

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032921\
 Data File : VX021579.D
 Acq On : 29 Mar 2021 16:02
 Operator : JC/MD
 Sample : M1785-13 10X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-2

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\82X032421W.M
 Title : SW846 8260

Signal : TIC: VX021579.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.464	732	744	756	rBV	48054	137121	11.45%	2.199%
2	5.629	761	772	783	rBV	68992	191604	16.00%	3.072%
3	6.029	829	840	852	rBV	52186	137059	11.45%	2.198%
4	6.829	966	976	989	rBV	118792	278954	23.30%	4.473%
5	8.694	1282	1293	1310	rBV	253581	396842	33.14%	6.363%
6	10.094	1525	1531	1539	rBV	279923	376562	31.45%	6.038%
7	10.236	1548	1555	1560	rBV	44304	60536	5.06%	0.971%
8	10.341	1564	1573	1579	rBV	47405	72157	6.03%	1.157%
9	10.630	1614	1622	1629	rBV5	7197	18487	1.54%	0.296%
10	10.818	1643	1654	1659	rVB3	13143	24386	2.04%	0.391%
11	10.894	1662	1667	1675	rBV3	10320	18733	1.56%	0.300%
12	11.000	1680	1685	1694	rBV	160969	201343	16.82%	3.229%
13	11.118	1700	1705	1711	rBV	245963	314623	26.28%	5.045%
14	11.342	1738	1743	1751	rBV	268627	324162	27.07%	5.198%
15	11.436	1751	1759	1763	rVV	76422	127358	10.64%	2.042%
16	11.489	1763	1768	1777	rVB	33353	49202	4.11%	0.789%
17	11.653	1785	1796	1801	rBV7	10175	20966	1.75%	0.336%
18	11.753	1805	1813	1815	rVV	23551	34973	2.92%	0.561%
19	11.789	1815	1819	1829	rVB	1012548	1197354	100.00%	19.199%
20	11.900	1829	1838	1840	rBV	35928	49163	4.11%	0.788%
21	11.930	1840	1843	1848	rVB	74404	87116	7.28%	1.397%
22	12.006	1853	1856	1860	rVV	13465	17488	1.46%	0.280%
23	12.059	1860	1865	1871	rVV	318479	399750	33.39%	6.410%
24	12.124	1871	1876	1882	rVV	288395	348516	29.11%	5.588%
25	12.218	1886	1892	1897	rVB	32861	42595	3.56%	0.683%
26	12.283	1897	1903	1910	rBV	254197	321300	26.83%	5.152%
27	12.353	1910	1915	1923	rVB2	49029	100336	8.38%	1.609%
28	12.442	1925	1930	1936	rBV	33451	42990	3.59%	0.689%
29	12.571	1946	1952	1954	rBV	37776	51743	4.32%	0.830%
30	12.595	1954	1956	1960	rVV	36407	35439	2.96%	0.568%
31	12.653	1960	1966	1969	rVV	127177	155950	13.02%	2.501%
32	12.683	1969	1971	1976	rVB	41022	44186	3.69%	0.709%
33	12.742	1976	1981	1988	rBV2	101626	170756	14.26%	2.738%
34	12.883	1999	2005	2010	rVB2	30349	46004	3.84%	0.738%
35	12.959	2010	2018	2021	rVV	62348	80677	6.74%	1.294%
36	13.001	2021	2025	2031	rVV	79784	96876	8.09%	1.553%

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Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 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\82X032421W.M
Title : SW846 8260

37	13.330	2076	2081	2087	rVB2	76521	106235	8.87%	1.703%
38	13.459	2099	2103	2110	rVB	15665	20159	1.68%	0.323%
39	13.818	2154	2164	2170	rBV	26347	36702	3.07%	0.589%

