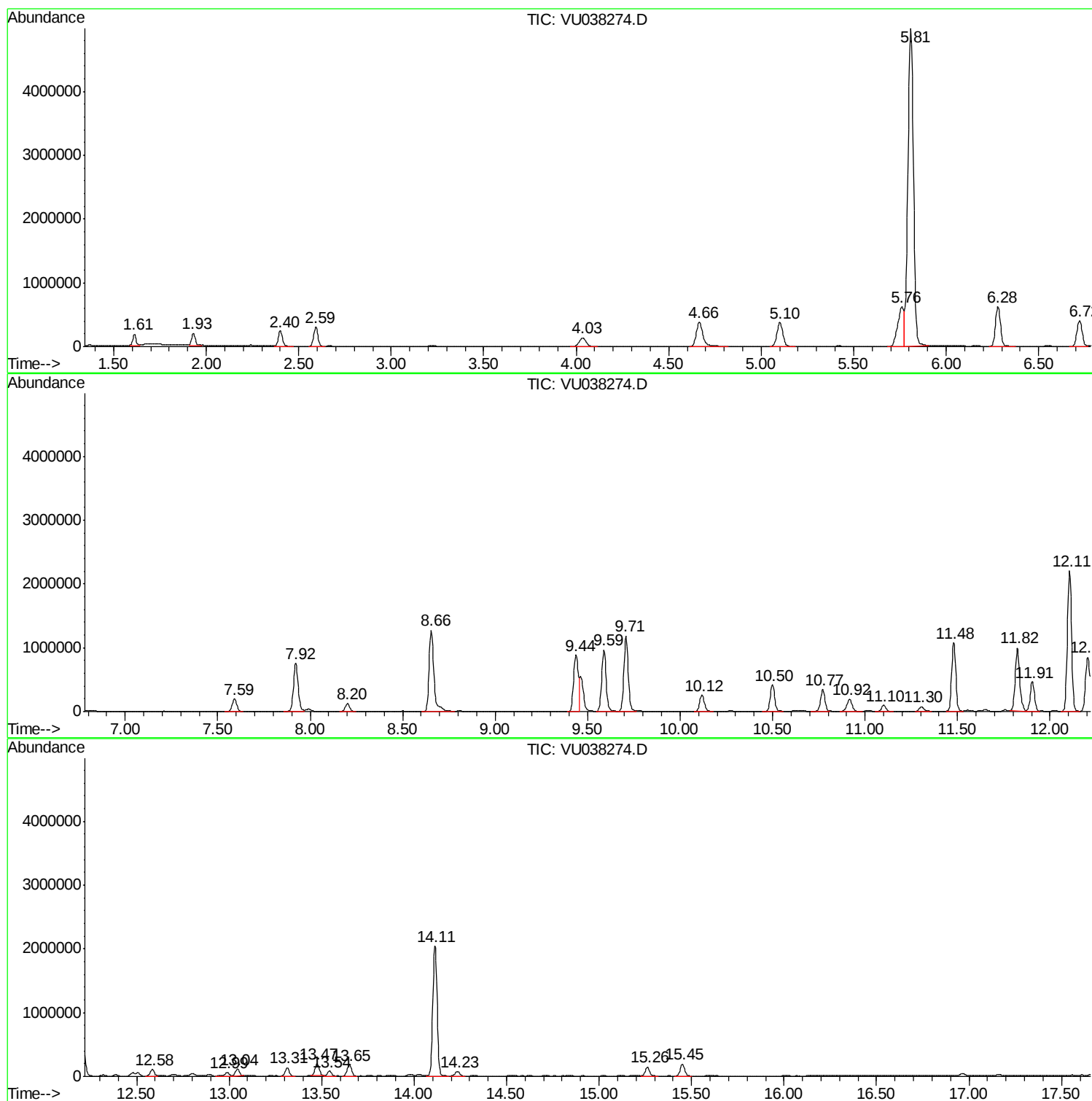
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR050720WMA.M
Quant Title : TRACE VOA SOM01.0

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



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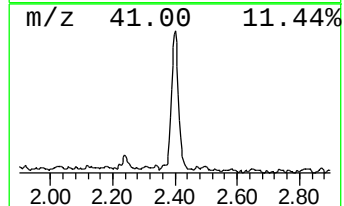
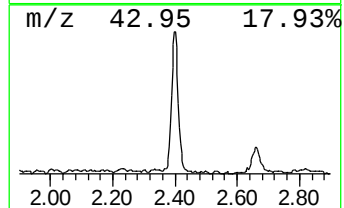
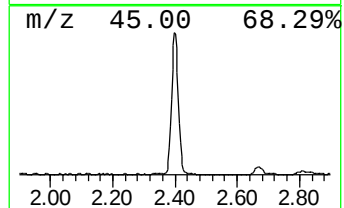
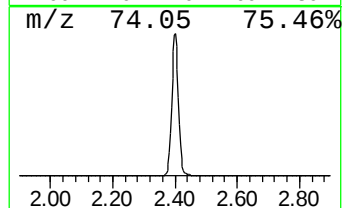
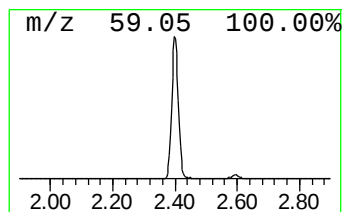
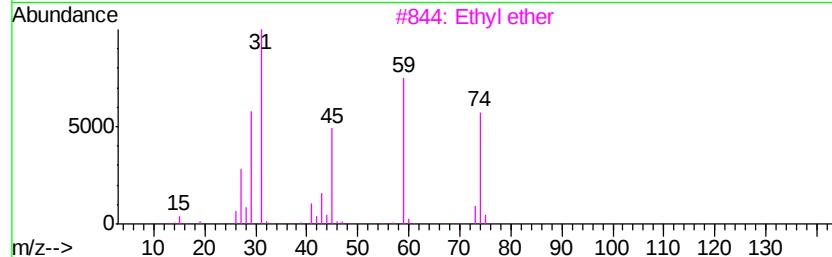
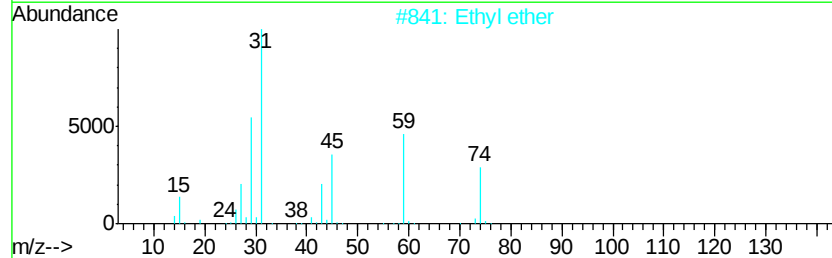
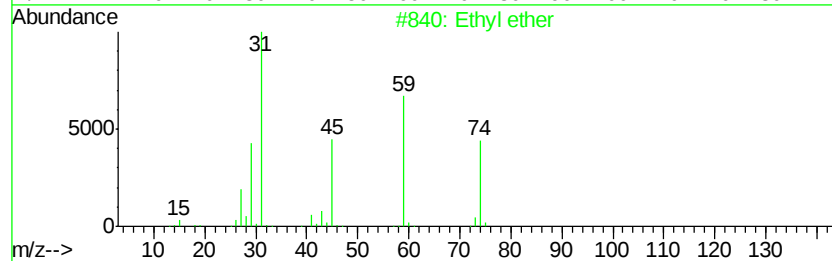
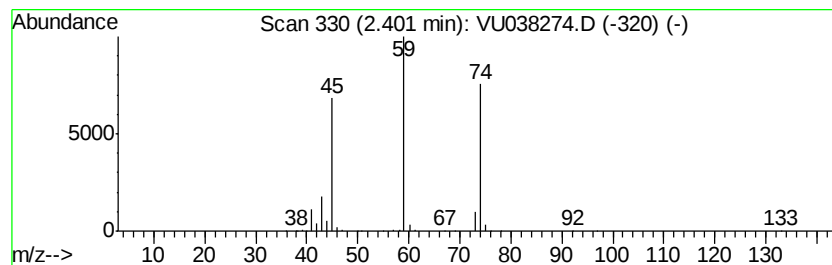
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR050720WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 1 Ethyl ether Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.40	1.60 ug/L	382681	1,4-Difluorobenzene	6.28

