

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU060620\
 Data File : VU038613.D
 Acq On : 06 Jun 2020 13:49
 Operator : SY/MD
 Sample : L2884-04DL 10X
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BFXD8DL

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\SOMUTR053120WMA.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	1.613	74	85	93	rBV	103614	120080	6.25%	0.674%
2	1.932	175	184	198	rBV	90538	116612	6.07%	0.655%
3	2.594	376	390	407	rBV	187156	312238	16.25%	1.753%
4	3.218	571	584	600	rVB2	12936	28148	1.47%	0.158%
5	3.391	621	638	652	rBV4	13427	29397	1.53%	0.165%
6	4.568	987	1004	1018	rBV3	9176	21810	1.14%	0.122%
7	4.674	1020	1037	1067	rBV	209607	675121	35.14%	3.791%
8	5.102	1154	1170	1184	rBV	195613	424721	22.11%	2.385%
9	5.420	1253	1269	1281	rBV8	18901	46847	2.44%	0.263%
10	5.497	1281	1293	1307	rVB4	19084	41643	2.17%	0.234%
11	5.761	1354	1375	1386	rBV3	379018	1014874	52.83%	5.699%
12	5.883	1404	1413	1424	rVB6	14118	28964	1.51%	0.163%
13	6.282	1521	1537	1555	rBV	395063	752747	39.19%	4.227%
14	6.565	1606	1625	1627	rBV	11370	24520	1.28%	0.138%
15	6.719	1659	1673	1688	rBV	241798	476048	24.78%	2.673%
16	6.790	1688	1695	1704	rVB5	11207	21935	1.14%	0.123%
17	7.591	1930	1944	1964	rBV	124271	219106	11.41%	1.230%
18	7.922	2025	2047	2060	rBV	449626	795778	41.43%	4.468%
19	7.992	2060	2069	2083	rVB	105300	182338	9.49%	1.024%
20	8.201	2122	2134	2146	rBV	81140	138061	7.19%	0.775%
21	8.655	2260	2275	2304	rBV	848369	1527646	79.53%	8.578%
22	9.436	2504	2518	2543	rBV2	568949	1081893	56.32%	6.075%
23	9.587	2554	2565	2582	rBV	235467	390113	20.31%	2.191%
24	9.709	2588	2603	2622	rBV	1152441	1920960	100.00%	10.787%
25	10.115	2715	2729	2755	rVB	769026	1234462	64.26%	6.932%
26	10.497	2837	2848	2863	rVB2	51602	78748	4.10%	0.442%
27	10.584	2863	2875	2889	rBV	84986	140993	7.34%	0.792%
28	10.771	2921	2933	2947	rBV2	217253	347973	18.11%	1.954%
29	10.918	2964	2979	2992	rVB	127387	224181	11.67%	1.259%
30	11.012	2996	3008	3013	rBV2	164206	279008	14.52%	1.567%
31	11.098	3025	3035	3050	rVB	122651	189759	9.88%	1.066%
32	11.243	3065	3080	3088	rBV2	18324	28671	1.49%	0.161%
33	11.304	3088	3099	3114	rVB	125713	198824	10.35%	1.116%
34	11.478	3141	3153	3169	rVB	440539	675296	35.15%	3.792%

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 Title : TRACE VOA SOM01.0

35	11.822	3247	3260	3275	rVV	631944	1037326	54.00%	5.825%
36	11.906	3276	3286	3297	rVB	178428	275314	14.33%	1.546%
37	12.105	3332	3348	3366	rBV2	131527	243546	12.68%	1.368%
38	12.205	3366	3379	3401	rVV2	605929	1186681	61.78%	6.663%
39	12.388	3425	3436	3448	rVB	29793	48507	2.53%	0.272%
40	12.478	3455	3464	3469	rBV	26814	41869	2.18%	0.235%
41	12.507	3470	3473	3480	rVV2	17371	21872	1.14%	0.123%
42	12.558	3480	3489	3492	rVV	32107	48580	2.53%	0.273%
43	12.584	3493	3497	3505	rVV	43525	59720	3.11%	0.335%
44	12.700	3522	3533	3553	rVV6	26847	76735	3.99%	0.431%
45	12.793	3553	3562	3568	rVV3	11803	20705	1.08%	0.116%
46	12.838	3568	3576	3585	rVV	45497	73575	3.83%	0.413%
47	12.893	3585	3593	3606	rVB2	29292	46583	2.42%	0.262%
48	12.992	3608	3624	3632	rBV2	68768	108679	5.66%	0.610%
49	13.047	3632	3641	3653	rVB2	51928	81052	4.22%	0.455%
50	13.317	3714	3725	3736	rVB	42831	66342	3.45%	0.373%
51	13.478	3765	3775	3787	rVB2	98087	165860	8.63%	0.931%
52	13.648	3817	3828	3840	rVB10	8547	19262	1.00%	0.108%
53	13.880	3886	3900	3910	rBV3	53148	106701	5.55%	0.599%
54	14.114	3961	3973	3985	rBV	27604	46676	2.43%	0.262%
55	14.844	4191	4200	4214	rBV	20014	32779	1.71%	0.184%
56	16.150	4594	4606	4625	rBV2	127818	210912	10.98%	1.184%

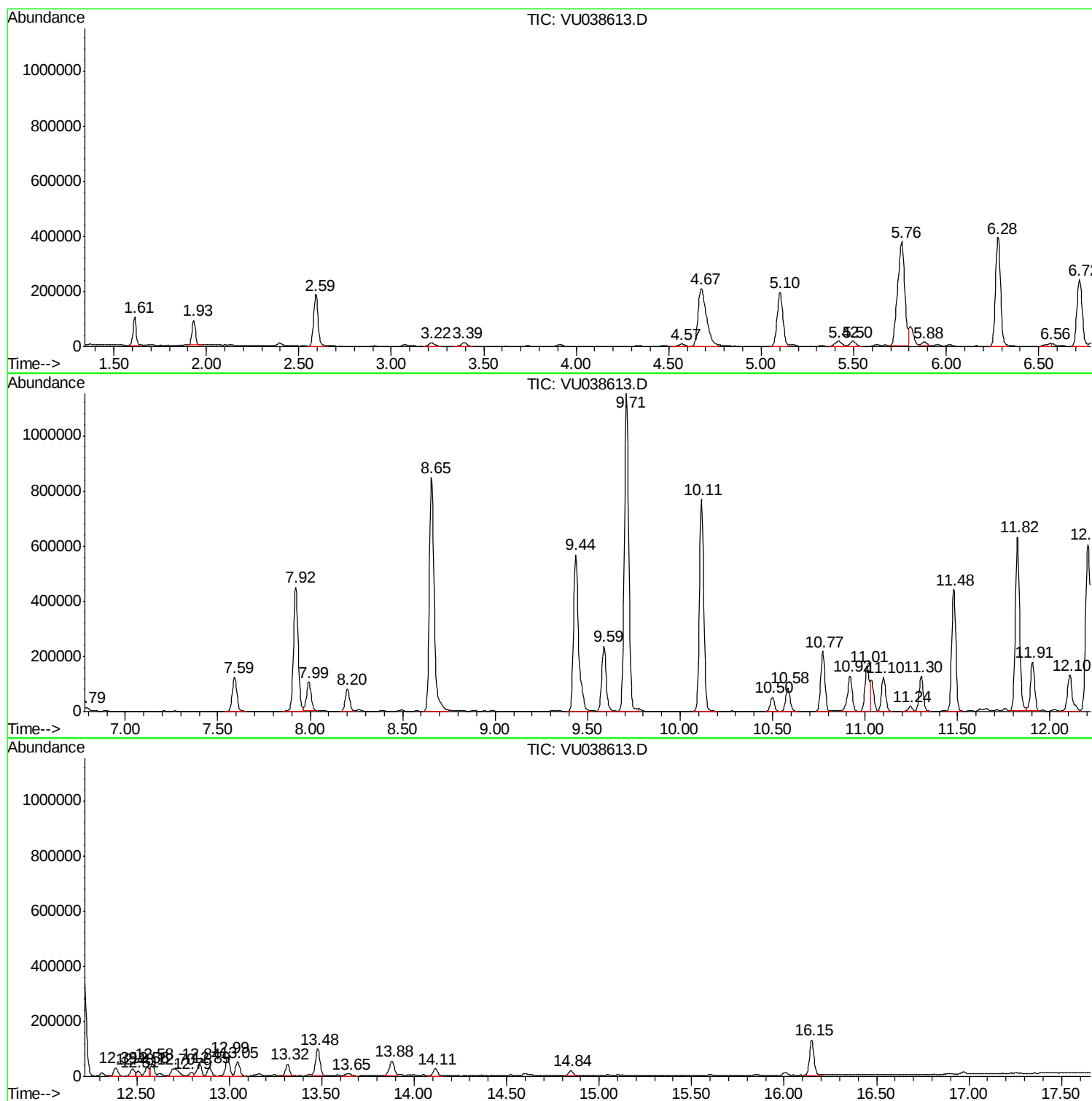
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR053120WMA.M
Quant Title : TRACE VOA SOM01.0