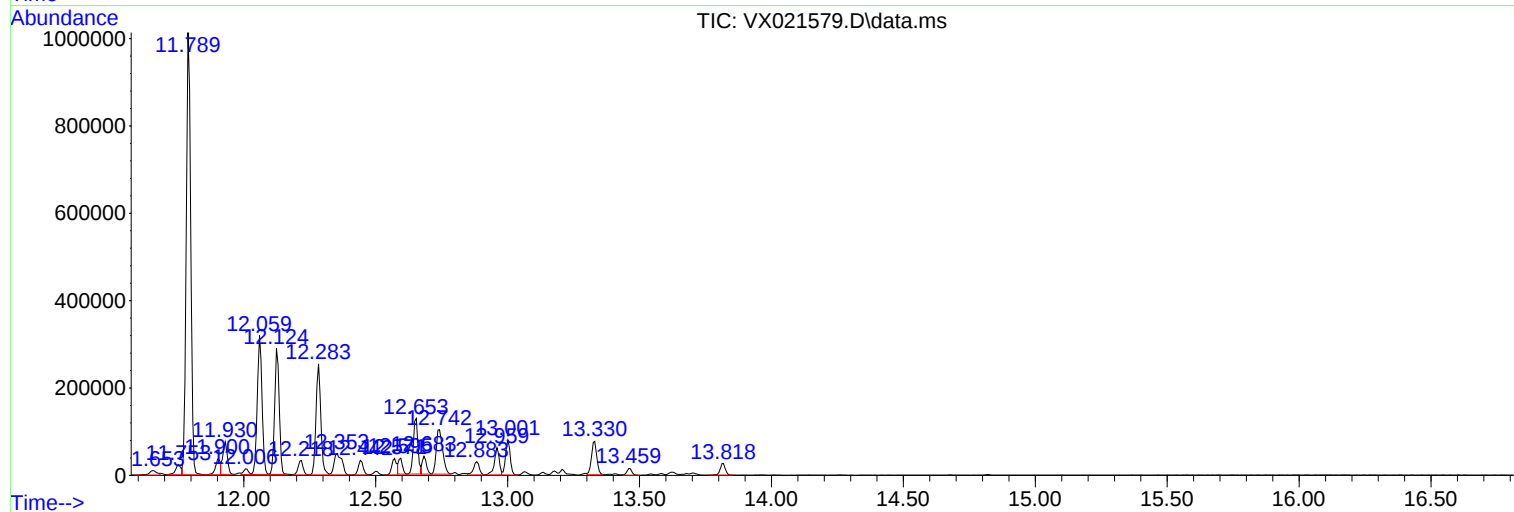
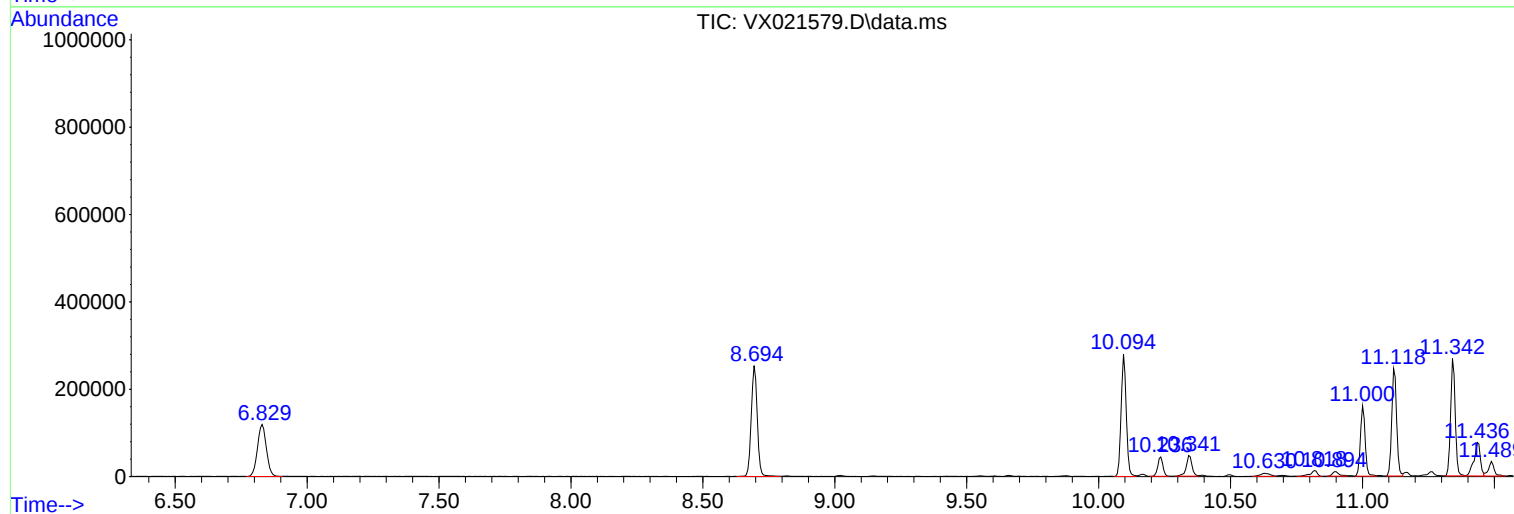
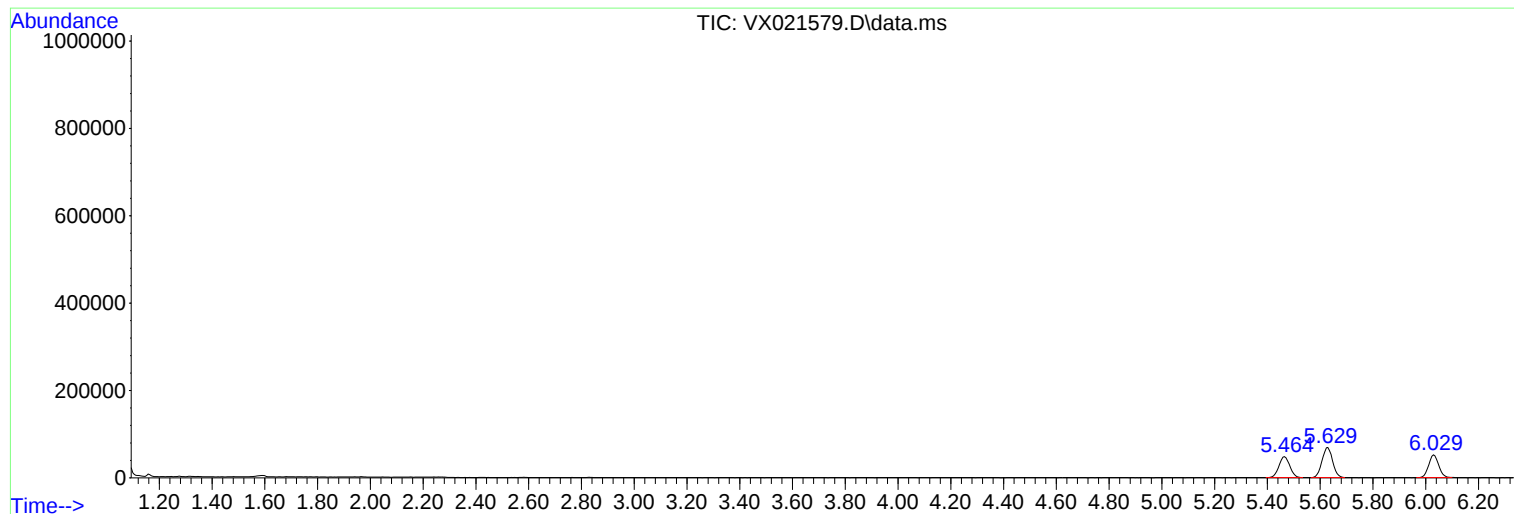
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X032421W.M
Quant Title : SW846 8260

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



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Instrument :
MSVOA_X
ClientSampleId :
MW-2

