

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX102220\
 Data File : VX019054.D
 Acq On : 23 Oct 2020 02:27
 Operator : JC/SP
 Sample : L4417-17
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-10

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X102120W.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.776	108	113	130	rVB	3759045	5004216	5.86%	1.109%
2	1.983	142	147	159	rVB	1349302	1938872	2.27%	0.430%
3	2.380	200	212	235	rVB	1525944	3041593	3.56%	0.674%
4	2.825	278	285	289	rBV	1942993	4269967	5.00%	0.946%
5	2.873	289	293	303	rVV	2811852	5848142	6.85%	1.296%
6	3.154	327	339	357	rVB	1972188	4458291	5.22%	0.988%
7	3.794	436	444	453	rVB	398876	963432	1.13%	0.213%
8	4.282	515	524	531	rVV	314877	965104	1.13%	0.214%
9	4.379	531	540	551	rVB	1858123	5284513	6.19%	1.171%
10	5.245	659	682	701	rVB7	166635	1049569	1.23%	0.233%
11	5.562	724	734	743	rBV5	660246	2703639	3.17%	0.599%
12	5.696	750	756	778	rVB3	629320	2252234	2.64%	0.499%
13	5.903	778	790	802	rBV2	700480	2066773	2.42%	0.458%
14	6.232	828	844	851	rBV3	328449	1244151	1.46%	0.276%
15	6.324	851	859	869	rVV2	956142	2943092	3.45%	0.652%
16	6.434	869	877	889	rVB	371318	1039252	1.22%	0.230%
17	6.830	934	942	948	rBV	468170	1193377	1.40%	0.264%
18	6.915	948	956	967	rVB3	480448	1546929	1.81%	0.343%
19	7.232	1001	1008	1016	rBV	452665	971716	1.14%	0.215%
20	7.440	1025	1042	1053	rBV4	956333	2635970	3.09%	0.584%
21	7.933	1112	1123	1132	rBV3	239999	880945	1.03%	0.195%
22	8.037	1132	1140	1151	rVB	572721	1198790	1.40%	0.266%
23	8.159	1151	1160	1163	rBV	1026666	2076769	2.43%	0.460%
24	8.189	1163	1165	1170	rVB	658347	958474	1.12%	0.212%
25	8.549	1217	1224	1234	rVB2	523375	982430	1.15%	0.218%
26	8.695	1242	1248	1255	rVB2	1268014	2159583	2.53%	0.478%
27	10.098	1474	1478	1482	rVB	976908	1255063	1.47%	0.278%
28	10.238	1496	1501	1507	rBV	18912543	24776996	29.01%	5.489%
29	10.354	1512	1520	1525	rBV	33717306	46755930	54.75%	10.358%
30	10.683	1569	1574	1584	rVB	1100541	1392349	1.63%	0.308%
31	11.006	1619	1627	1632	rBV	18051409	22599225	26.46%	5.006%
32	11.122	1641	1646	1651	rBV	975180	1191533	1.40%	0.264%
33	11.347	1667	1683	1690	rBV	17894339	22459358	26.30%	4.975%
34	11.445	1690	1699	1703	rBV	21261119	27250223	31.91%	6.037%

