

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX073123\
 Data File : VX036719.D
 Acq On : 31 Jul 2023 14:47
 Operator : JC/MD
 Sample : 03808-05
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 1712

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

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\82X072123W.M
 Title : SW846 8260

Signal : TIC: VX036719.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.385	692	705	721	rBV	93210	271997	3.72%	2.138%
2	5.550	721	732	751	rVB2	153838	435997	5.96%	3.427%
3	5.958	788	799	813	rBV	100194	266306	3.64%	2.093%
4	6.763	921	931	947	rBV	251039	601379	8.23%	4.726%
5	8.647	1233	1240	1246	rBV	509680	826918	11.31%	6.499%
6	8.720	1246	1252	1273	rVB	4803177	7309909	100.00%	57.451%
7	10.055	1465	1471	1489	rBV	530850	729144	9.97%	5.731%
8	11.079	1634	1639	1653	rBV	438321	551653	7.55%	4.336%
9	12.018	1788	1793	1800	rBV	515213	678381	9.28%	5.332%
10	13.292	1993	2002	2007	rBV2	180672	262885	3.60%	2.066%
11	13.420	2018	2023	2032	rVB	166114	204446	2.80%	1.607%
12	13.926	2099	2106	2110	rBV2	84313	112143	1.53%	0.881%
13	14.213	2148	2153	2159	rBV	86092	126412	1.73%	0.994%
14	14.371	2174	2179	2184	rVB	70463	85670	1.17%	0.673%
15	14.481	2191	2197	2202	rBV2	196653	260421	3.56%	2.047%