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



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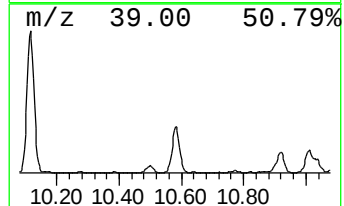
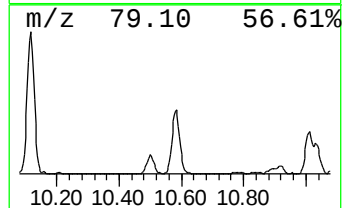
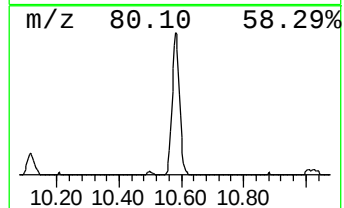
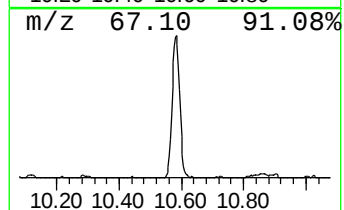
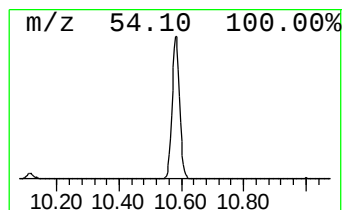
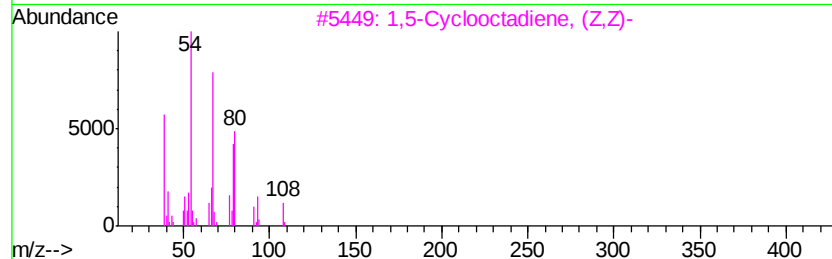
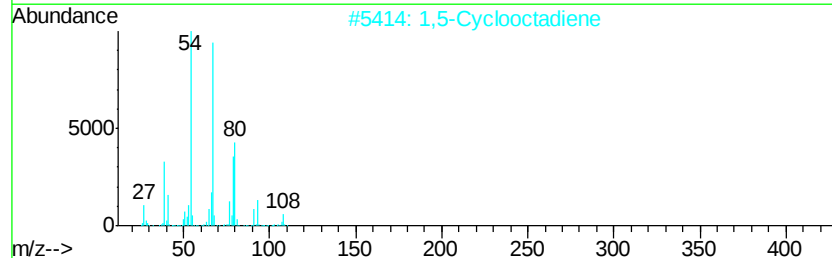
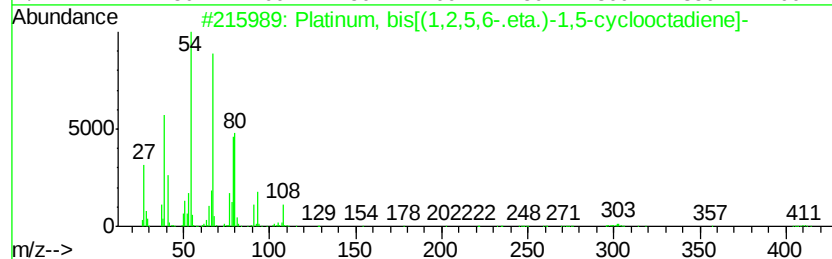
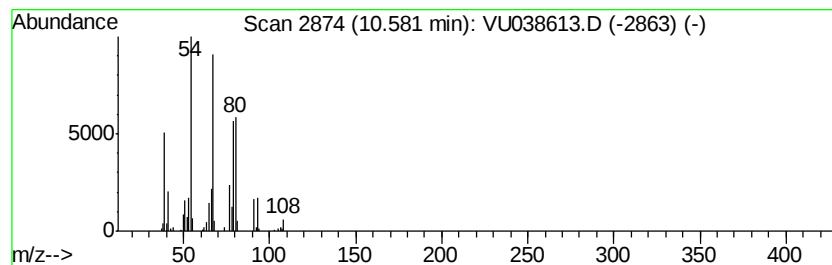
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 2 Platinum, bis[(1,2,5,6-.eta... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.58	0.65 ug/L	140993	Chlorobenzene-d5	9.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Platinum, bis[(1,2,5,6-.eta.)-1,...	411	C16H24Pt	012130-66-4	90
2		1,5-Cyclooctadiene	108	C8H12	000111-78-4	87
3		1,5-Cyclooctadiene, (Z,Z)-	108	C8H12	001552-12-1	83
4		1,5-Cyclooctadiene	108	C8H12	000111-78-4	72
5		1,5-Cyclooctadiene, (E,Z)-	108	C8H12	005259-71-2	72