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethyl ether		74	C4H10O	000060-29-7	91
2	Ethyl ether		74	C4H10O	000060-29-7	90
3	Ethyl ether		74	C4H10O	000060-29-7	90
4	Ethyl ether		74	C4H10O	000060-29-7	72
5	Hydrazine, trimethyl-		74	C3H10N2	001741-01-1	59



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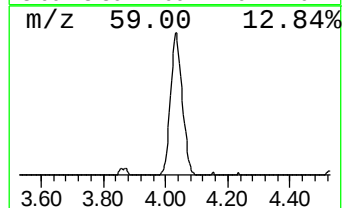
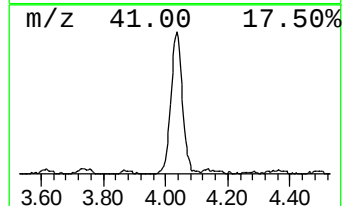
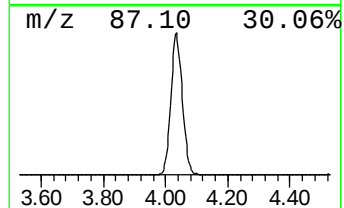
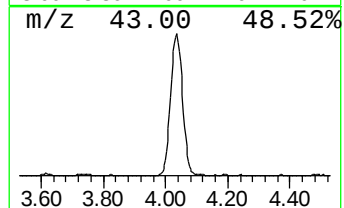
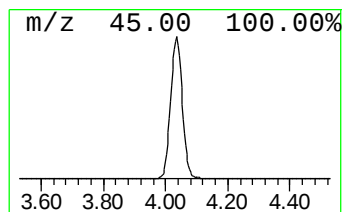
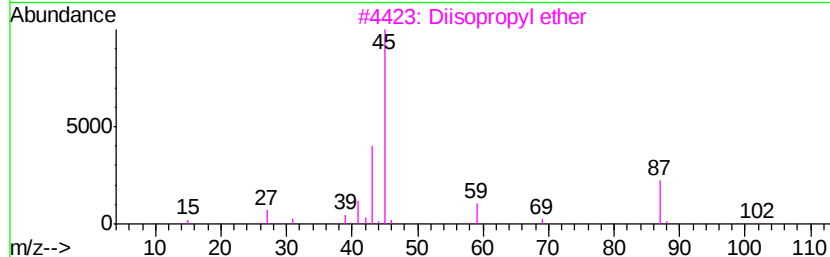
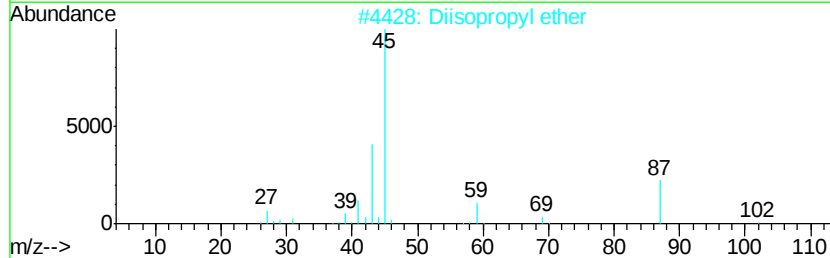
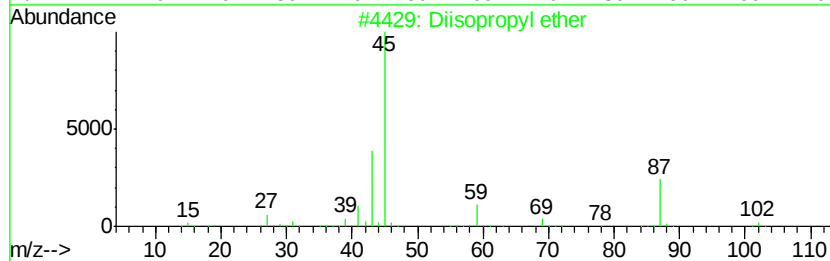
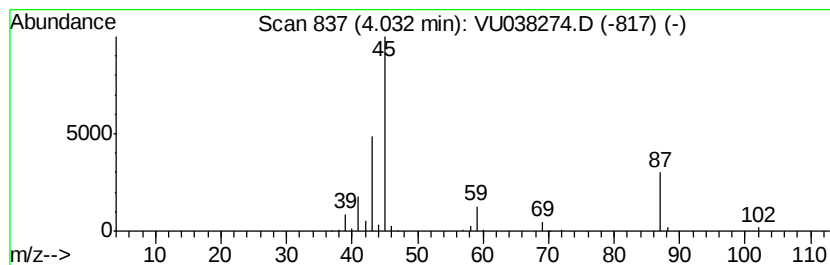
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 Diisopropyl ether Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.03	1.49 ug/L	355171	1,4-Difluorobenzene	6.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Diisopropyl ether	102	C6H14O	000108-20-3	86
2		Diisopropyl ether	102	C6H14O	000108-20-3	83
3		Diisopropyl ether	102	C6H14O	000108-20-3	83
4		1-Propanol, 2-(1-methylethoxy)-	118	C6H14O2	003944-37-4	72
5		Ether, 2-chloro-1-methylethyl is...	136	C6H13ClO	098277-76-0	64