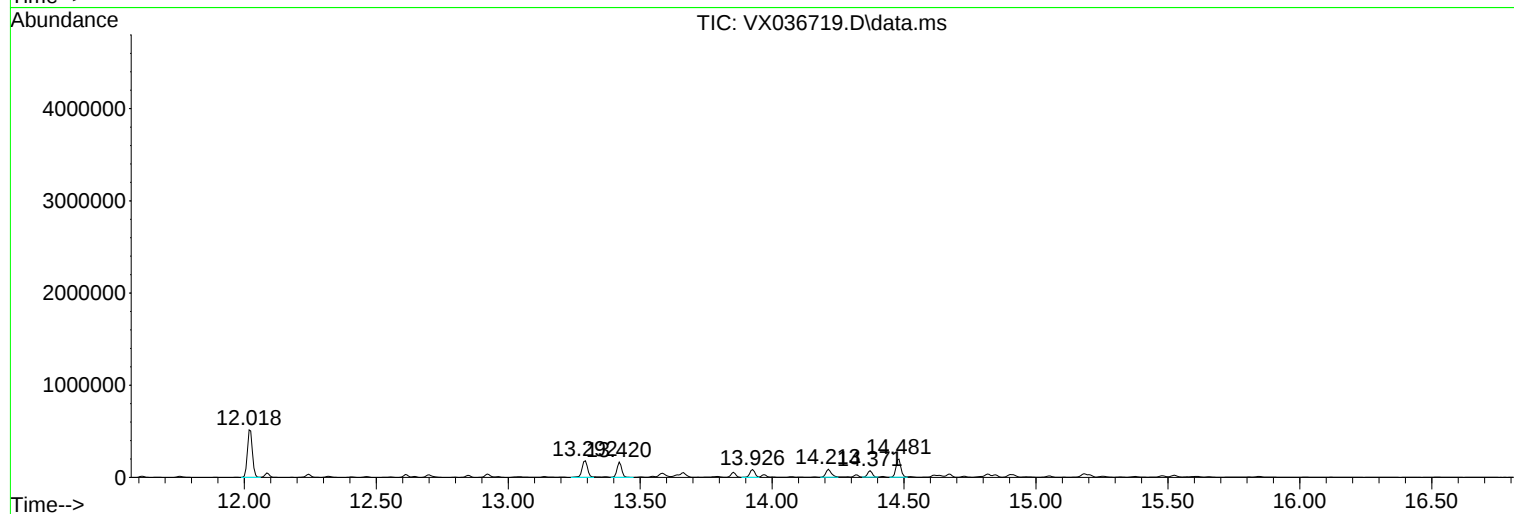
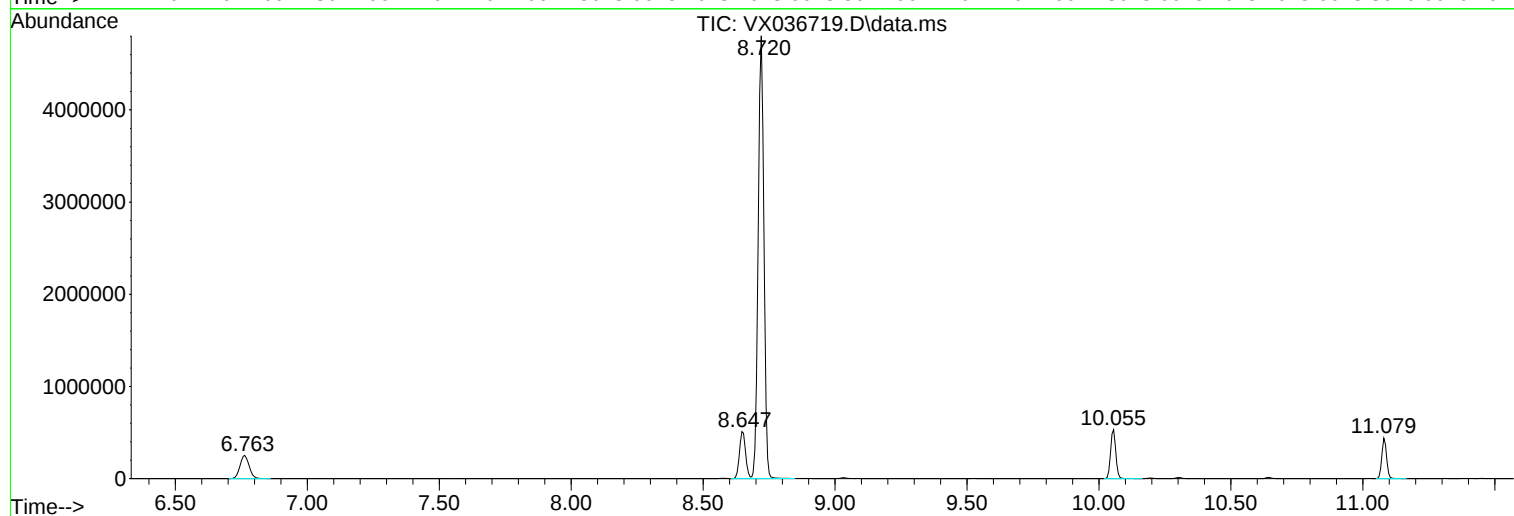
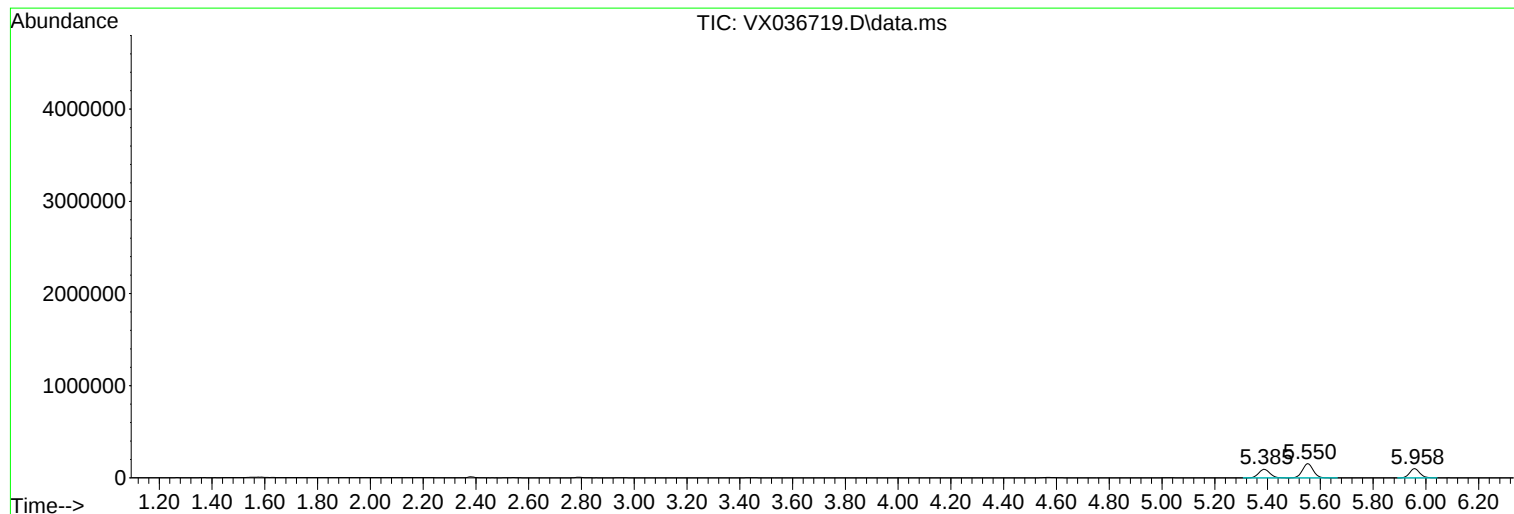
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