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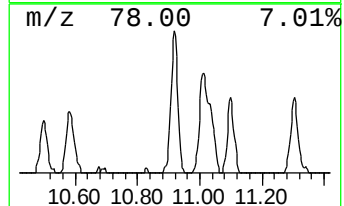
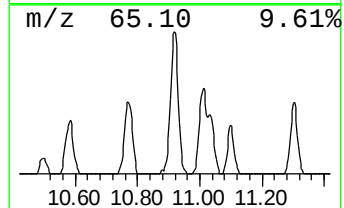
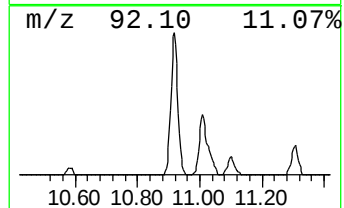
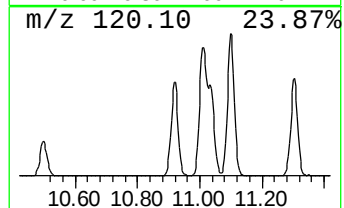
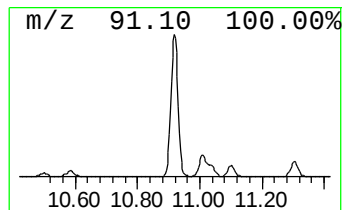
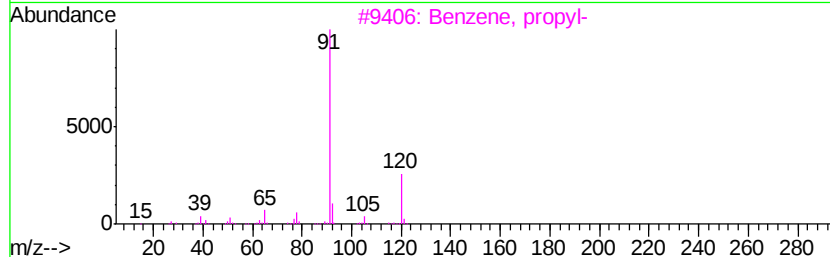
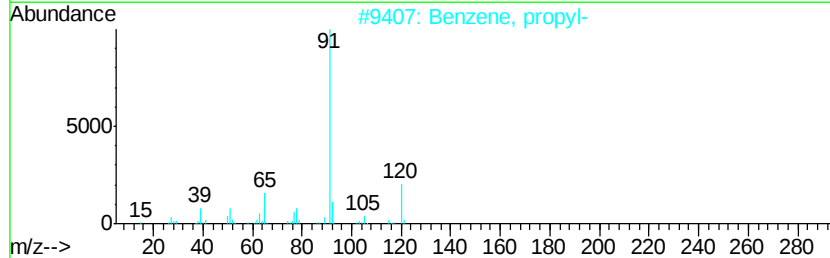
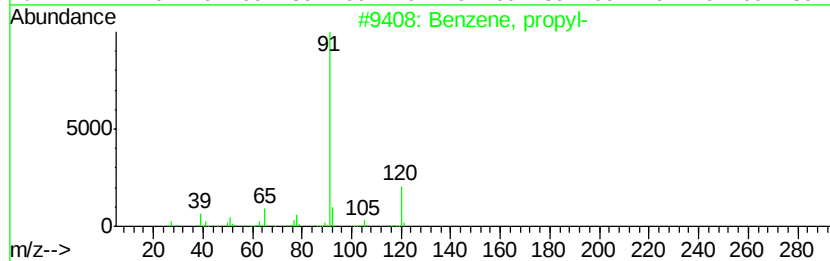
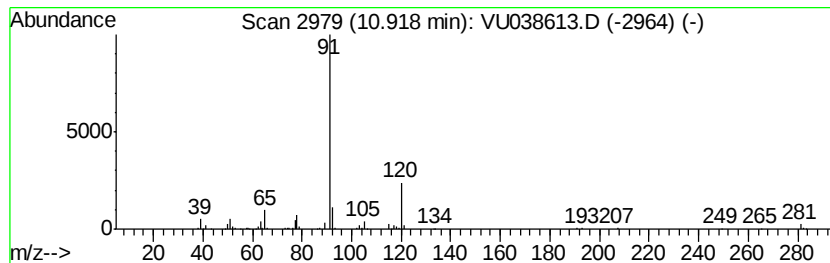
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR053120WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 3 Benzene, propyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.92	1.08 ug/L	224181	1,4-Dichlorobenzene-d4	11.83

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	87
3		Benzene, propyl-	120	C9H12	000103-65-1	80
4		N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	72
5		Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	64



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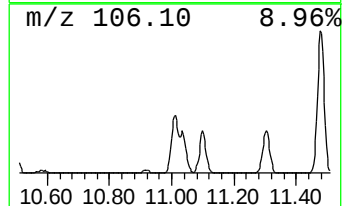
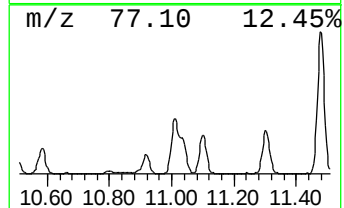
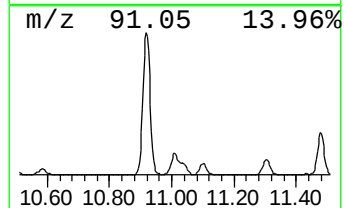
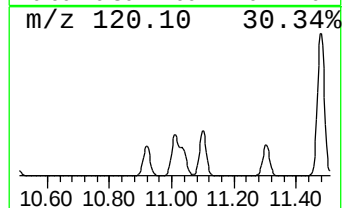
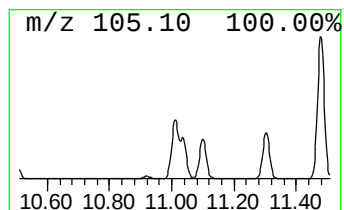
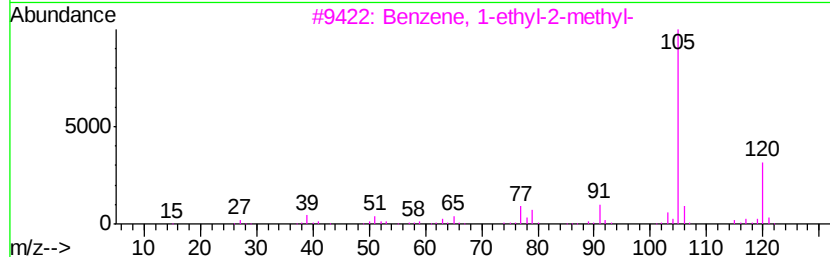
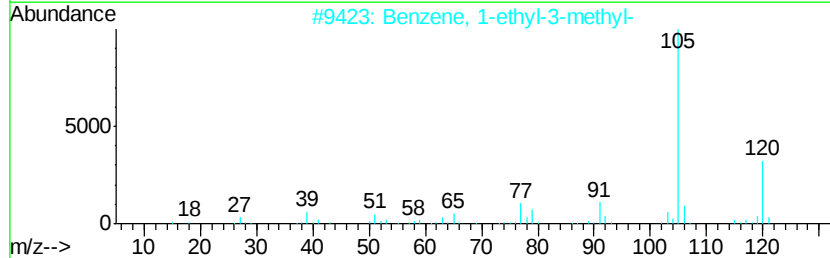
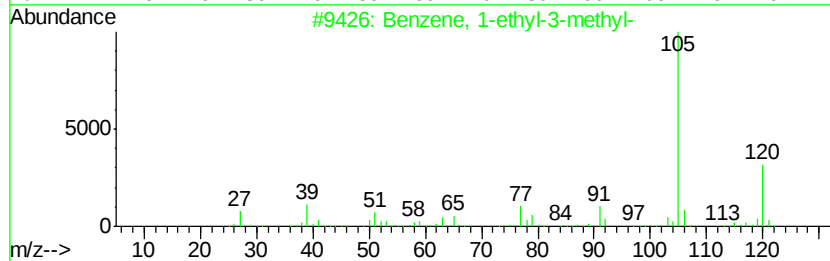
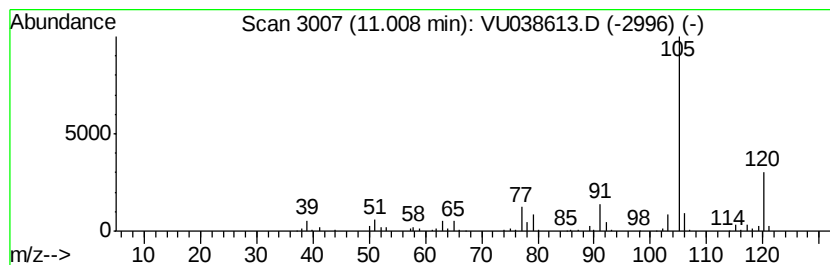
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 4 Benzene, 1-ethyl-3-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.01	1.34 ug/L	279008	1,4-Dichlorobenzene-d4	11.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
2		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
3		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
4		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