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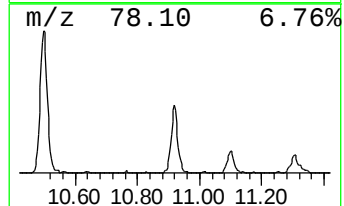
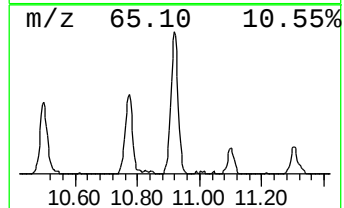
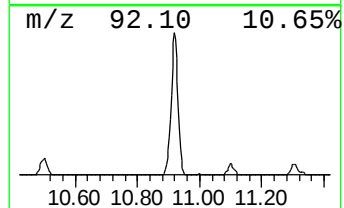
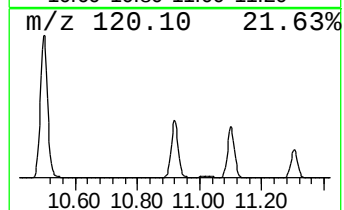
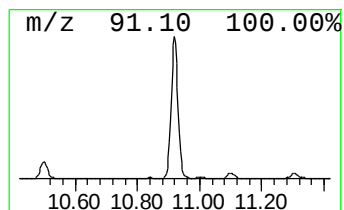
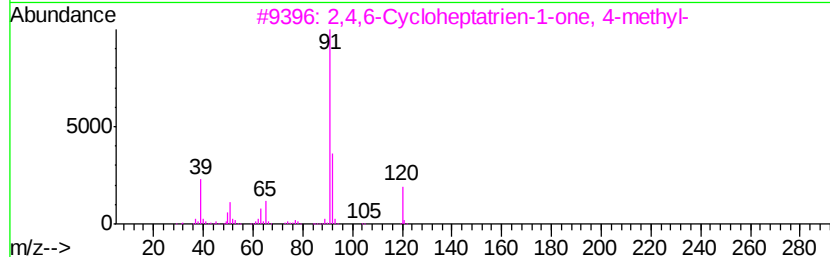
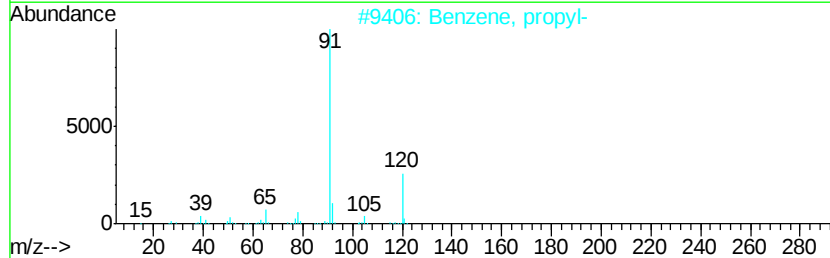
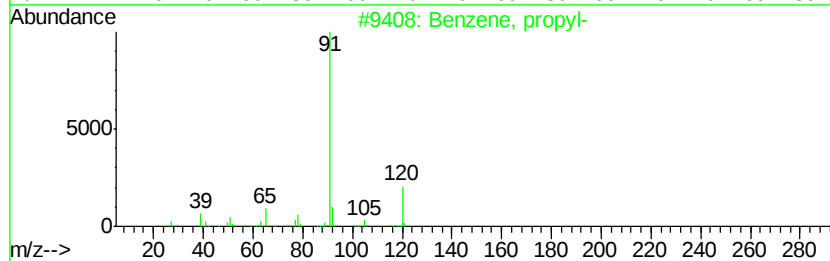
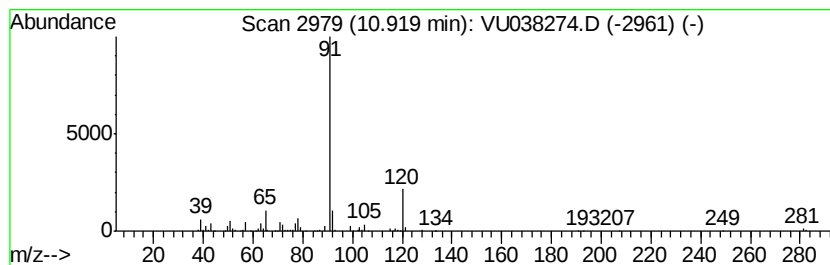
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 4 Benzene, propyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.92	1.07 ug/L	366316	1,4-Dichlorobenzene-d4	11.82

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	76
2			Benzene, propyl-	120	C9H12	000103-65-1	68
3			2,4,6-Cycloheptatrien-1-one, 4-m...	120	C8H8O	1000327-34-9	47
4			4[h]-Pyridone, 1-benzyl-3,5-dich...	281	C14H13Cl2NO	1000252-55-6	43
5			Benzeneacetaldehyde	120	C8H8O	000122-78-1	43



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Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
BE4T8DL

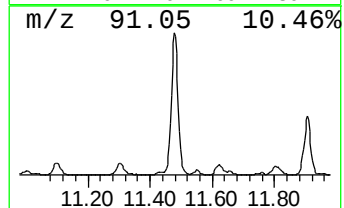
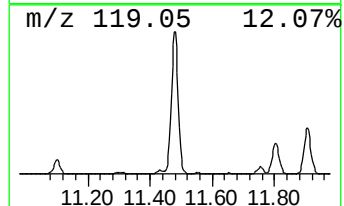
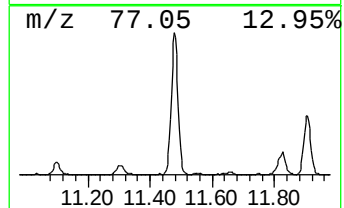
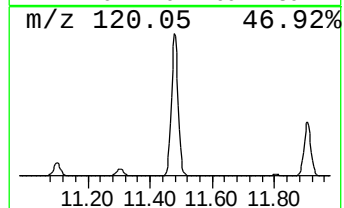
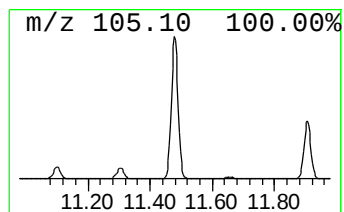
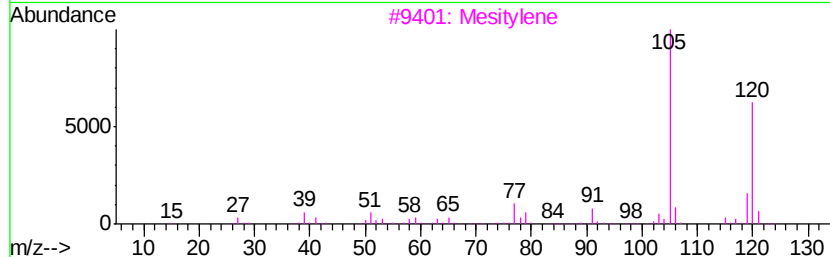
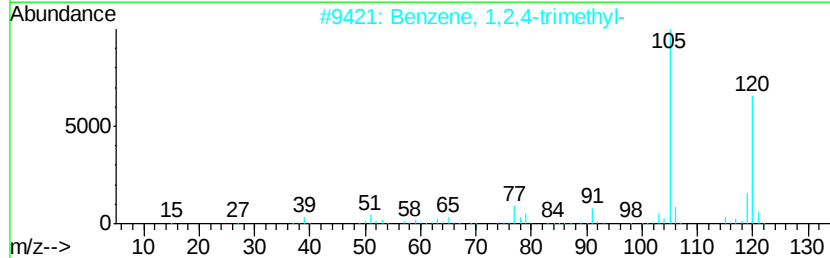
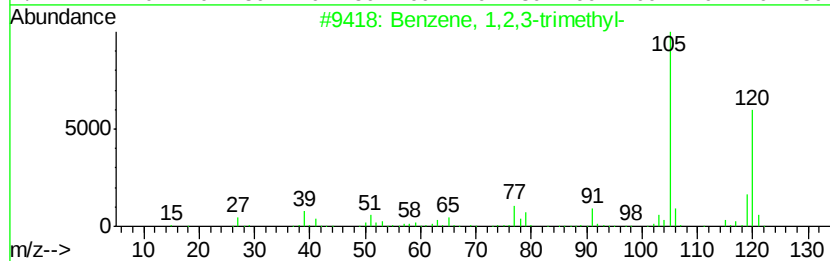
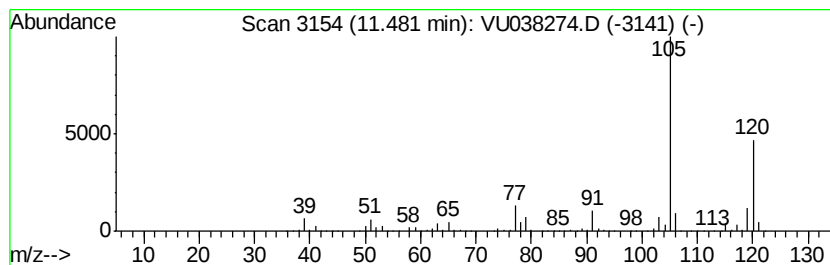
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1,2,3-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.48	4.80 ug/L	1639380	1,4-Dichlorobenzene-d4	11.82