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Title : SW846 8260

35	11.494	1703	1707	1712	rVB	23488388	29792592	34.89%	6.600%
36	11.658	1728	1734	1744	rVB	17502392	22825401	26.73%	5.056%
37	11.805	1751	1758	1763	rBV2	51800676	85395096	100.00%	18.917%
38	11.933	1776	1779	1785	rVB	719232	871585	1.02%	0.193%
39	12.061	1795	1800	1806	rVB2	1387298	1892363	2.22%	0.419%
40	12.128	1806	1811	1816	rBV	16882381	21222641	24.85%	4.701%
41	12.286	1831	1837	1844	rBV	13647878	17244973	20.19%	3.820%
42	12.359	1844	1849	1856	rVV2	5487812	8262603	9.68%	1.830%
43	12.500	1868	1872	1878	rVB	1566451	1871224	2.19%	0.415%
44	12.573	1878	1884	1886	rBV	3211635	4405574	5.16%	0.976%
45	12.597	1886	1888	1892	rVB	2698339	2717784	3.18%	0.602%
46	12.652	1892	1897	1901	rBV	5334698	6207548	7.27%	1.375%
47	12.737	1907	1911	1920	rBV2	1862039	2762906	3.24%	0.612%
48	12.890	1931	1936	1941	rVB	1068018	1319175	1.54%	0.292%
49	12.963	1941	1948	1951	rBV	3256691	3953244	4.63%	0.876%
50	13.000	1951	1954	1960	rVB	4519341	5495119	6.43%	1.217%
51	13.207	1984	1988	1993	rVB	2334057	2709488	3.17%	0.600%
52	13.329	2004	2008	2014	rVB2	4702725	6516133	7.63%	1.443%
53	13.463	2025	2030	2038	rVB	776521	952517	1.12%	0.211%
54	13.816	2082	2088	2093	rBV	9048148	10921422	12.79%	2.419%
55	14.676	2217	2229	2240	rBV	3529913	4291238	5.03%	0.951%
56	14.816	2247	2252	2258	rBV	1891875	2427178	2.84%	0.538%

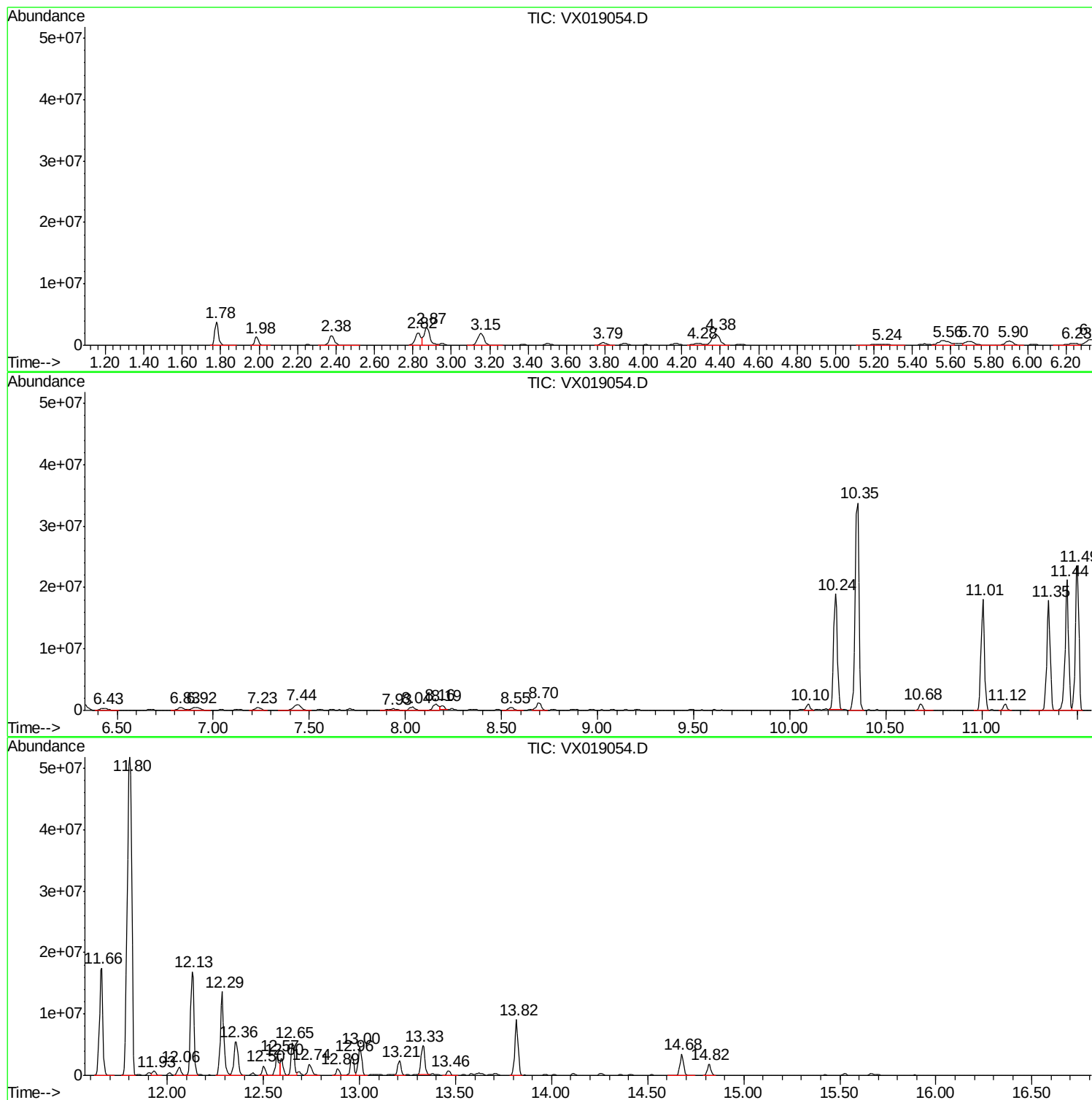
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Quant Title : SW846 8260