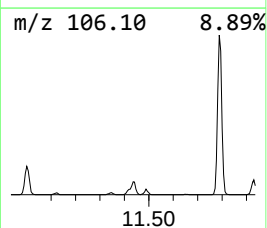
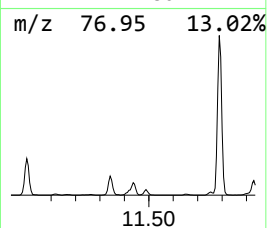
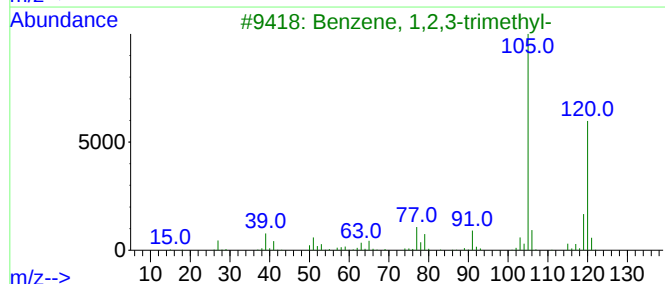
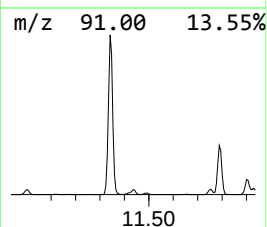
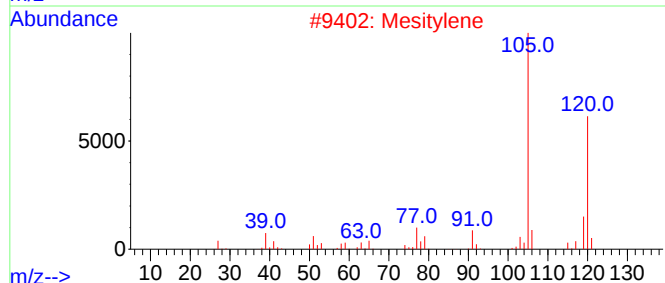
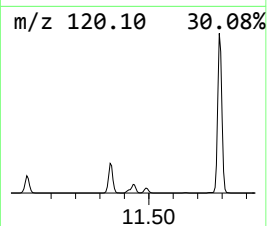
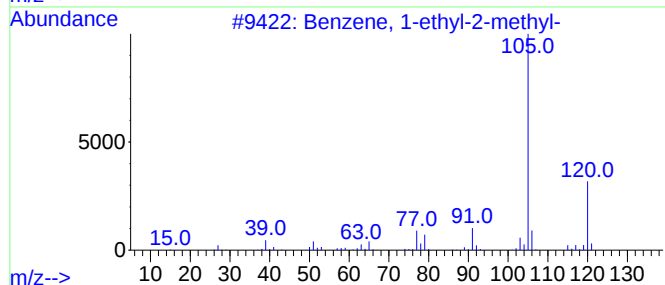
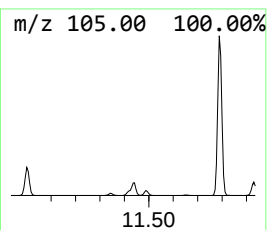
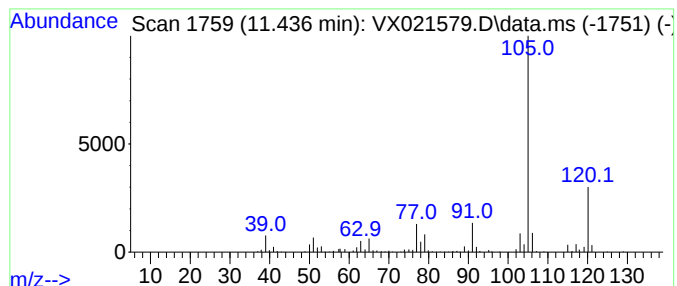
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X032421W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.436	15.93 ug/l	127358	1,4-Dichlorobenzene-d4	12.059

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Mesitylene	120	C9H12	000108-67-8	91
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2

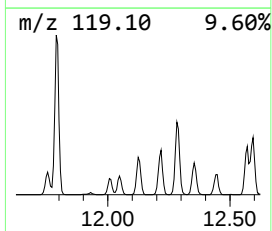
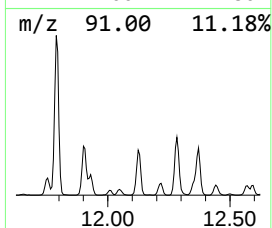
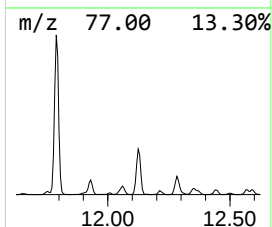
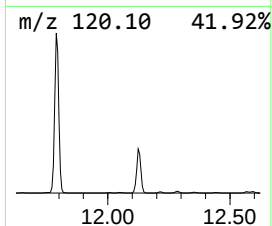
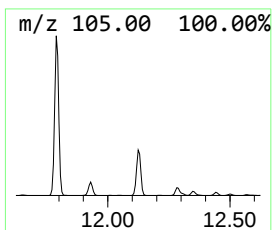
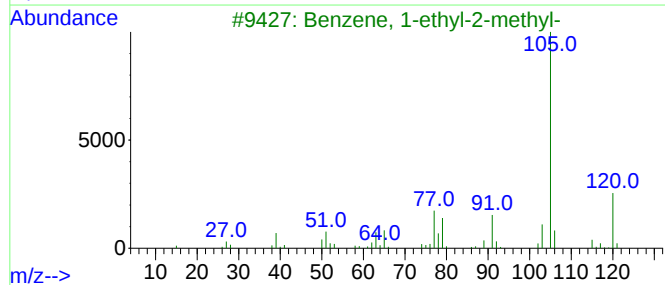
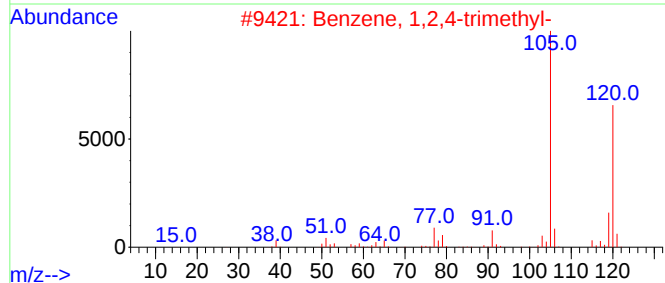
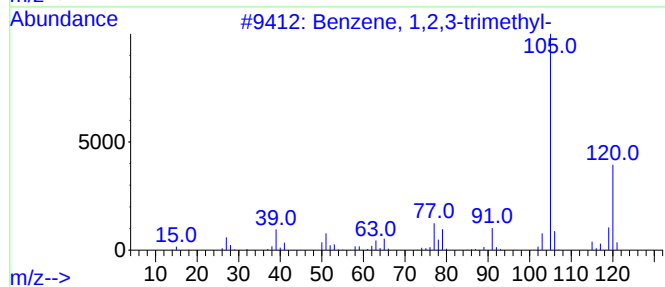
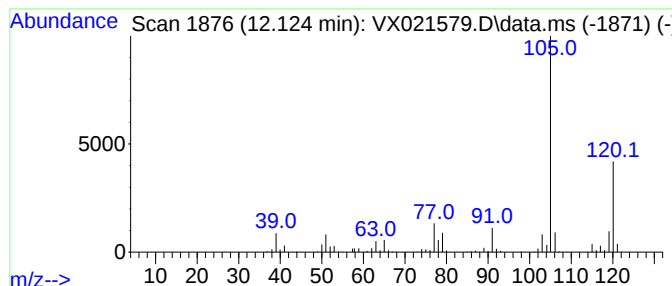
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X032421W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.124	43.59 ug/l	348516	1,4-Dichlorobenzene-d4	12.059

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			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-ethyl-2-methyl-	120	C9H12	000611-14-3	93
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032921\
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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
MW-2