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



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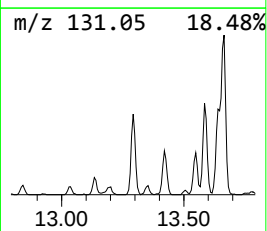
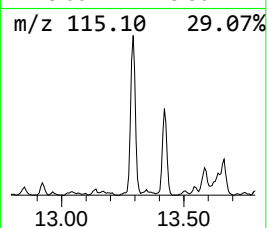
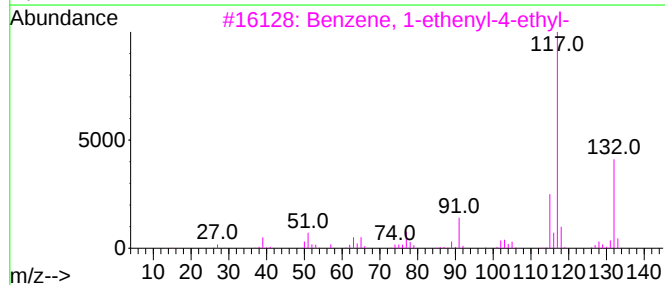
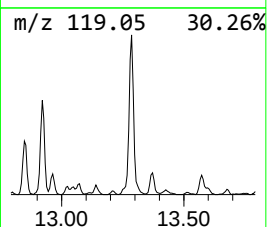
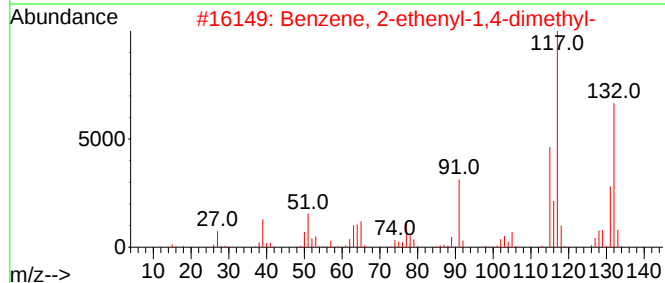
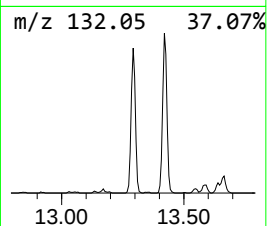
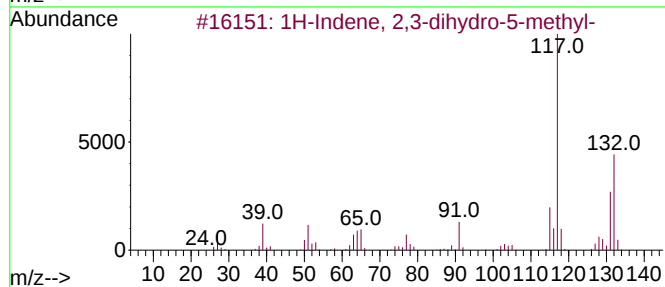
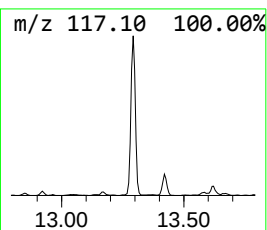
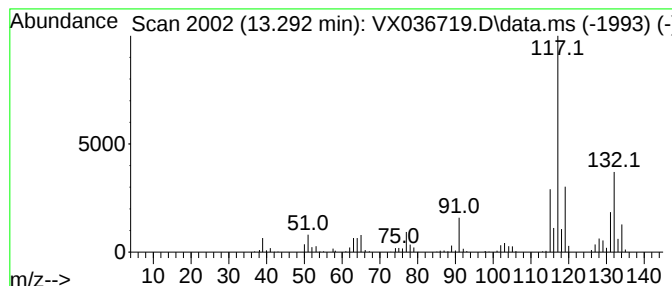
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