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



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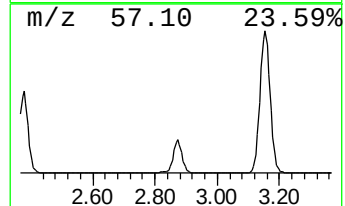
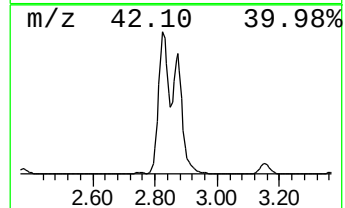
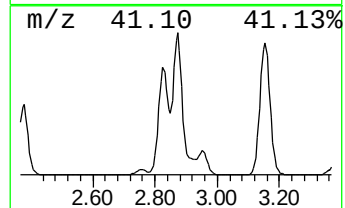
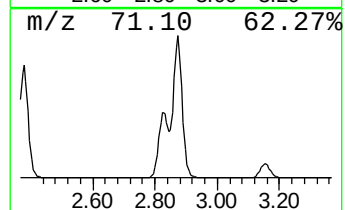
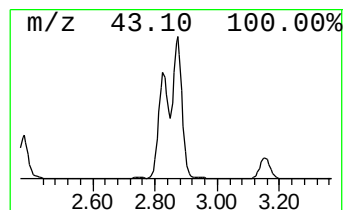
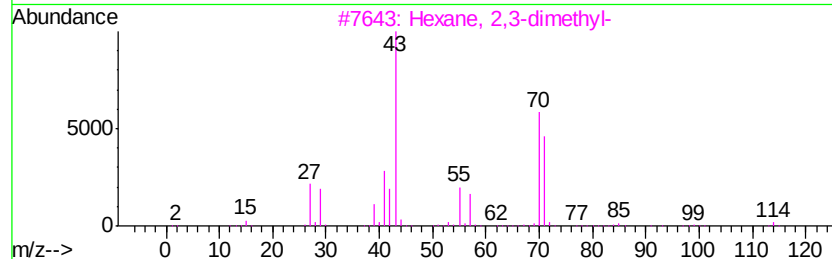
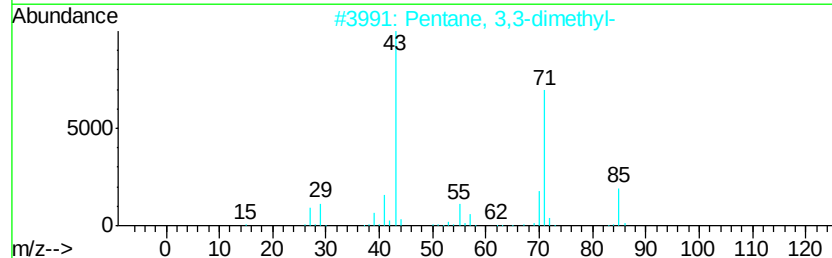
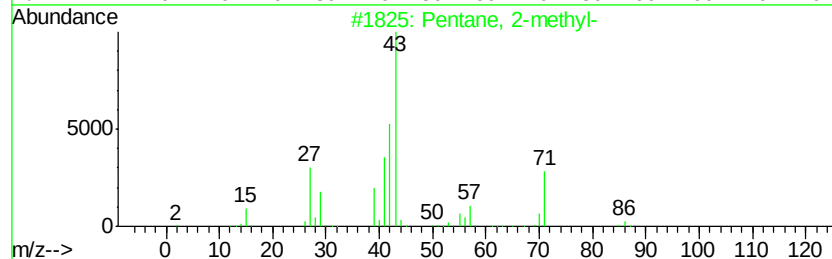
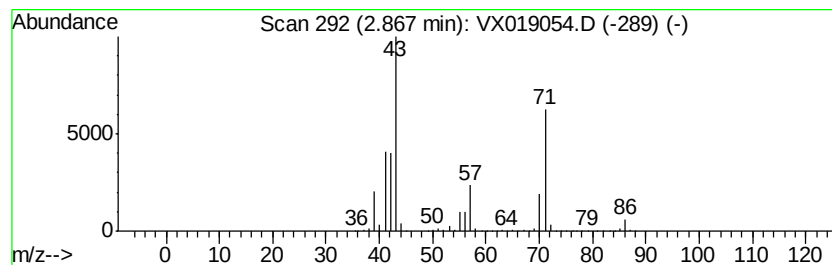
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X102120W.M
Quant Title : SW846 8260