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



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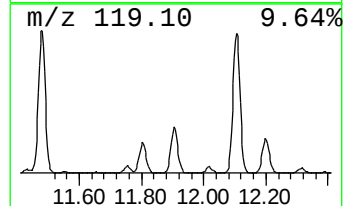
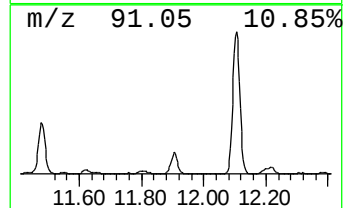
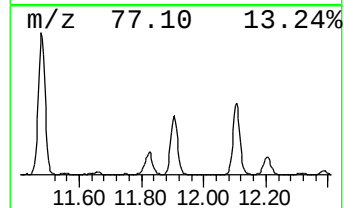
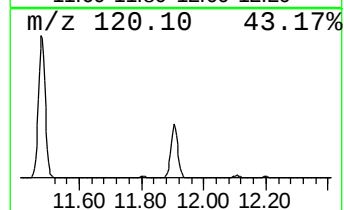
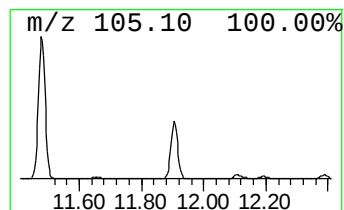
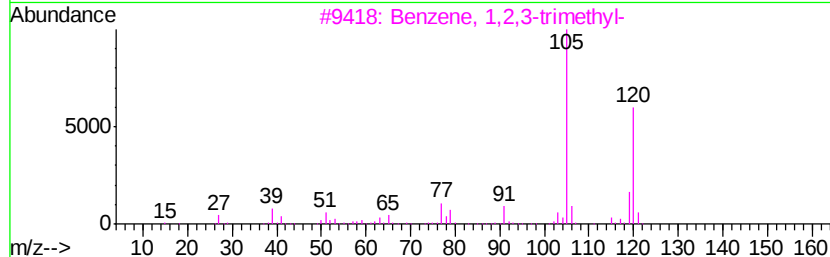
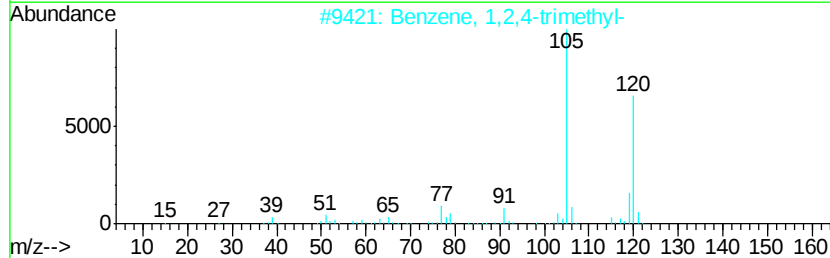
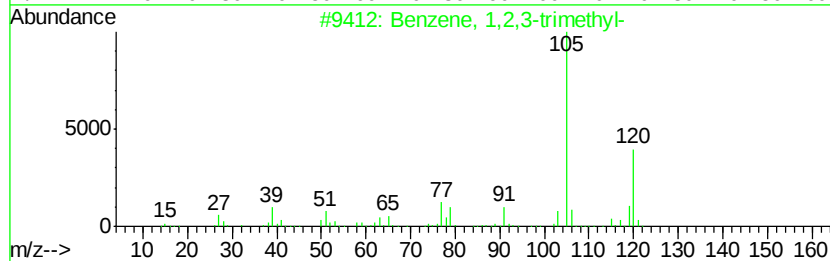
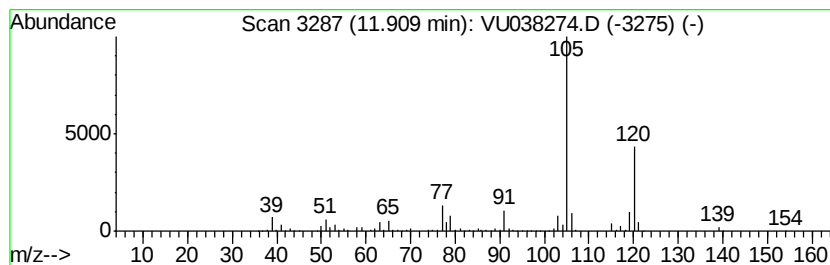
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 6 Benzene, 1,2,4-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.91	2.13 ug/L	728695	1,4-Dichlorobenzene-d4	11.82

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



Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU052120\
Data File : VU038274.D
Acq On : 21 May 2020 19:21
Operator : JC/MD
Sample : L2724-12DL 40X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BE4T8DL