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

Peak Number 1 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	19.38 ug/l	262885	1,4-Dichlorobenzene-d4	12.024

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12		000874-35-1	95
2	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12		002039-89-6	95
3	Benzene, 1-ethenyl-4-ethyl-	132	C10H12		003454-07-7	94
4	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12		000824-22-6	93
5	3-Phenylbut-1-ene	132	C10H12		000934-10-1	87



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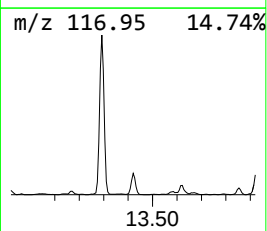
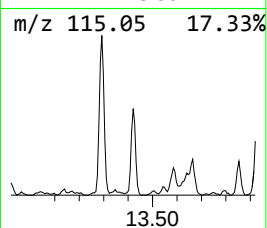
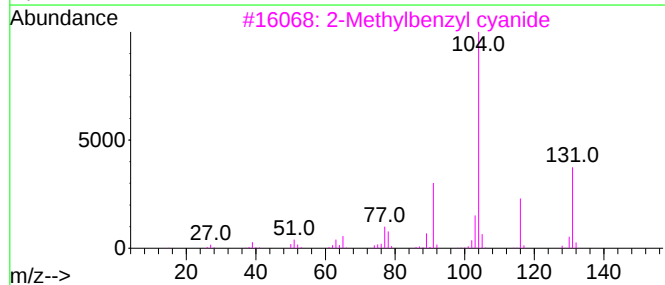
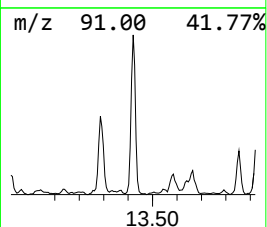
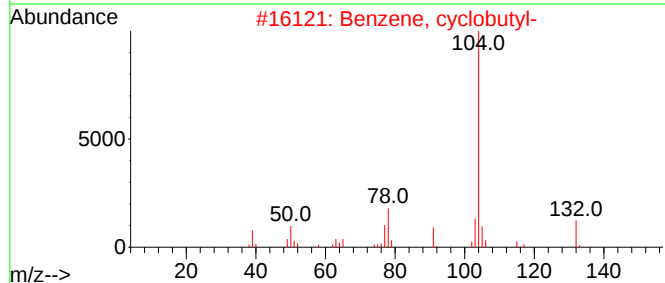
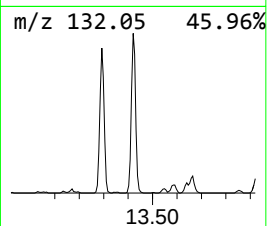
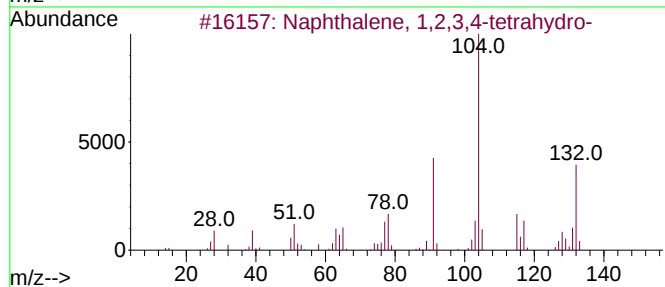
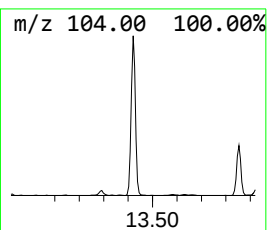
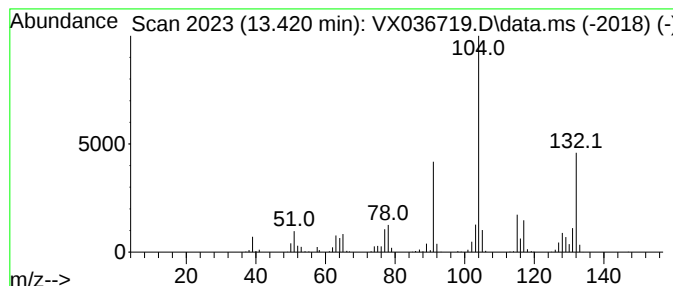
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