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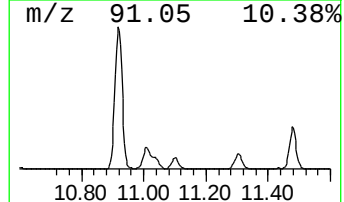
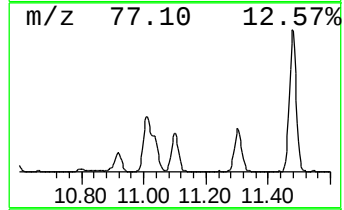
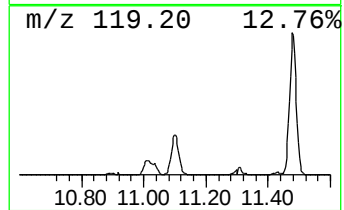
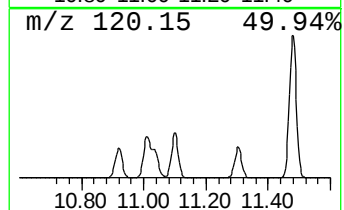
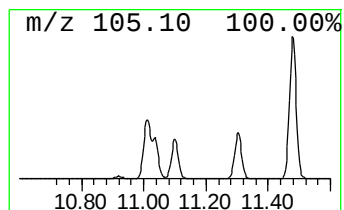
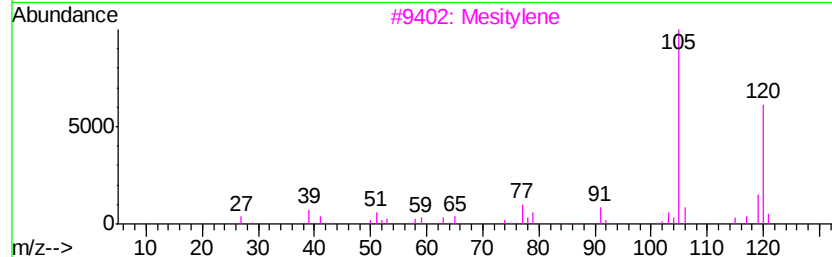
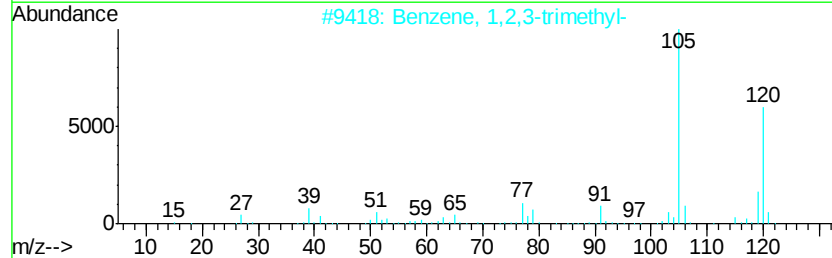
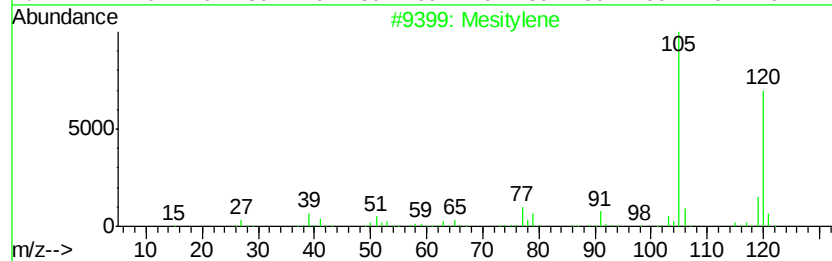
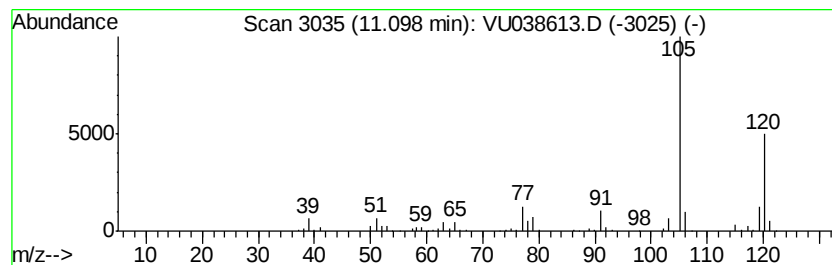
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 5 Mesitylene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.10	0.91 ug/L	189759	1,4-Dichlorobenzene-d4	11.83