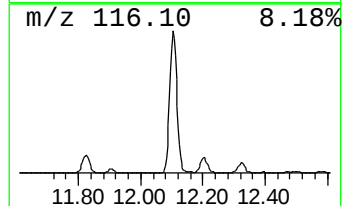
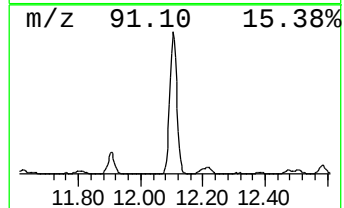
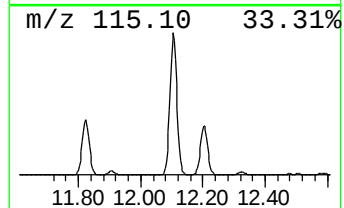
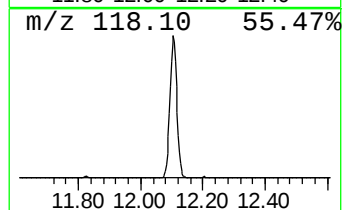
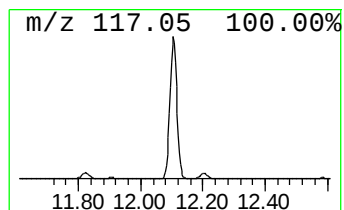
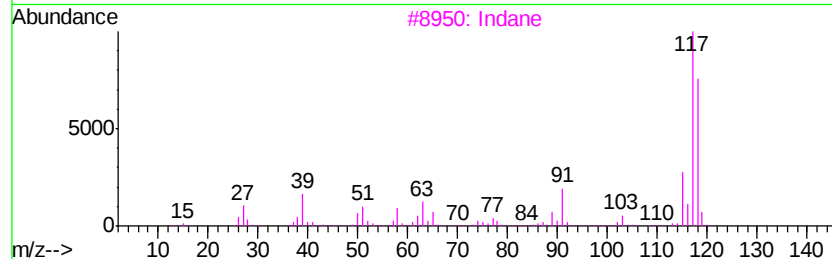
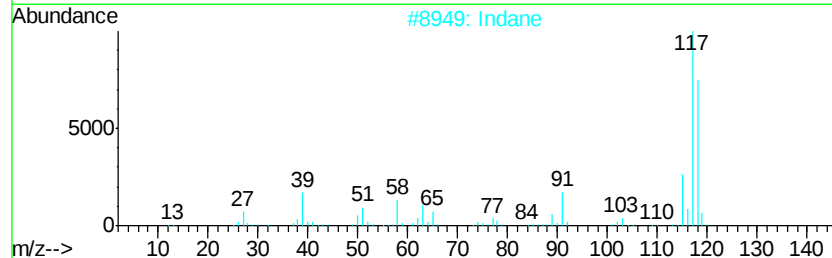
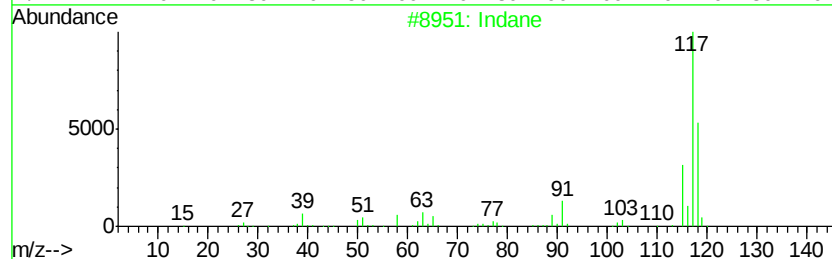
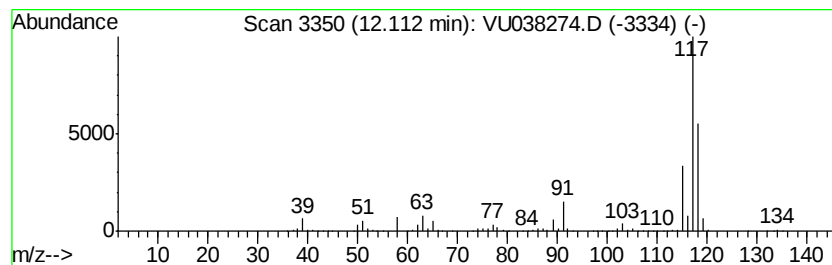
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 7 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.11	10.10 ug/L	3450530	1,4-Dichlorobenzene-d4	11.82

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	93
2	Indane		118	C9H10	000496-11-7	91
3	Indane		118	C9H10	000496-11-7	90
4	Deltacyclene		118	C9H10	007785-10-6	72
5	Indane		118	C9H10	000496-11-7	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU052120\
Data File : VU038274.D
Acq On : 21 May 2020 19:21
Operator : JC/MD
Sample : L2724-12DL 40X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BE4T8DL

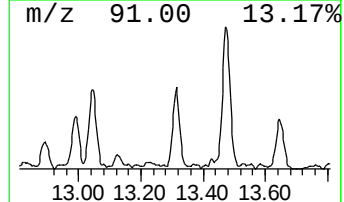
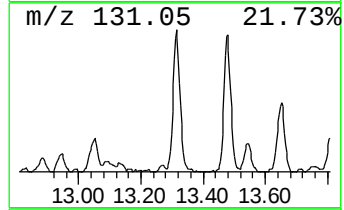
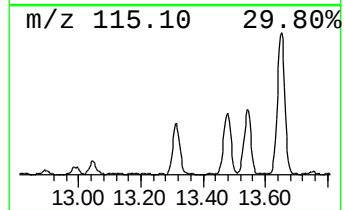
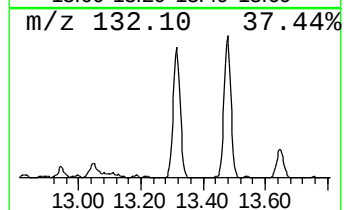
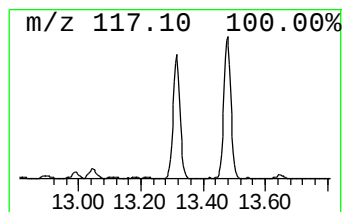
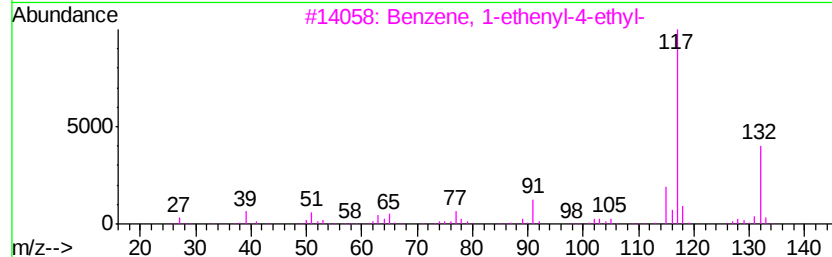
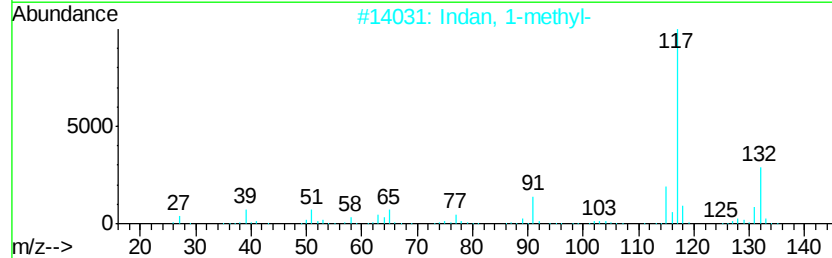
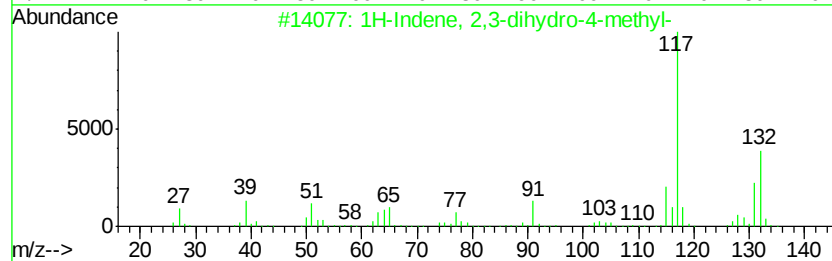
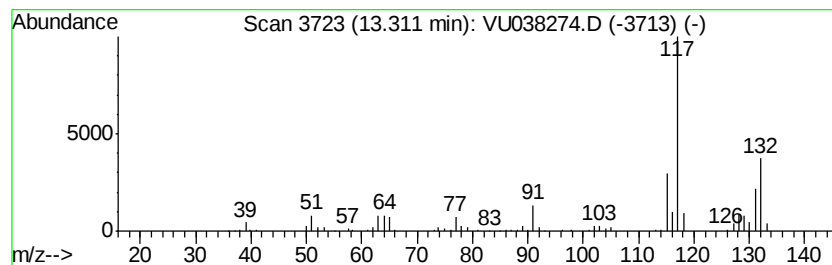
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 8 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.31	0.61 ug/L	206943	1,4-Dichlorobenzene-d4	11.82