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

Peak Number 1 Pentane, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.87	129.83 ug/l	5848140	Pentafluorobenzene	5.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2-methyl-	86	C6H14	000107-83-5	72
2		Pentane, 3,3-dimethyl-	100	C7H16	000562-49-2	47
3		Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	47
4		Sulfurous acid, hexyl 2-pentyl e...	236	C11H24O3S	1000309-15-6	42
5		Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	38



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ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
MW-10

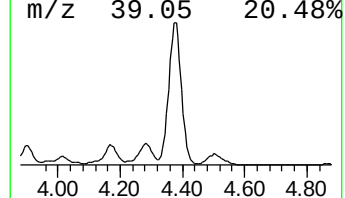
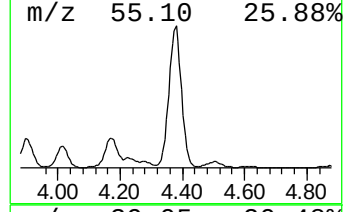
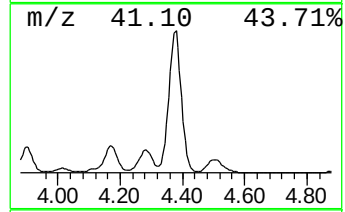
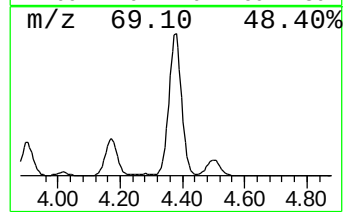
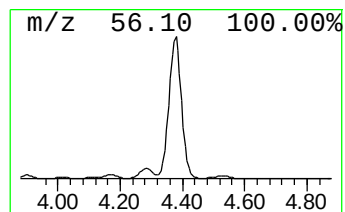
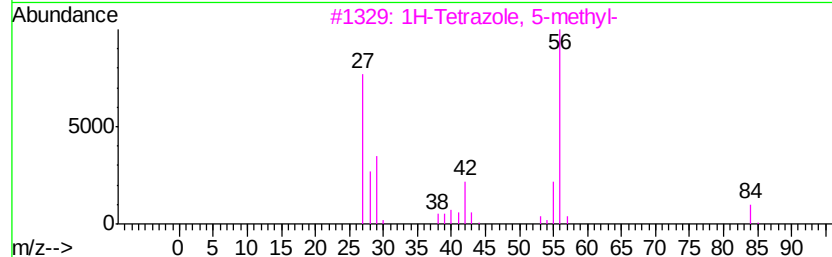
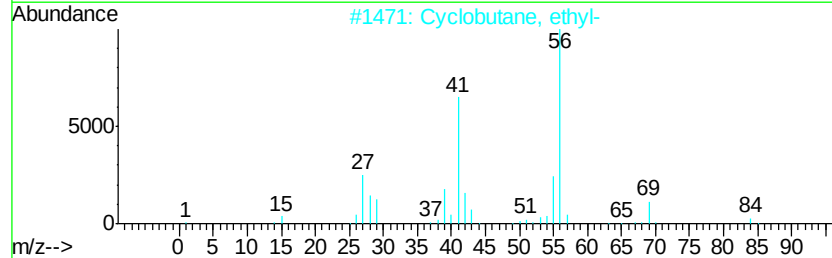
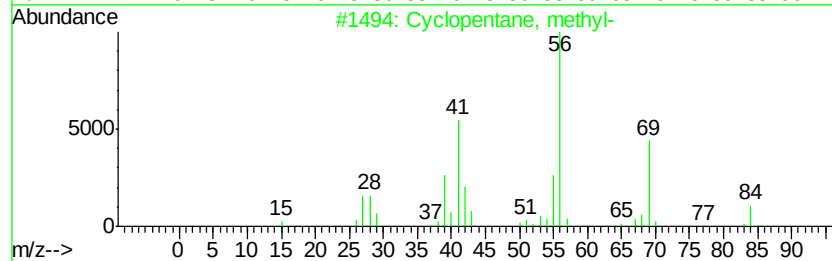
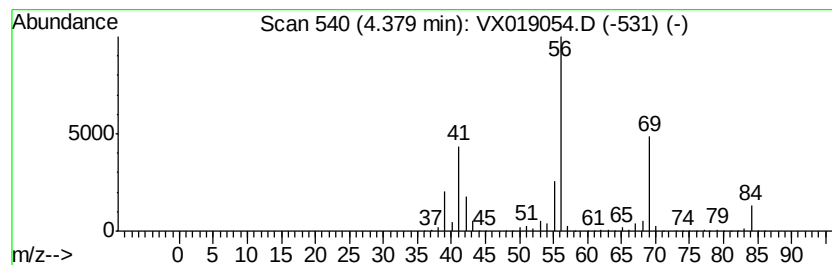
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X102120W.M
Quant Title : SW846 8260

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

Peak Number 2 Cyclopentane, methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.38	117.32 ug/l	5284510	Pentafluorobenzene	5.64

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2		Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
3		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64
4		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	52
5		1-Octene, 7-methyl-	126	C9H18	013151-06-9	43



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ALS Vial : 45 Sample Multiplier: 1