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



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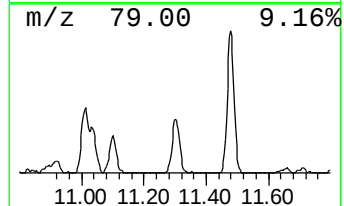
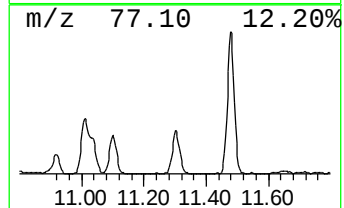
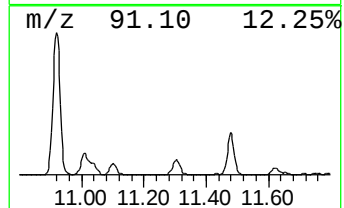
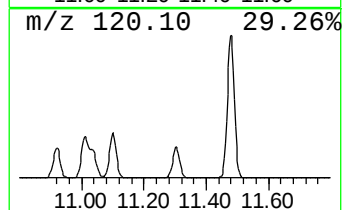
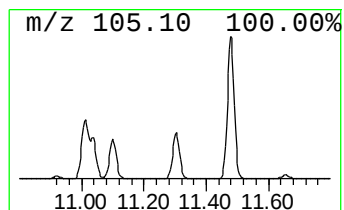
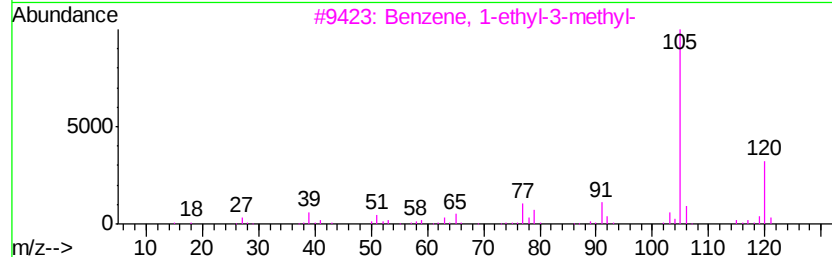
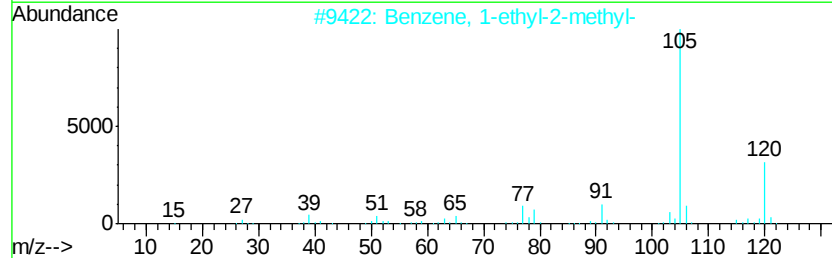
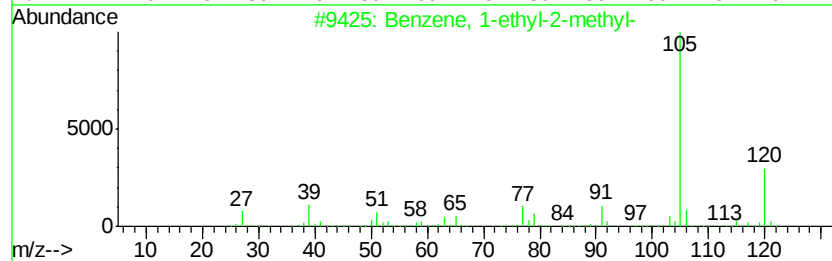
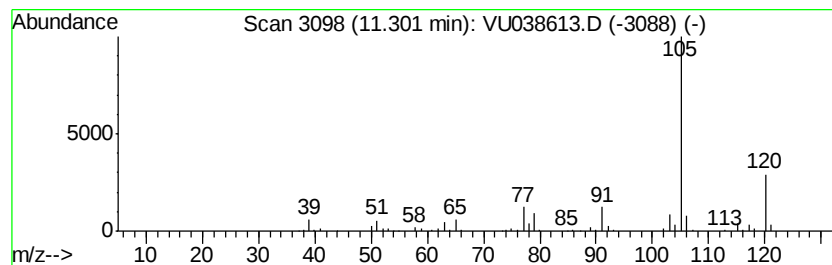
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MSVOA_U
ClientSampleId :
BFXD8DL

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR053120WMA.M
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 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.30	0.96 ug/L	198824	1,4-Dichlorobenzene-d4	11.83
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethyl-2-methyl-	120 C9H12	000611-14-3 93
2		Benzene, 1-ethyl-2-methyl-	120 C9H12	000611-14-3 93
3		Benzene, 1-ethyl-3-methyl-	120 C9H12	000620-14-4 93
4		Benzene, 1-ethyl-4-methyl-	120 C9H12	000622-96-8 93
5		Benzene, 1,2,4-trimethyl-	120 C9H12	000095-63-6 91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU060620\
Data File : VU038613.D
Acq On : 06 Jun 2020 13:49
Operator : SY/MD
Sample : L2884-04DL 10X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BFXD8DL

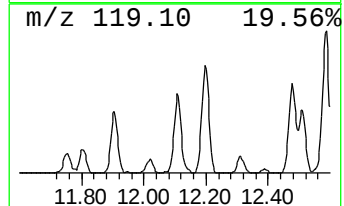
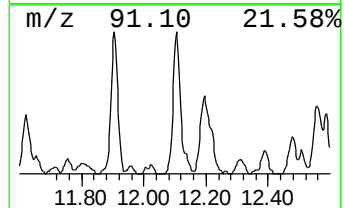
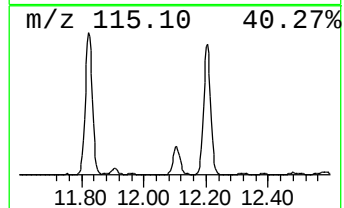
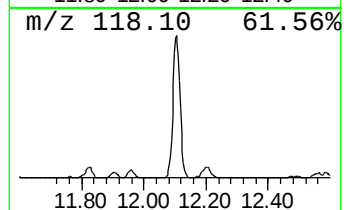
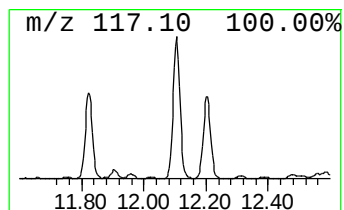
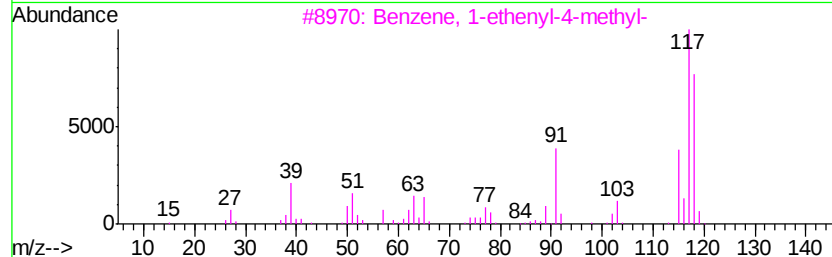
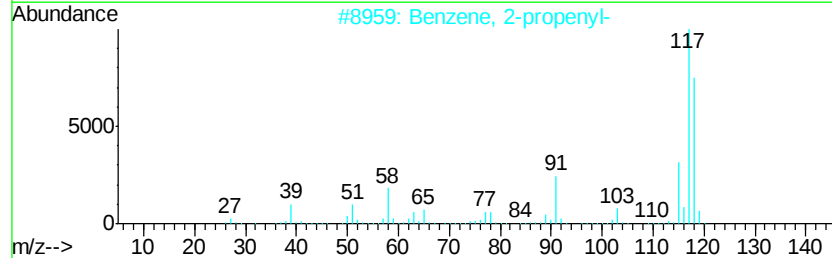
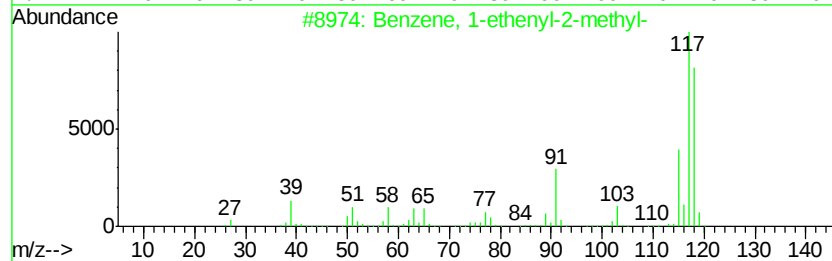
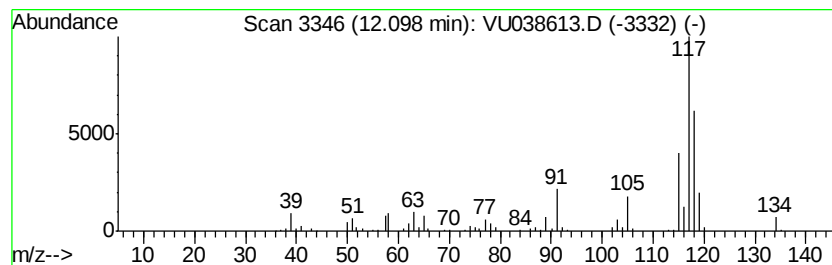
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 9 Benzene, 1-ethenyl-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.10	1.17 ug/L	243546	1,4-Dichlorobenzene-d4	11.83