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	91
2		Indan, 1-methyl-	132	C10H12	000767-58-8	81
3		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	74
5		Indan, 1-methyl-	132	C10H12	000767-58-8	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU052120\
Data File : VU038274.D
Acq On : 21 May 2020 19:21
Operator : JC/MD
Sample : L2724-12DL 40X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BE4T8DL

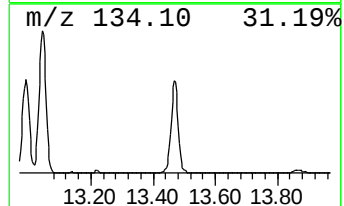
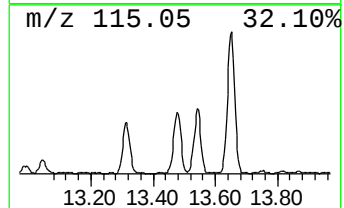
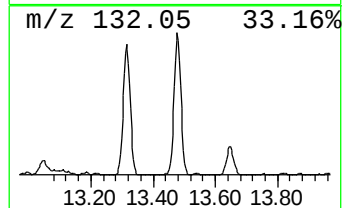
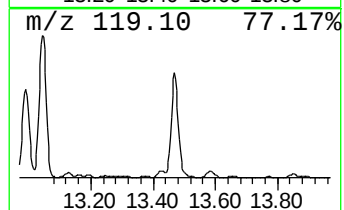
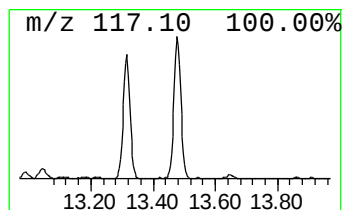
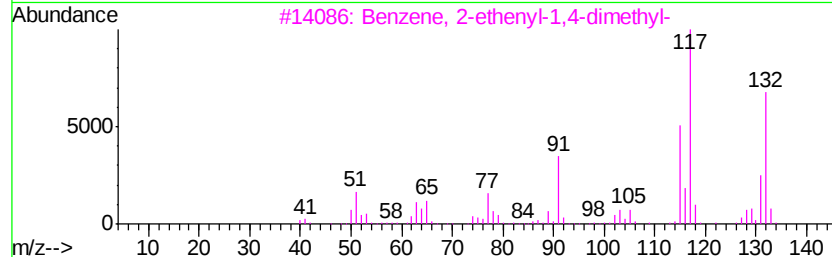
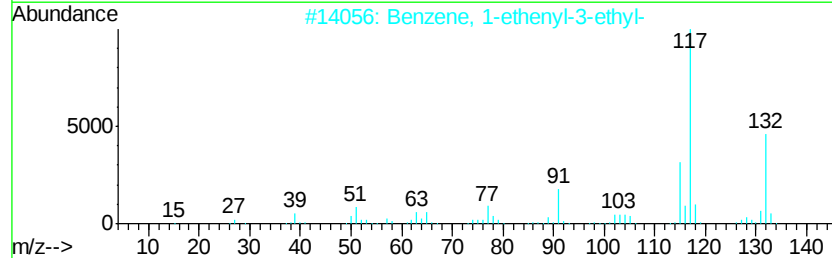
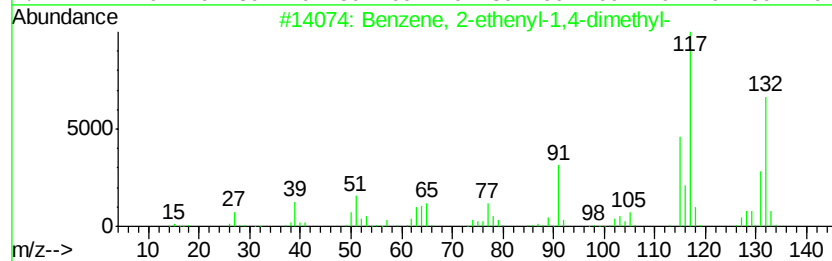
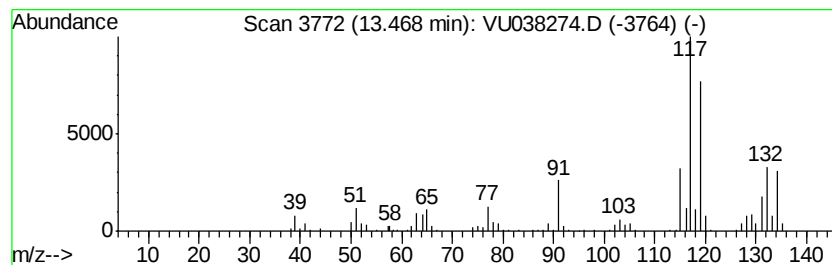
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 9 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.47	0.92 ug/L	315835	1,4-Dichlorobenzene-d4	11.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
3		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	86
4		2,4-Dimethylstyrene	132	C10H12	002234-20-0	86
5		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU052120\
Data File : VU038274.D
Acq On : 21 May 2020 19:21
Operator : JC/MD
Sample : L2724-12DL 40X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 19 Sample Multiplier: 1