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TIC Integration Parameters: LSCINT.P

Peak Number 2 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.420	15.07 ug/l	204446	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
2			Benzene, cyclobutyl-	132	C10H12	004392-30-7	59
3			2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	43
4			N-Ethylbenzimidoyl bromide	211	C9H10BrN	041182-87-0	40
5			Benzenemethanol, 2-methyl-	122	C8H10O	000089-95-2	38



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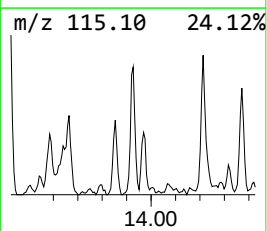
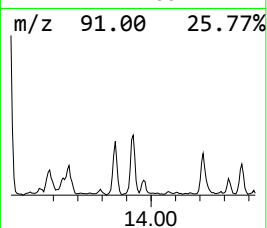
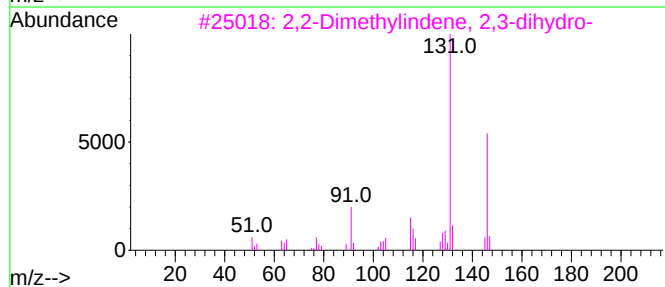
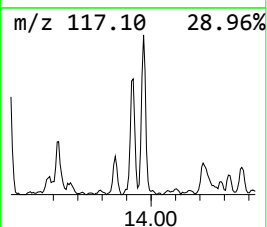
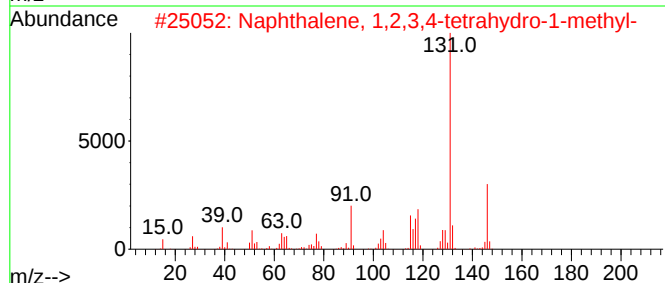
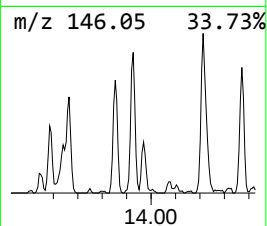
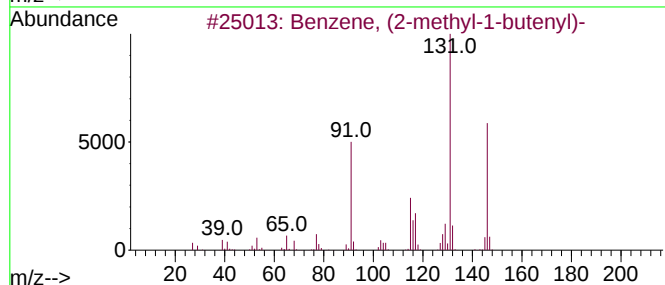
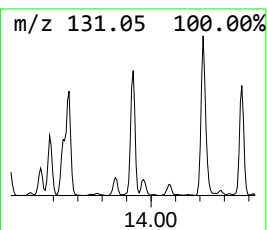
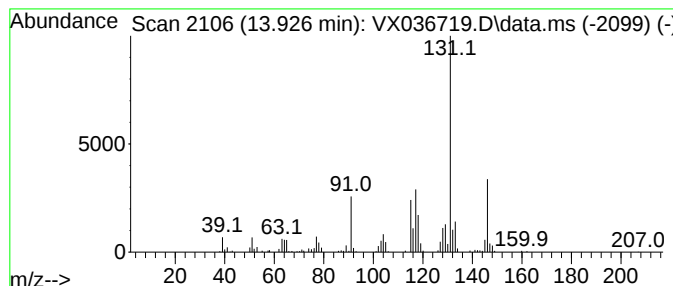
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TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, (2-methyl-1-butenyl)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.927	8.27 ug/l	112143	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	95
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	94
3			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	90
4			1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	87
5			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	76