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	83
2		Benzene, 2-propenyl-	118	C9H10	000300-57-2	78
3		Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	64
4		Indane	118	C9H10	000496-11-7	64
5		Indane	118	C9H10	000496-11-7	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU060620\
Data File : VU038613.D
Acq On : 06 Jun 2020 13:49
Operator : SY/MD
Sample : L2884-04DL 10X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BFXD8DL

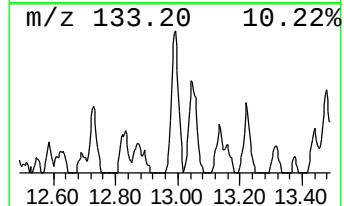
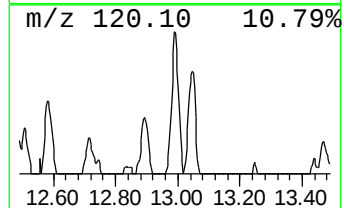
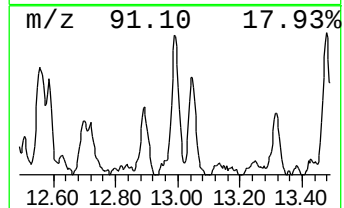
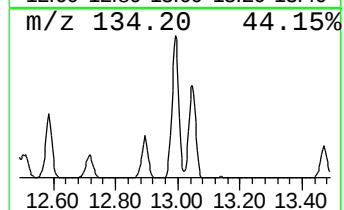
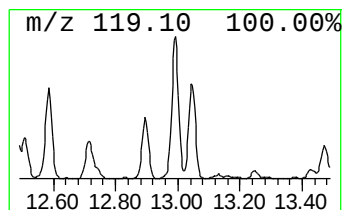
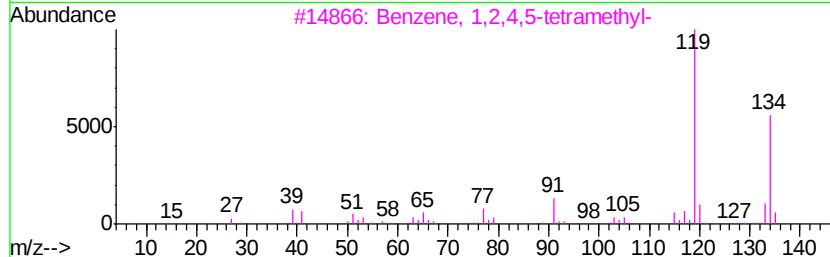
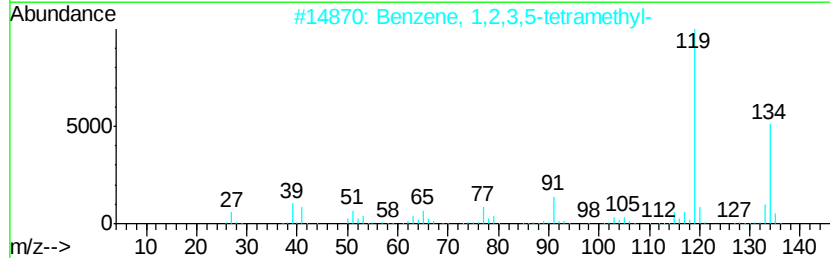
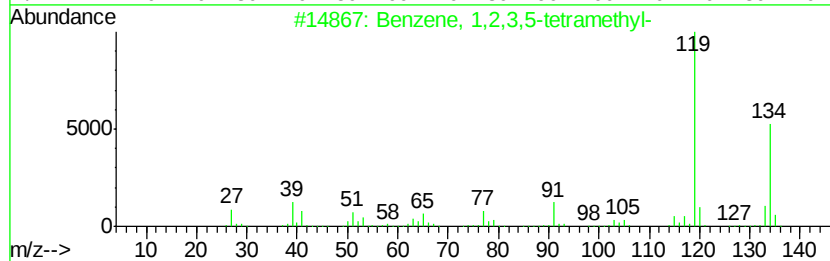
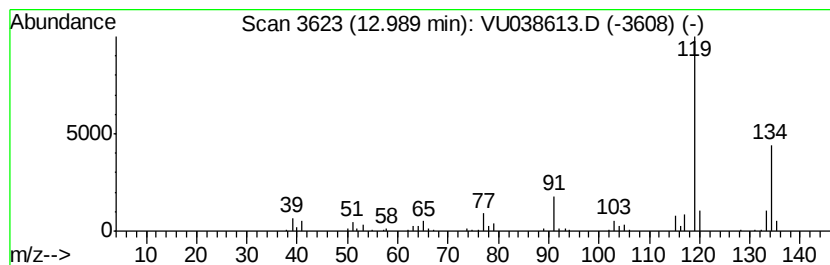
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 10 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.99	0.52 ug/L	108679	1,4-Dichlorobenzene-d4	11.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
2		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
3		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94
4		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
5		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU060620\
Data File : VU038613.D
Acq On : 06 Jun 2020 13:49
Operator : SY/MD
Sample : L2884-04DL 10X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BFXD8DL