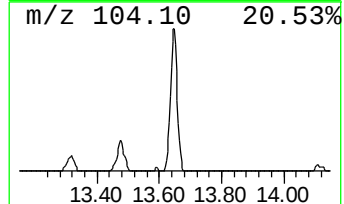
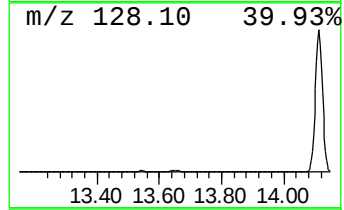
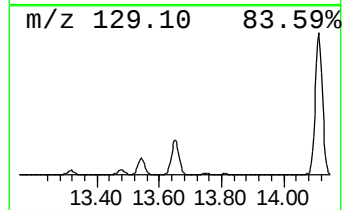
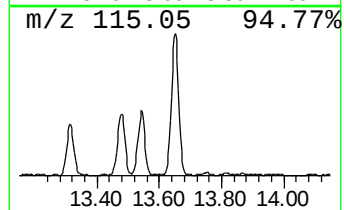
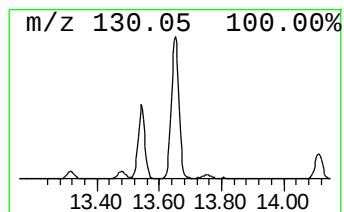
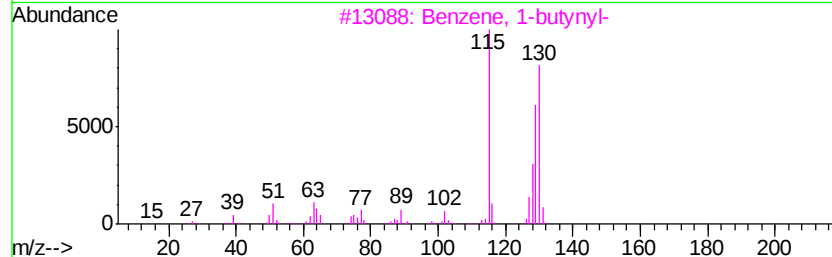
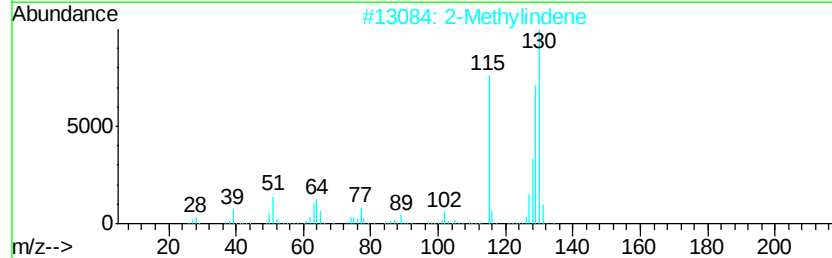
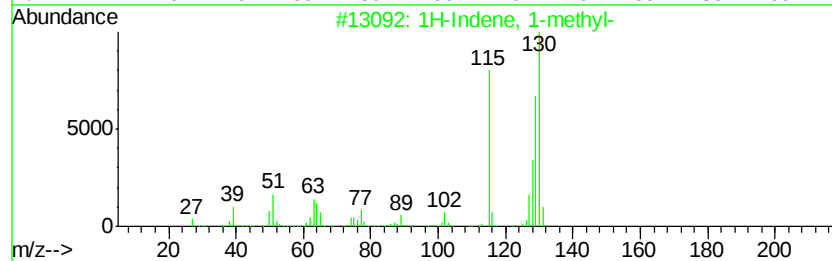
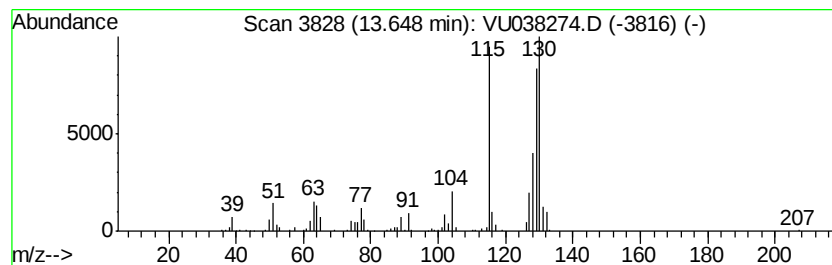
Instrument :
MSVOA_U
ClientSampleId :
BE4T8DL

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Quant Title : TRACE VOA SOM01.0

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

Peak Number 10 1H-Indene, 1-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.65	0.91 ug/L	310461	1,4-Dichlorobenzene-d4	11.82
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1H-Indene, 1-methyl-	130 C10H10	000767-59-9 96
2		2-Methylindene	130 C10H10	002177-47-1 96
3		Benzene, 1-butynyl-	130 C10H10	000622-76-4 95
4		Benzene,1-methyl-1,2-propadienyl-	130 C10H10	022433-39-2 94
5		1H-Indene, 3-methyl-	130 C10H10	000767-60-2 94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU052120\
Data File : VU038274.D
Acq On : 21 May 2020 19:21
Operator : JC/MD
Sample : L2724-12DL 40X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BE4T8DL

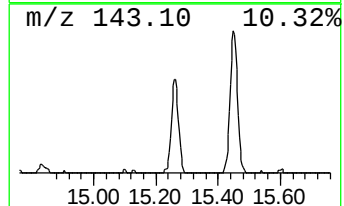
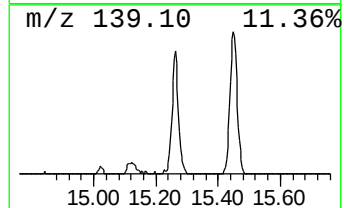
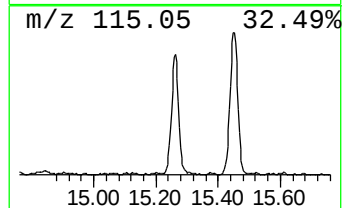
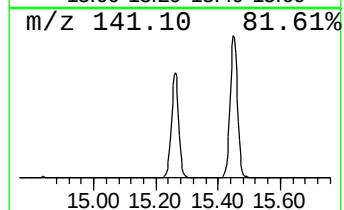
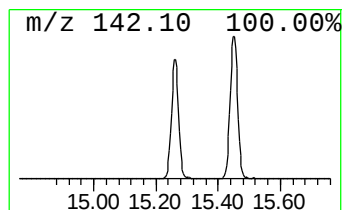
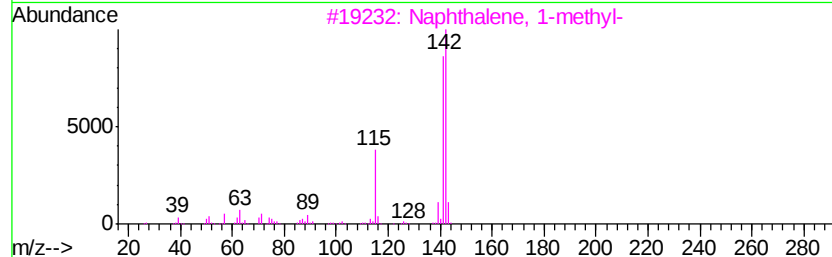
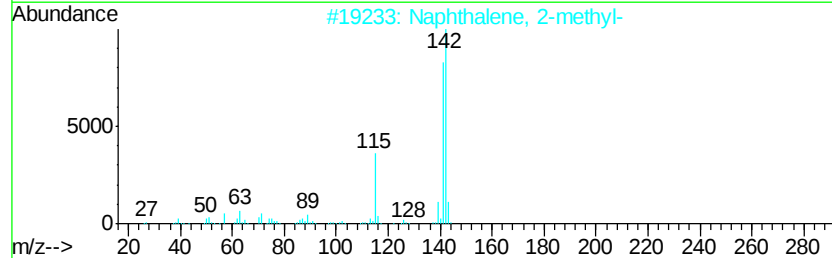
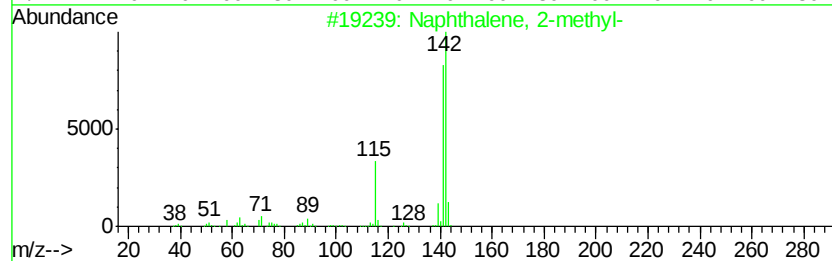
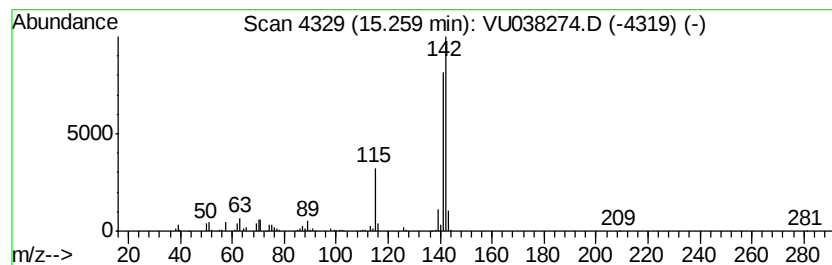
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 12 Naphthalene, 1-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.26	0.70 ug/L	238341	1,4-Dichlorobenzene-d4	11.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	95
4		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94
5		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU052120\
Data File : VU038274.D
Acq On : 21 May 2020 19:21
Operator : JC/MD
Sample : L2724-12DL 40X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BE4T8DL