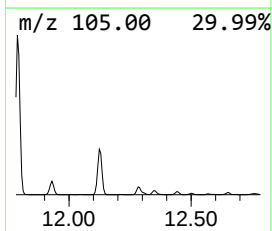
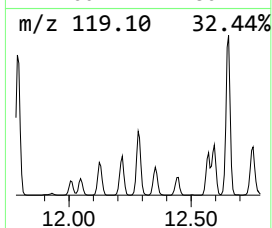
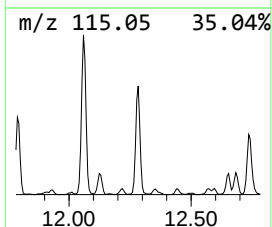
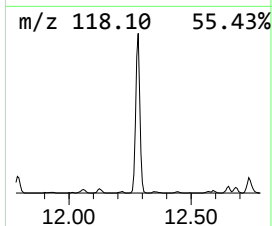
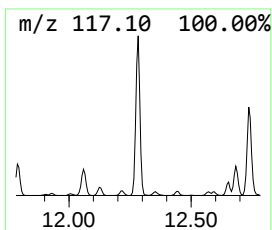
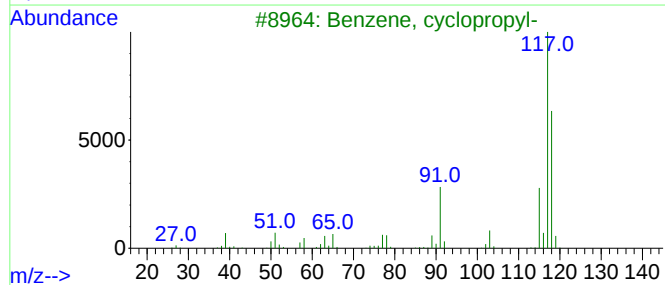
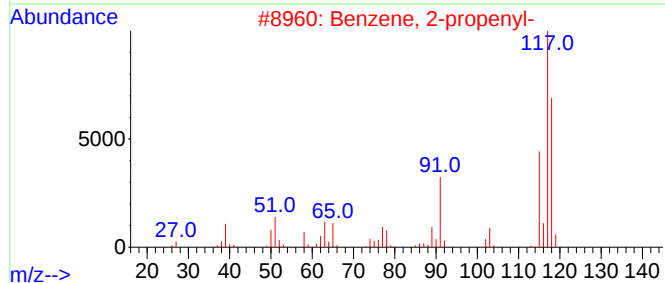
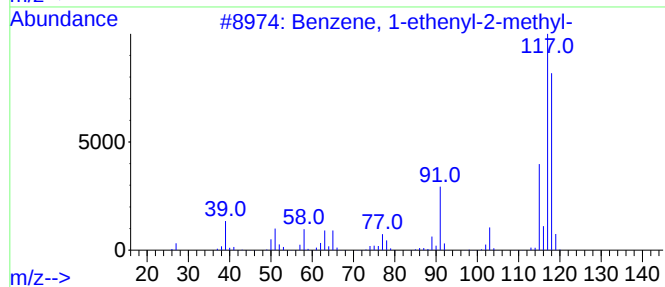
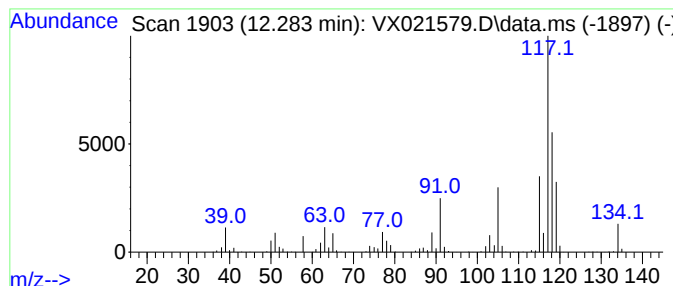
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X032421W.M
Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1-ethenyl-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.283	40.19 ug/l	321300	1,4-Dichlorobenzene-d4	12.059

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	93
2			Benzene, 2-propenyl-	118	C9H10	000300-57-2	70
3			Benzene, cyclopropyl-	118	C9H10	000873-49-4	70
4			Indane	118	C9H10	000496-11-7	64
5			Benzene, 1-propenyl-	118	C9H10	000637-50-3	64



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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2

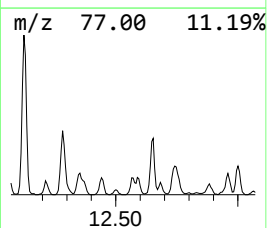
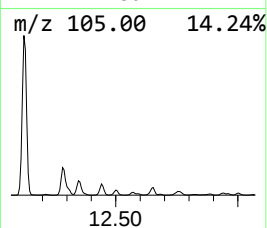
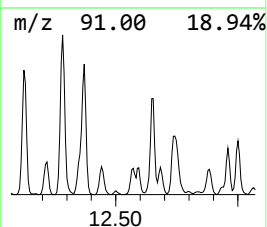
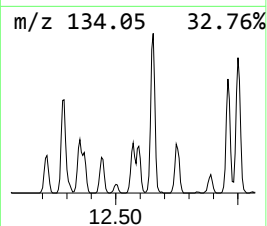
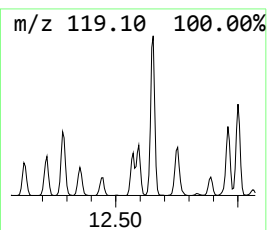
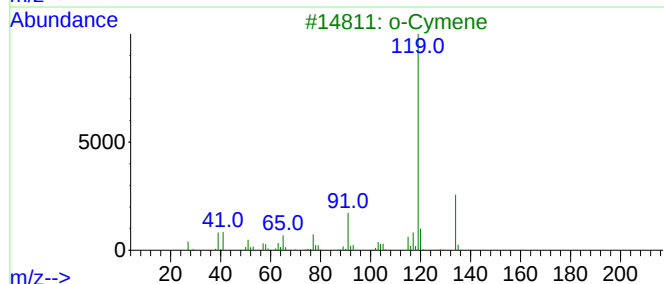
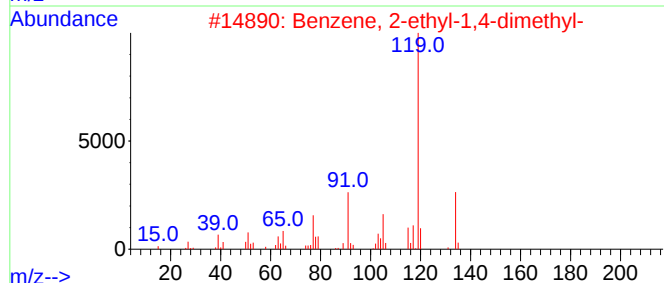
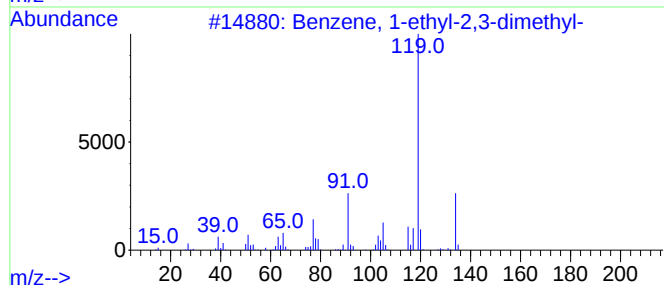
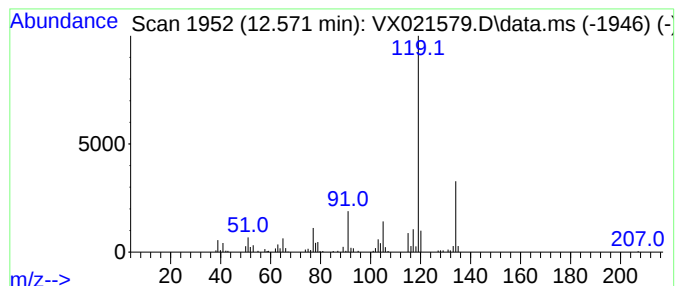
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X032421W.M
Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.571	6.47 ug/l	51743	1,4-Dichlorobenzene-d4	12.059