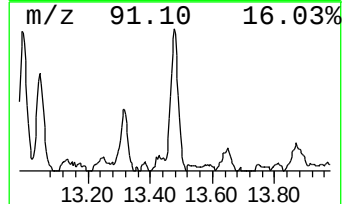
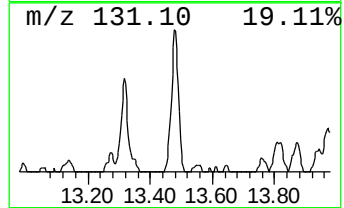
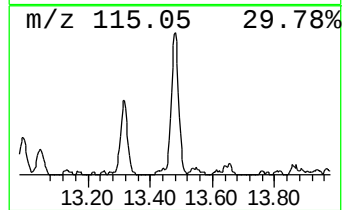
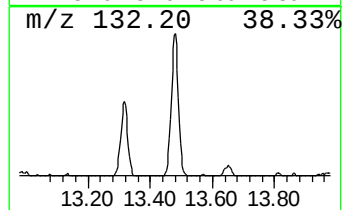
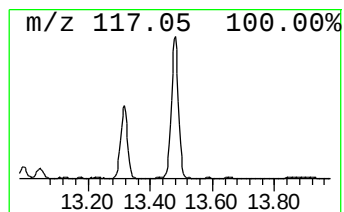
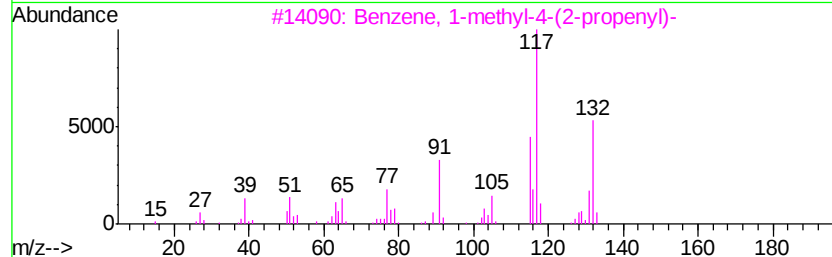
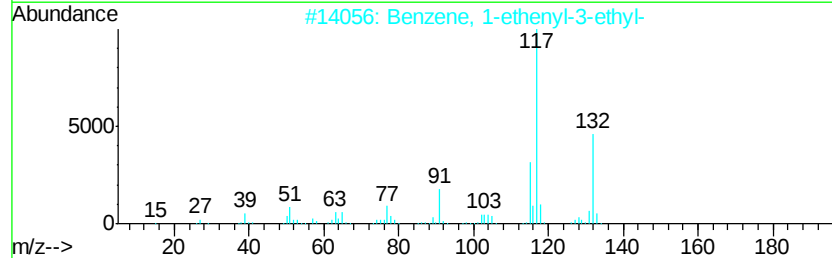
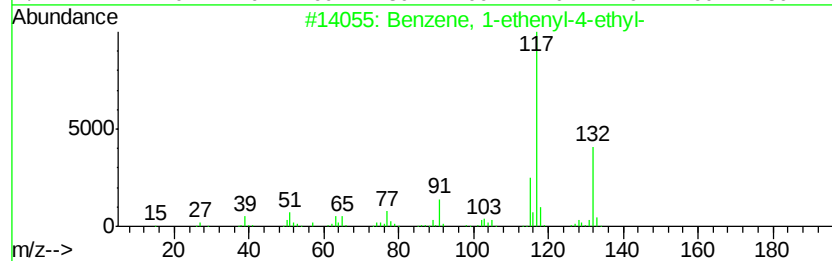
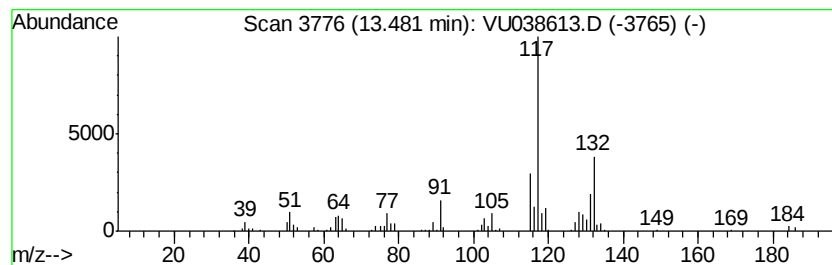
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 11 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.48	0.80 ug/L	165860	1,4-Dichlorobenzene-d4	11.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	93
2		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	93
3		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	90
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
5		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	81



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU060620\
Data File : VU038613.D
Acq On : 06 Jun 2020 13:49
Operator : SY/MD
Sample : L2884-04DL 10X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BFXD8DL

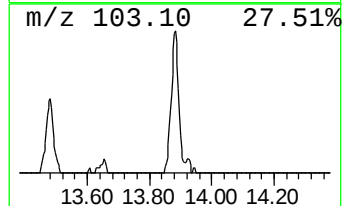
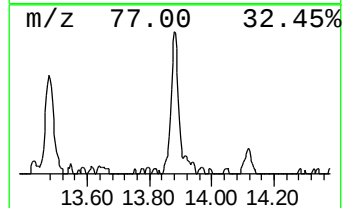
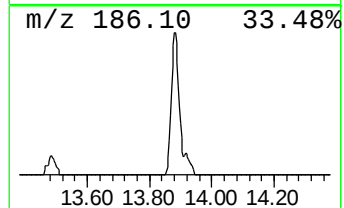
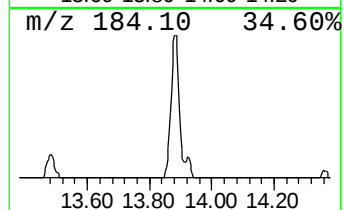
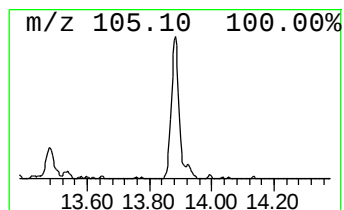
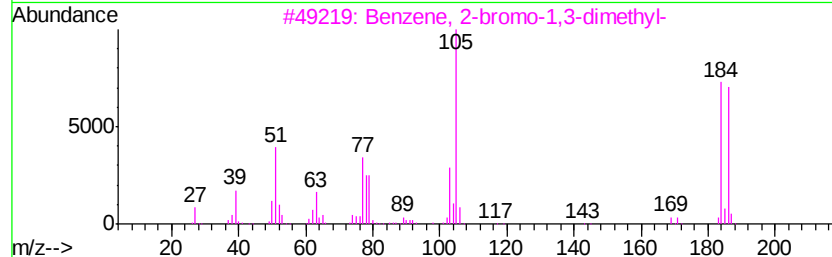
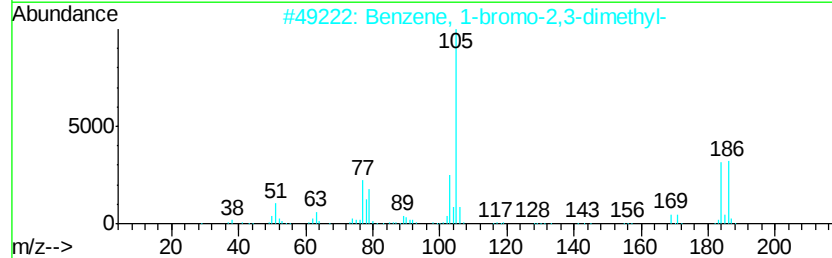
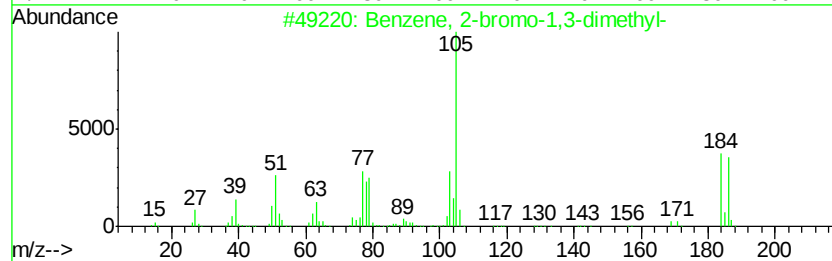
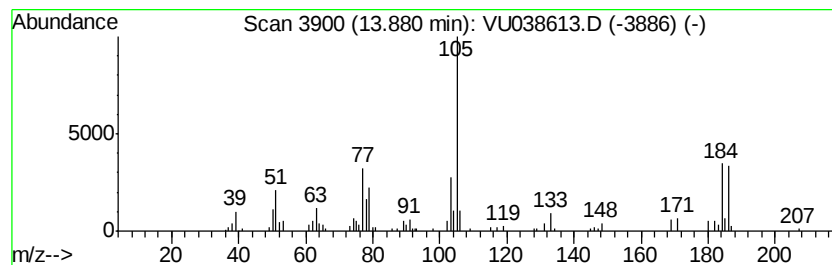
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 12 Benzene, 2-bromo-1,3-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.88	0.51 ug/L	106701	1,4-Dichlorobenzene-d4	11.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-bromo-1,3-dimethyl-	184	C8H9Br	000576-22-7	98
2		Benzene, 1-bromo-2,3-dimethyl-	184	C8H9Br	000576-23-8	98
3		Benzene, 2-bromo-1,3-dimethyl-	184	C8H9Br	000576-22-7	97
4		Benzene, 2-bromo-1,4-dimethyl-	184	C8H9Br	000553-94-6	97
5		Benzene, 1-bromo-2,4-dimethyl-	184	C8H9Br	000583-70-0	97