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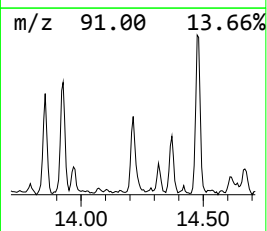
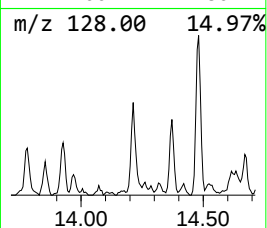
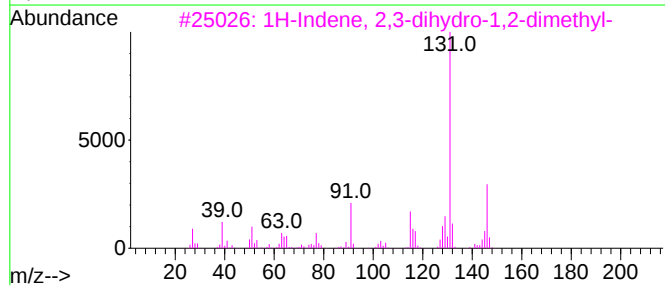
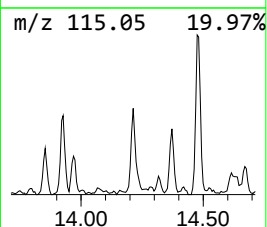
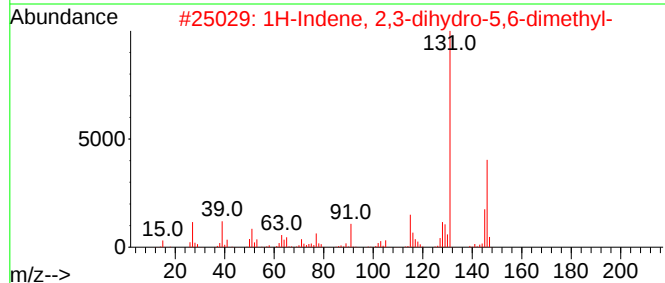
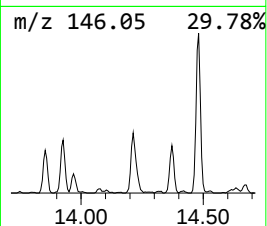
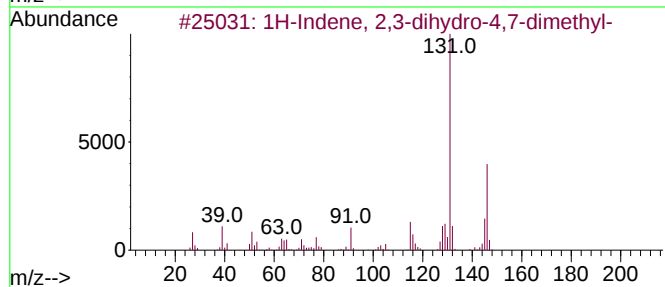
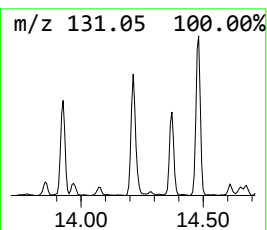
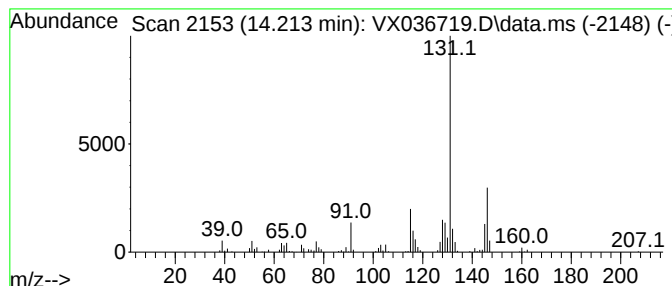
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TIC Integration Parameters: LSCINT.P

Peak Number 4 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.213	9.32 ug/l	126412	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	95
2			1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	93
3			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	90
4			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	90
5			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	90