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3			o-Cymene	134	C10H14	000527-84-4	95
4			p-Cymene	134	C10H14	000099-87-6	95
5			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95



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Instrument :
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ClientSampleId :
MW-2

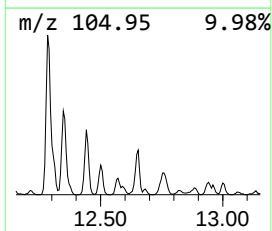
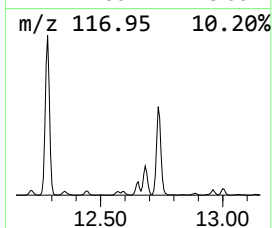
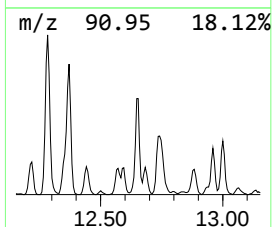
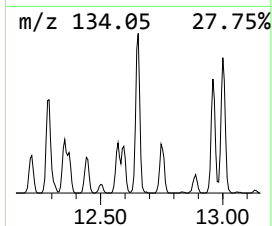
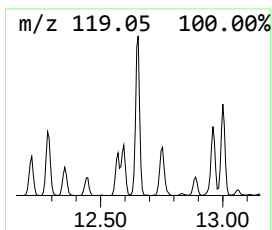
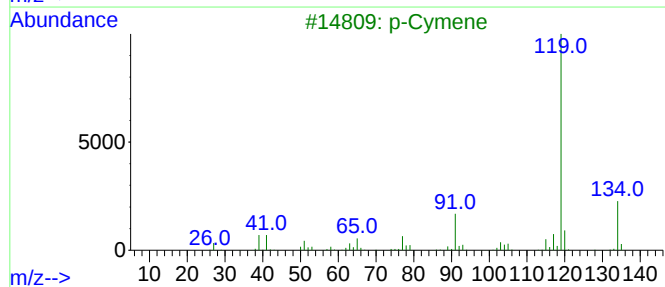
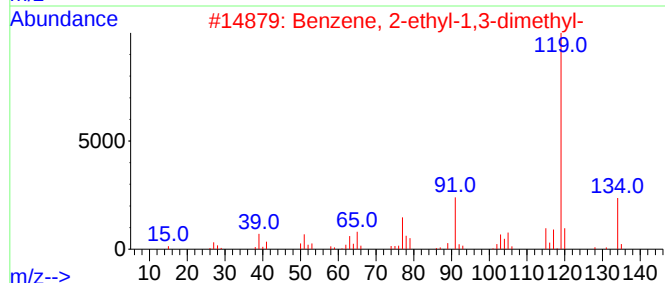
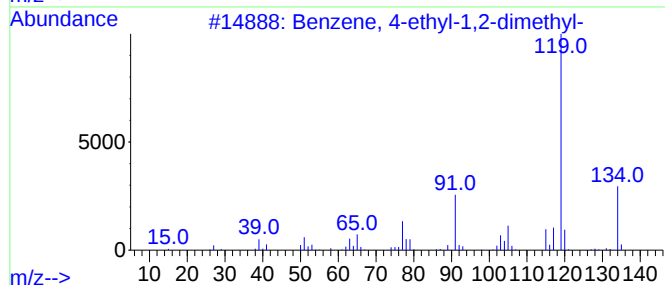
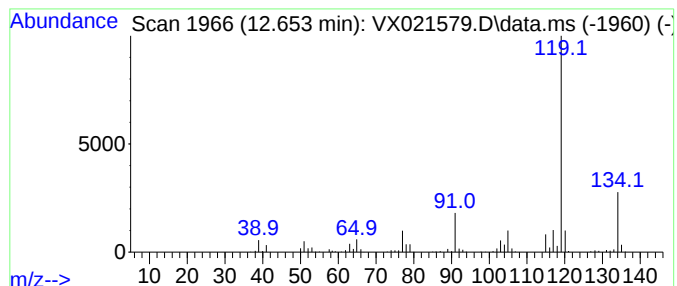
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.653	19.51 ug/l	155950	1,4-Dichlorobenzene-d4	12.059

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
3			p-Cymene	134	C10H14	000099-87-6	95
4			o-Cymene	134	C10H14	000527-84-4	95
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032921\
Data File : VX021579.D
Acq On : 29 Mar 2021 16:02
Operator : JC/MD
Sample : M1785-13 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2