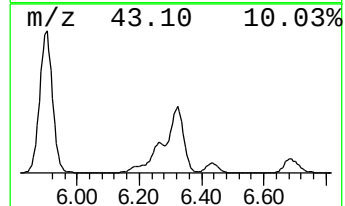
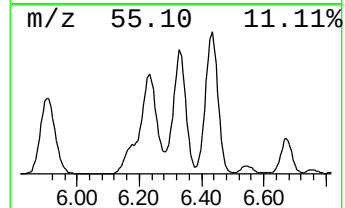
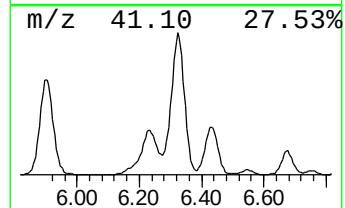
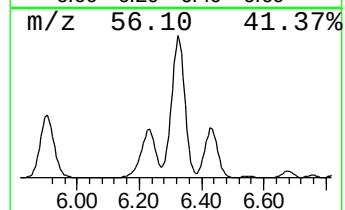
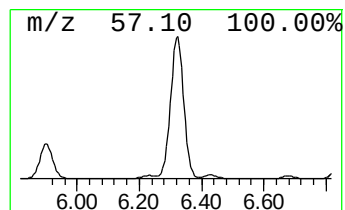
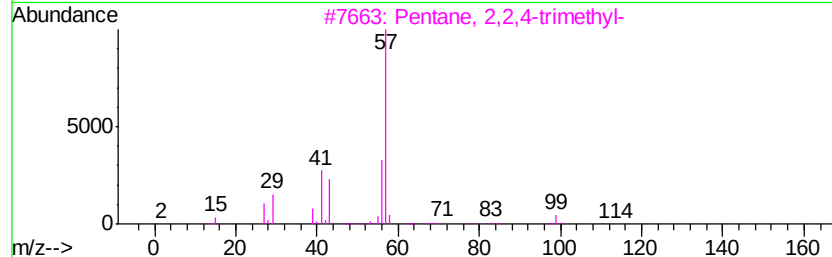
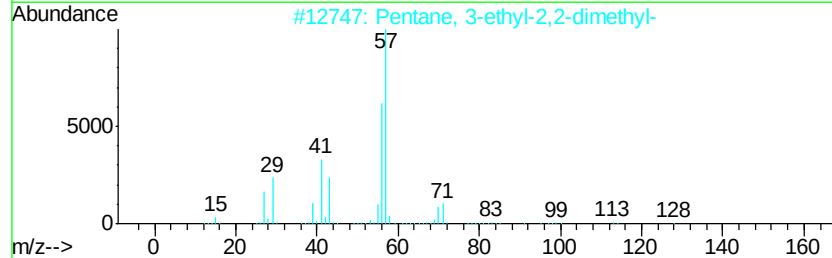
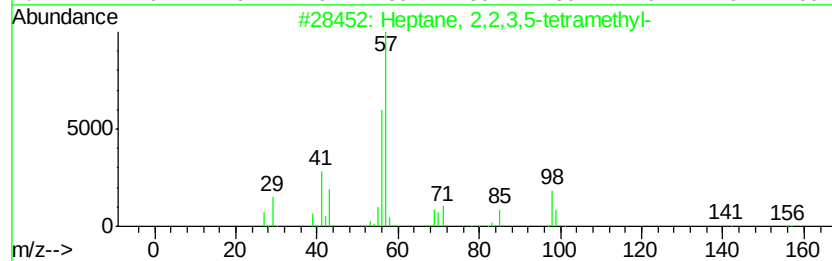
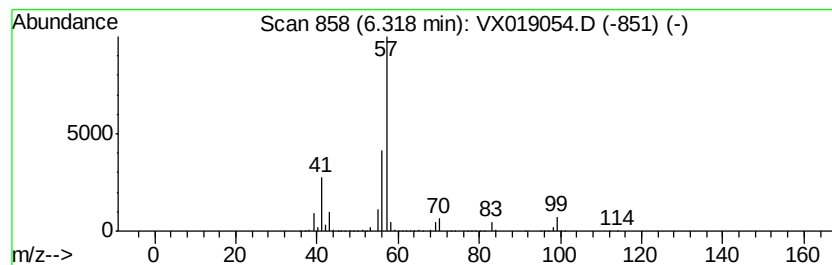
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Heptane, 2,2,3,5-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.32	123.31 ug/l	2943090	1,4-Difluorobenzene	6.83
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heptane, 2,2,3,5-tetramethyl-	156 C11H24	061868-42-6	78
2	Pentane, 3-ethyl-2,2-dimethyl-	128 C9H20	016747-32-3	64
3	Pentane, 2,2,4-trimethyl-	114 C8H18	000540-84-1	59
4	Pentane, 2,2,3,4-tetramethyl-	128 C9H20	001186-53-4	56
5	Hexane, 2,2-dimethyl-	114 C8H18	000590-73-8	53