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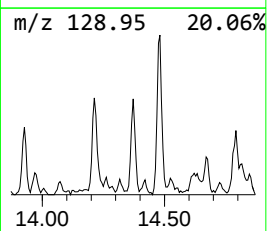
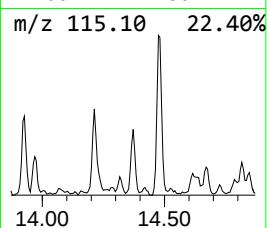
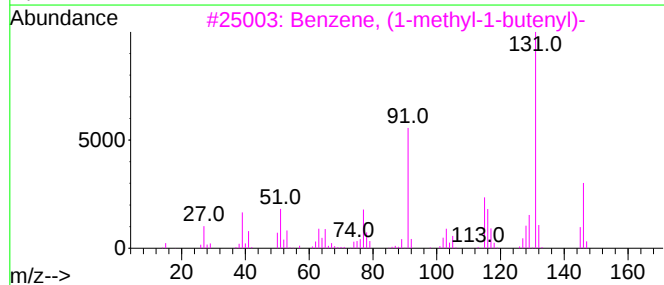
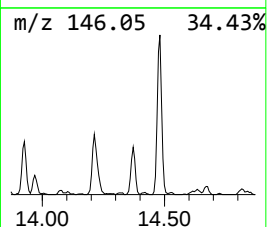
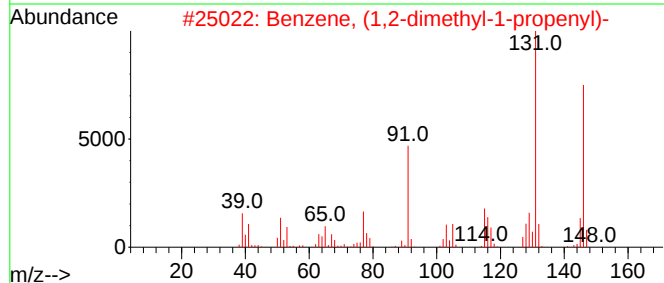
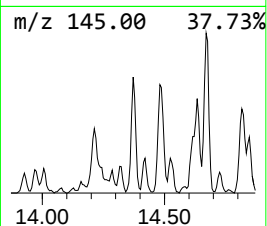
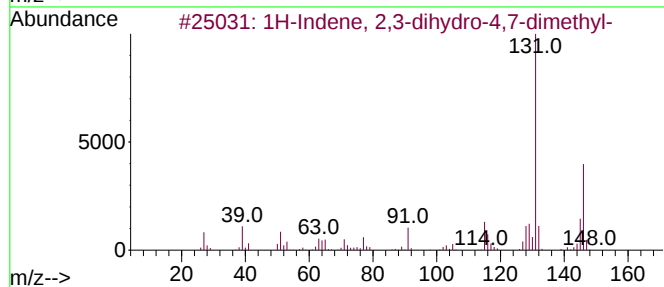
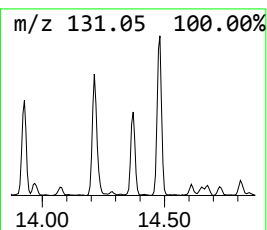
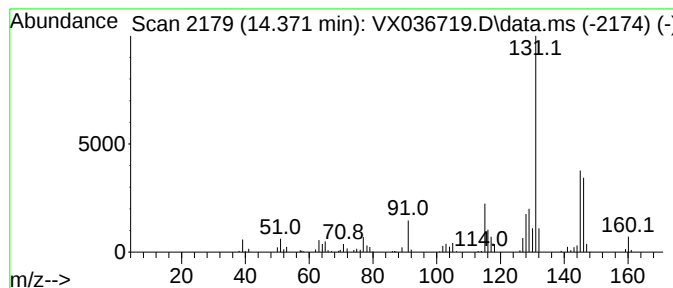
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, (1,2-dimethyl-1-pr... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.371	6.31 ug/l	85670	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	81
2			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	76
3			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	76
4			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	72
5			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	68