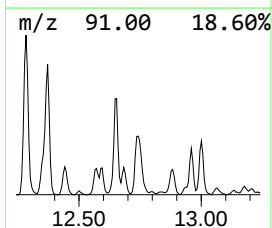
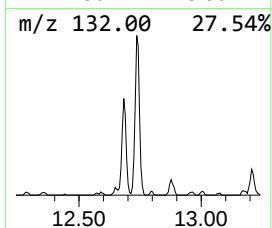
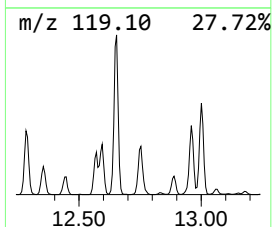
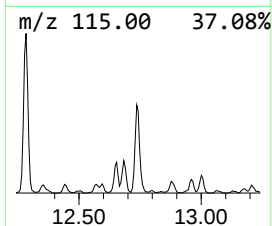
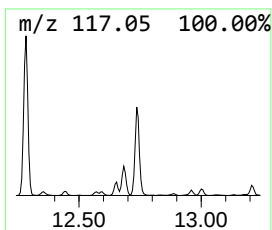
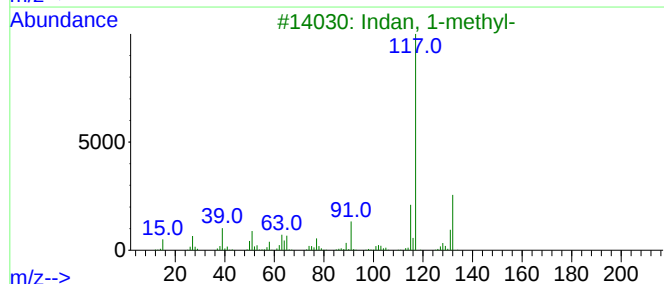
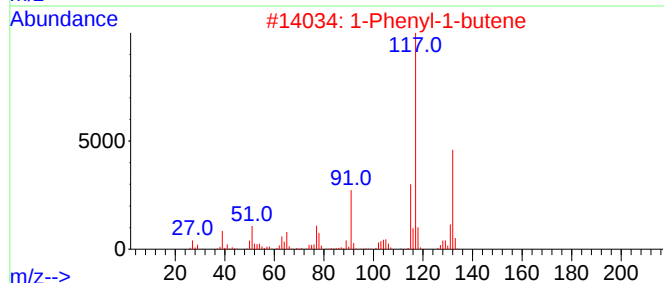
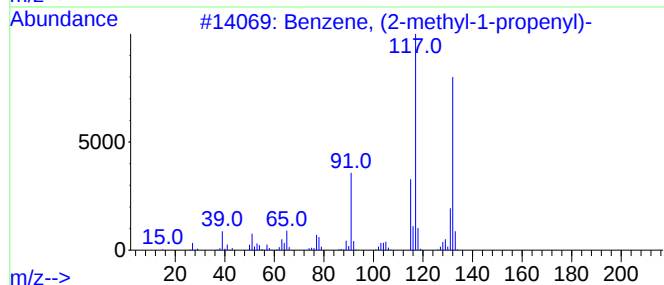
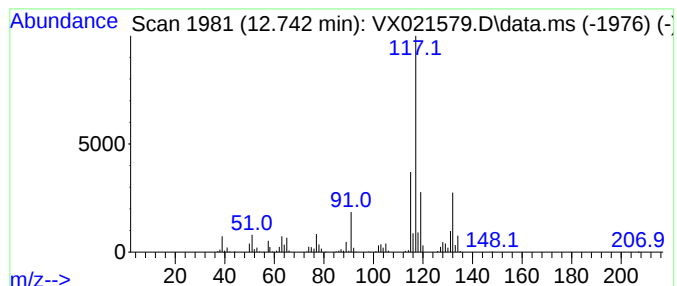
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, (2-methyl-1-propen... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.742	21.36 ug/l	170756	1,4-Dichlorobenzene-d4	12.059

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	83
2			1-Phenyl-1-butene	132	C10H12	000824-90-8	81
3			Indan, 1-methyl-	132	C10H12	000767-58-8	76
4			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	74
5			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032921\
Data File : VX021579.D
Acq On : 29 Mar 2021 16:02
Operator : JC/MD
Sample : M1785-13 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2

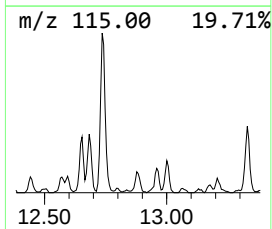
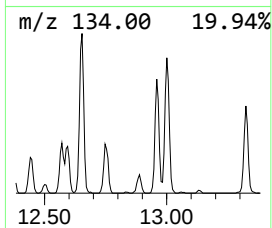
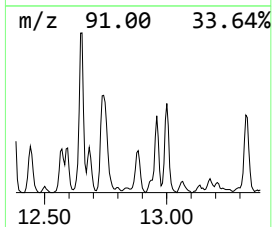
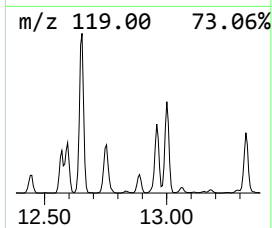
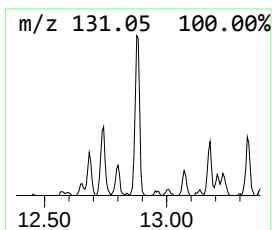
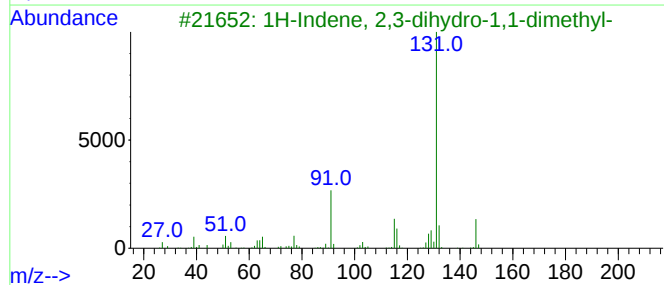
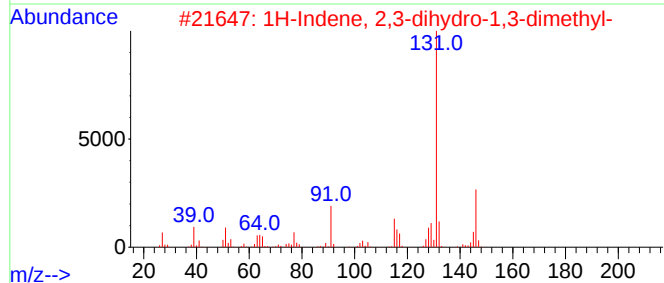
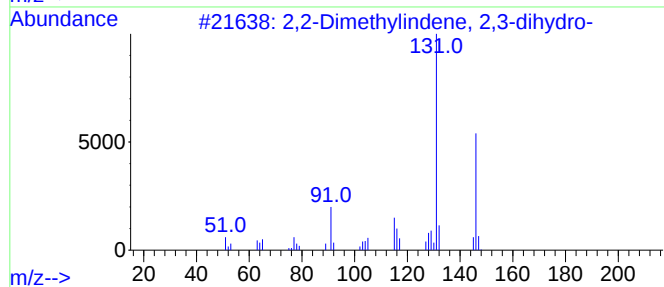
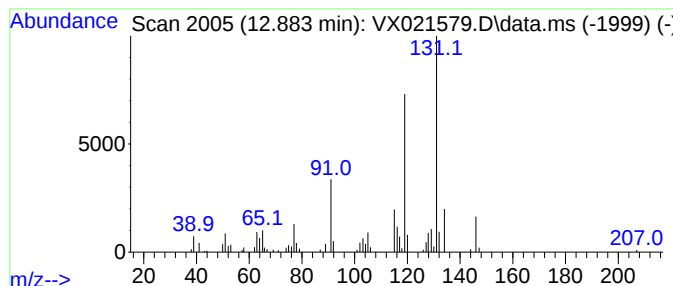
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Quant Title : SW846 8260

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

Peak Number 7 2,2-Dimethylindene, 2,3-dih... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.883	5.75 ug/l	46004	1,4-Dichlorobenzene-d4	12.059

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	94
2			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	86
3			1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	64
4			Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	58
5			Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032921\
Data File : VX021579.D
Acq On : 29 Mar 2021 16:02
Operator : JC/MD
Sample : M1785-13 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2