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU060620\
Data File : VU038613.D
Acq On : 06 Jun 2020 13:49
Operator : SY/MD
Sample : L2884-04DL 10X
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 1 Sample Multiplier: 1

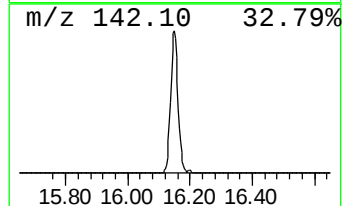
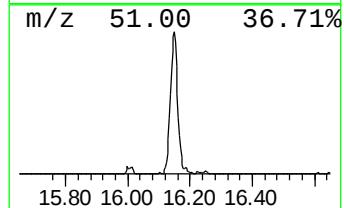
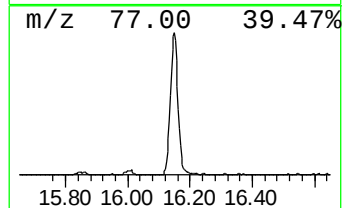
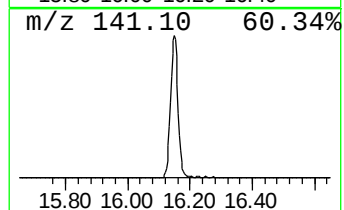
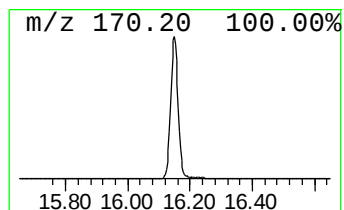
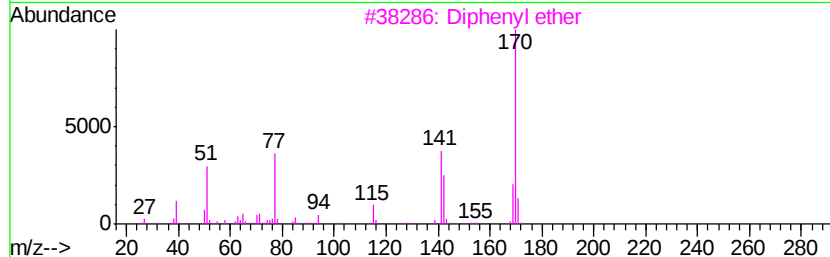
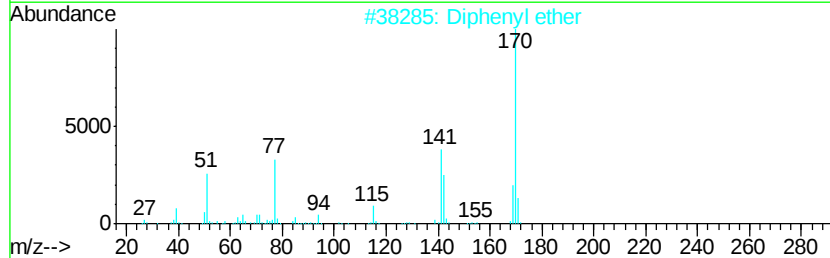
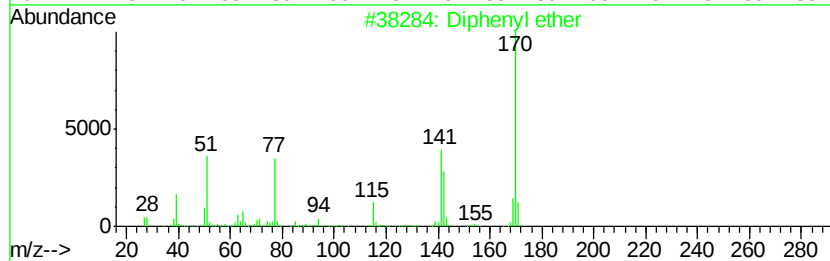
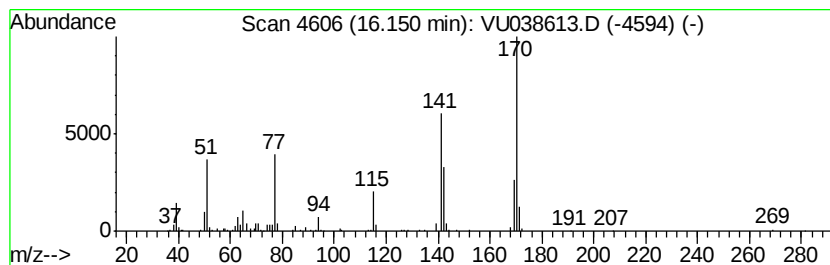
Instrument :
MSVOA_U
ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR053120WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 13 Diphenyl ether Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.15	1.02 ug/L	210912	1,4-Dichlorobenzene-d4	11.83	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diphenyl ether	170	C12H10O	000101-84-8	90
2	Diphenyl ether	170	C12H10O	000101-84-8	87
3	Diphenyl ether	170	C12H10O	000101-84-8	87
4	Spiro[cyclopropane-1,1'(4'H)-nap...	170	C12H10O	033498-24-7	76
5	Diphenyl ether	170	C12H10O	000101-84-8	64



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU060620\
 Data File : VU038613.D
 Acq On : 06 Jun 2020 13:49
 Operator : SY/MD
 Sample : L2884-04DL 10X
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BFXD8DL

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR053120WMA.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
Platinum, bis[(1,...	10.58	0.7	ug/L	140993	2	9.44	1081890	5.0
Benzene, propyl-	10.92	1.1	ug/L	224181	3	11.83	1037330	5.0
Benzene, 1-ethyl-...	11.01	1.3	ug/L	279008	3	11.83	1037330	5.0
Mesitylene	11.10	0.9	ug/L	189759	3	11.83	1037330	5.0
Benzene, 1-ethyl-...	11.30	1.0	ug/L	198824	3	11.83	1037330	5.0
Benzene, 1-etheny...	12.10	1.2	ug/L	243546	3	11.83	1037330	5.0
Benzene, 1,2,3,5-...	12.99	0.5	ug/L	108679	3	11.83	1037330	5.0
Benzene, 1-etheny...	13.48	0.8	ug/L	165860	3	11.83	1037330	5.0
Benzene, 2-bromo-...	13.88	0.5	ug/L	106701	3	11.83	1037330	5.0
Diphenyl ether	16.15	1.0	ug/L	210912	3	11.83	1037330	5.0