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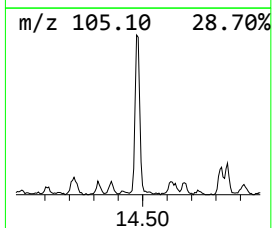
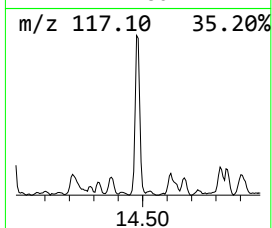
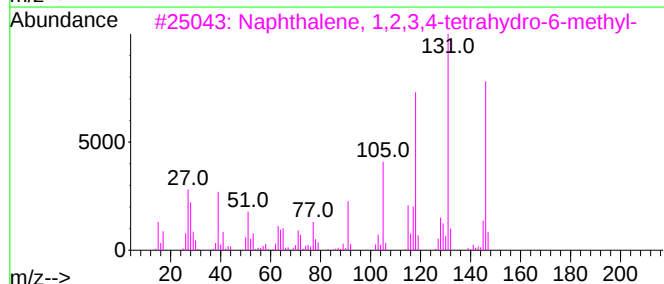
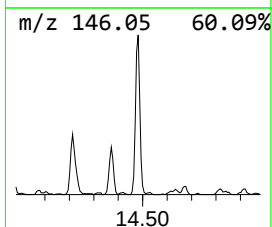
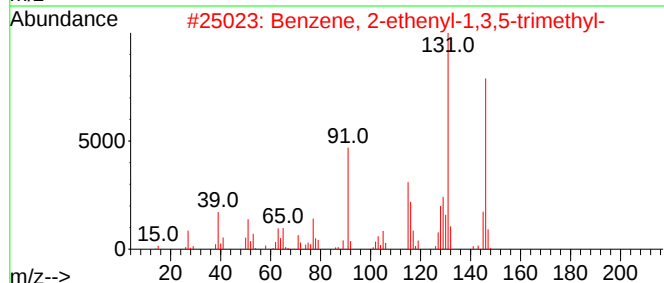
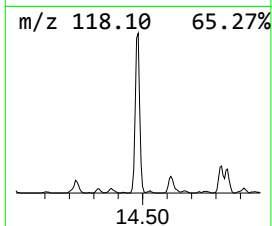
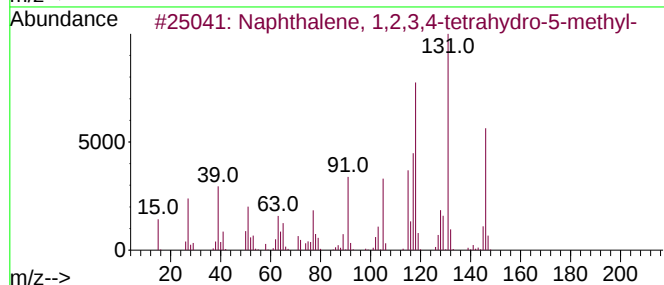
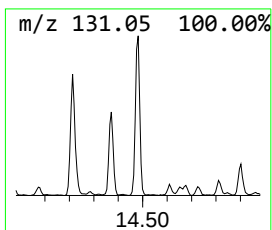
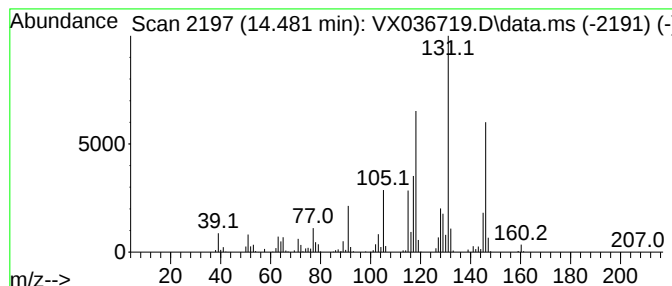
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

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

Peak Number 6 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.481	19.19 ug/l	260421	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	94
2			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	89
3			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	87
4			1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	76
5			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX073123\
Data File : VX036719.D
Acq On : 31 Jul 2023 14:47
Operator : JC/MD
Sample : 03808-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
1712

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1H-Indene, 2,3-...	13.292	19.4	ug/l	262885	4	12.024	678381	50.0
Naphthalene, 1,...	13.420	15.1	ug/l	204446	4	12.024	678381	50.0
Benzene, (2-met...	13.927	8.3	ug/l	112143	4	12.024	678381	50.0
1H-Indene, 2,3-...	14.213	9.3	ug/l	126412	4	12.024	678381	50.0
Benzene, (1,2-d...	14.371	6.3	ug/l	85670	4	12.024	678381	50.0
Naphthalene, 1,...	14.481	19.2	ug/l	260421	4	12.024	678381	50.0