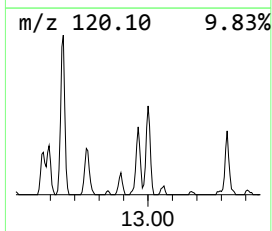
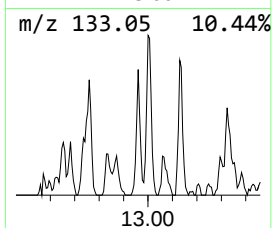
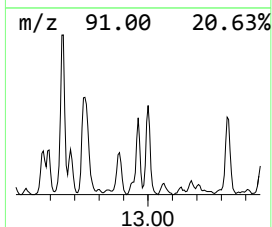
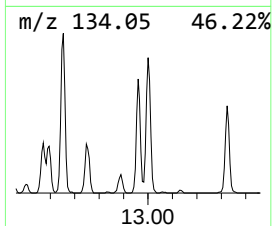
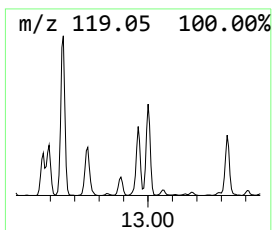
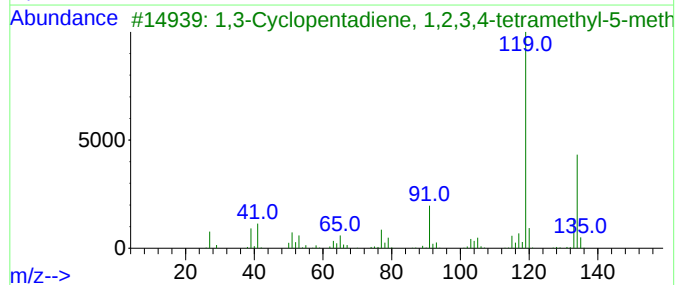
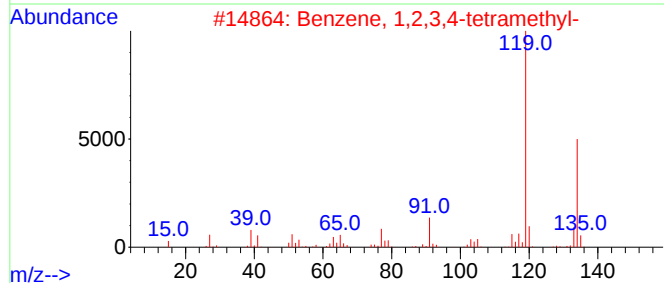
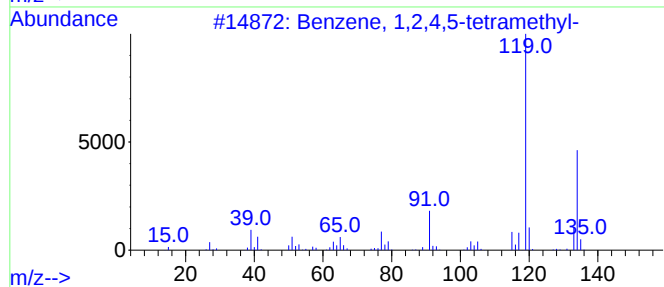
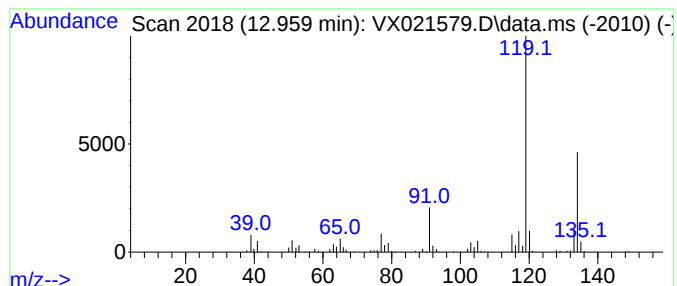
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X032421W.M
Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.959	10.09 ug/l	80677	1,4-Dichlorobenzene-d4	12.059

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
3			1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	94
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
5			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032921\
Data File : VX021579.D
Acq On : 29 Mar 2021 16:02
Operator : JC/MD
Sample : M1785-13 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2

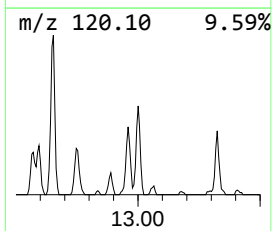
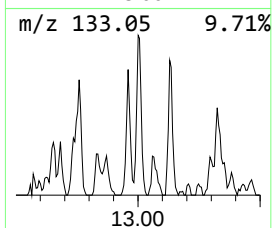
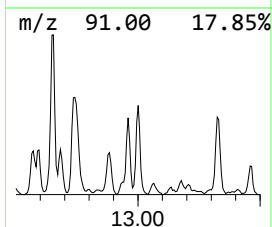
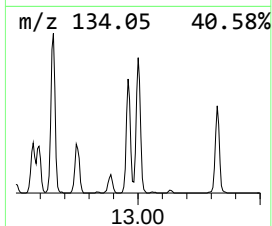
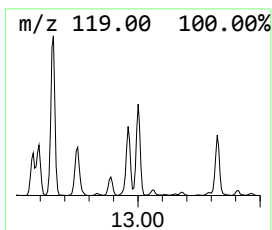
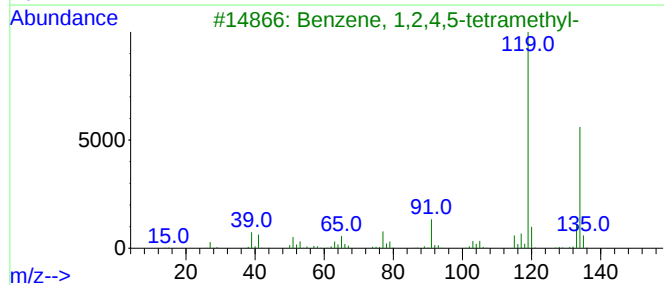
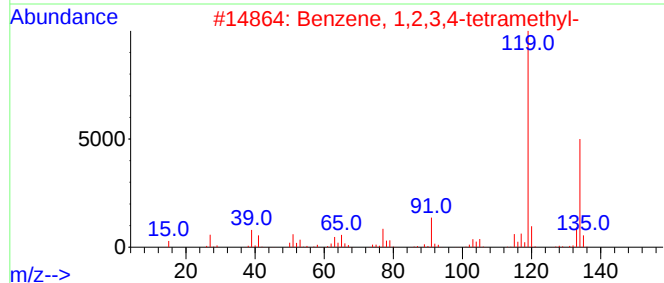
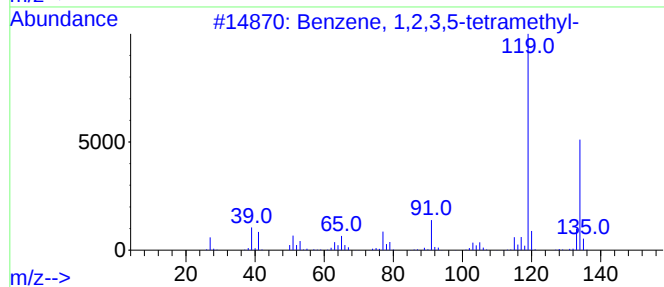
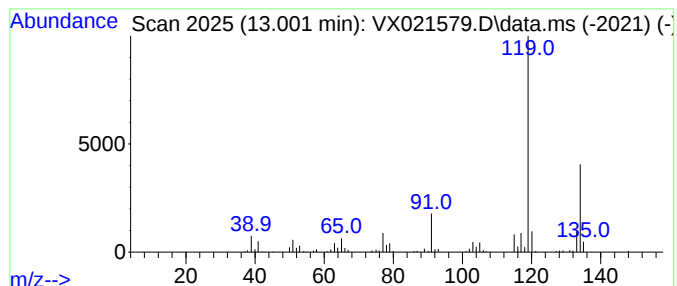
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X032421W.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.001	12.12 ug/l	96876	1,4-Dichlorobenzene-d4	12.059