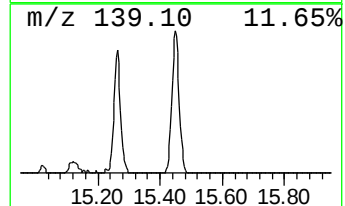
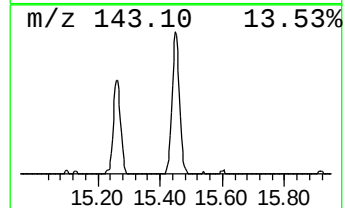
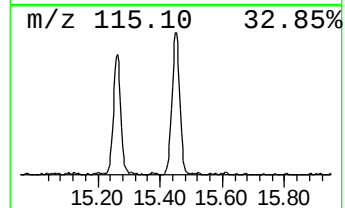
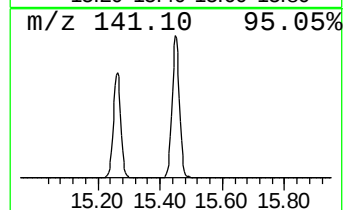
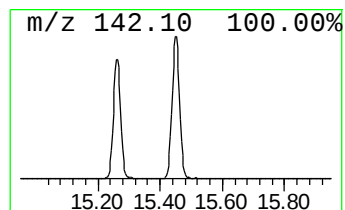
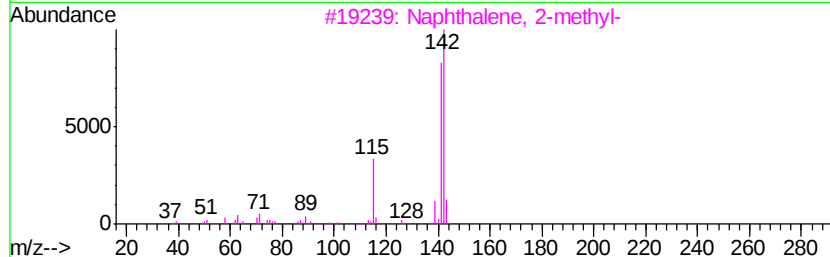
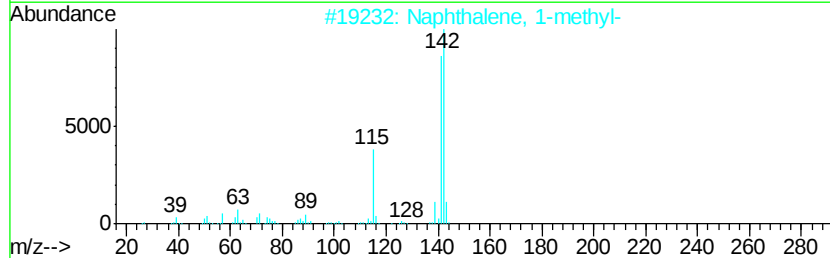
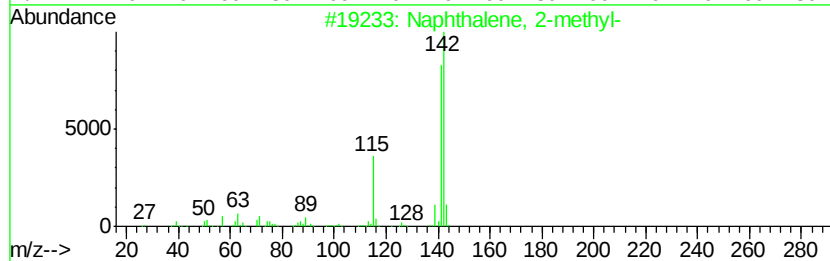
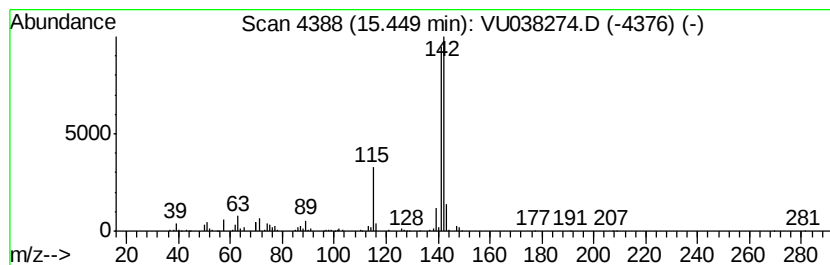
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 13 1,4-Methanonaphthalene, 1,4... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.45	0.97 ug/L	330399	1,4-Dichlorobenzene-d4	11.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	95
4		1,4-Methanonaphthalene, 1,4-dihv...	142	C11H10	004453-90-1	94
5		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU052120\
 Data File : VU038274.D
 Acq On : 21 May 2020 19:21
 Operator : JC/MD
 Sample : L2724-12DL 40X
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BE4T8DL

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR050720WMA.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
Ethyl ether	2.40	1.6	ug/L	382681	1	6.28	1192870	5.0
Diisopropyl ether	4.03	1.5	ug/L	355171	1	6.28	1192870	5.0
Benzene, propyl-	10.92	1.1	ug/L	366316	3	11.82	1707790	5.0
Benzene, 1,2,3-tr...	11.48	4.8	ug/L	1639380	3	11.82	1707790	5.0
Benzene, 1,2,4-tr...	11.91	2.1	ug/L	728695	3	11.82	1707790	5.0
Indane	12.11	10.1	ug/L	3450530	3	11.82	1707790	5.0
1H-Indene, 2,3-di...	13.31	0.6	ug/L	206943	3	11.82	1707790	5.0
Benzene, 2-etheny...	13.47	0.9	ug/L	315835	3	11.82	1707790	5.0
1H-Indene, 1-methyl-	13.65	0.9	ug/L	310461	3	11.82	1707790	5.0
Naphthalene, 1-me...	15.26	0.7	ug/L	238341	3	11.82	1707790	5.0
1,4-Methanonaphth...	15.45	1.0	ug/L	330399	3	11.82	1707790	5.0