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
3			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
4			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032921\
Data File : VX021579.D
Acq On : 29 Mar 2021 16:02
Operator : JC/MD
Sample : M1785-13 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2

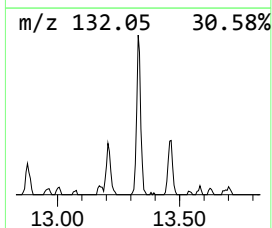
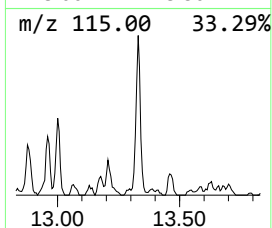
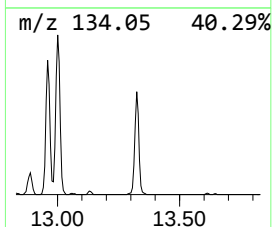
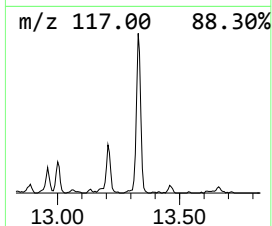
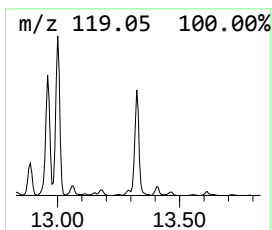
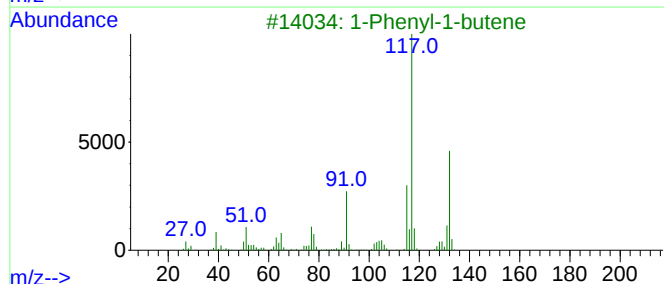
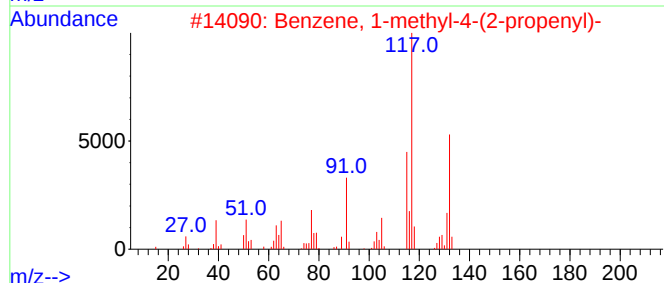
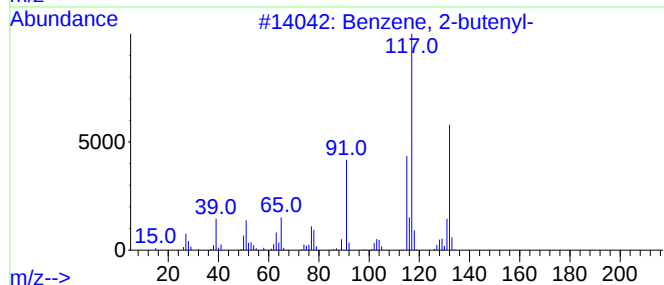
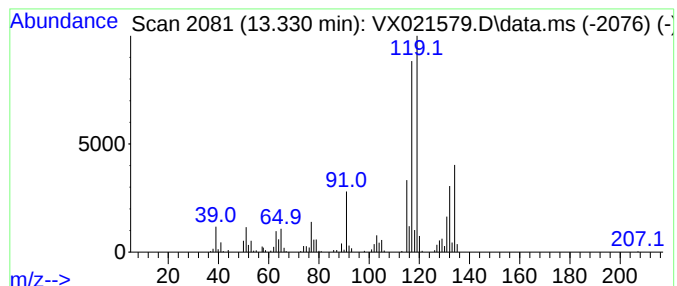
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Quant Title : SW846 8260

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

Peak Number 10 Benzene, 2-butenyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.330	13.29 ug/l	106235	1,4-Dichlorobenzene-d4	12.059

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-butenyl-	132	C10H12	001560-06-1	89
2			Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	70
3			1-Phenyl-1-butene	132	C10H12	000824-90-8	55
4			3-Phenylbut-1-ene	132	C10H12	000934-10-1	55
5			Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032921\
Data File : VX021579.D
Acq On : 29 Mar 2021 16:02
Operator : JC/MD
Sample : M1785-13 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X032421W.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Benzene, 1,2,3-...	12.124	43.6	ug/l	348516	4	12.059	399750	50.0
Benzene, 1-ethe...	12.283	40.2	ug/l	321300	4	12.059	399750	50.0
Benzene, 1-ethy...	12.571	6.5	ug/l	51743	4	12.059	399750	50.0
Benzene, 4-ethy...	12.653	19.5	ug/l	155950	4	12.059	399750	50.0
Benzene, (2-met...	12.742	21.4	ug/l	170756	4	12.059	399750	50.0
2,2-Dimethylind...	12.883	5.8	ug/l	46004	4	12.059	399750	50.0
Benzene, 1,2,4,...	12.959	10.1	ug/l	80677	4	12.059	399750	50.0
Benzene, 1,2,3,...	13.001	12.1	ug/l	96876	4	12.059	399750	50.0
Benzene, 2-bute...	13.330	13.3	ug/l	106235	4	12.059	399750	50.0