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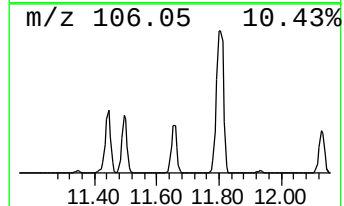
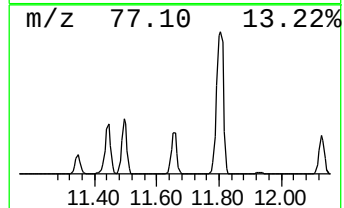
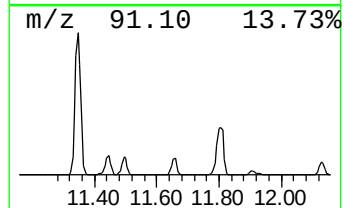
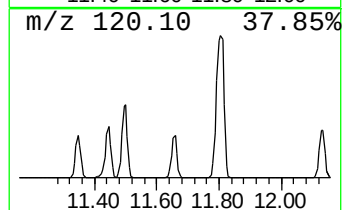
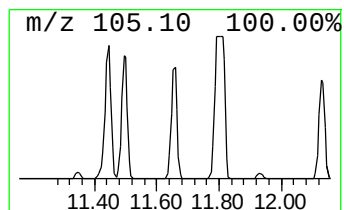
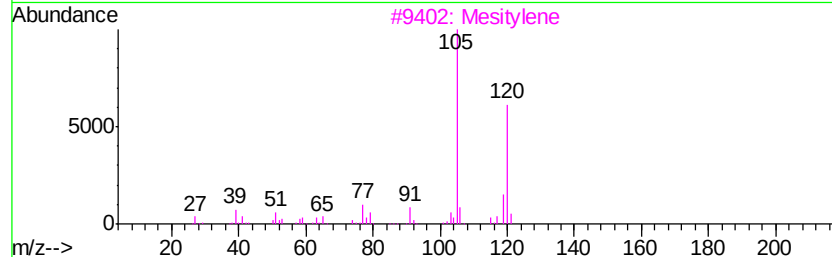
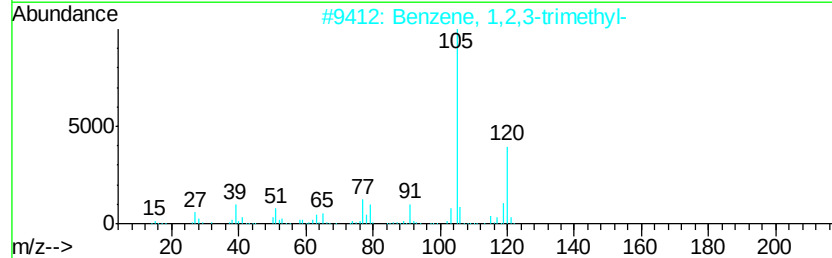
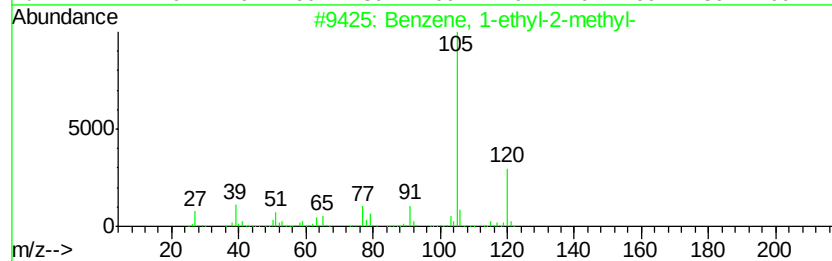
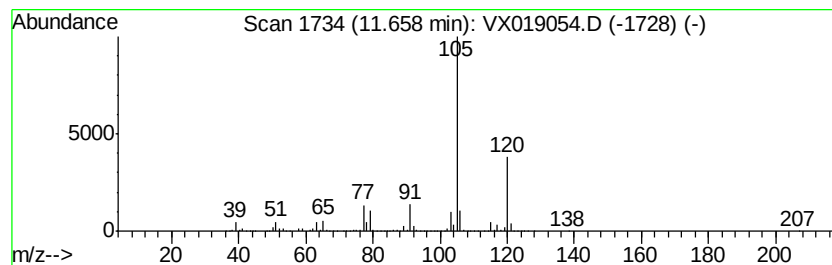
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 1-ethyl-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.66	603.09 ug/l	22825400	1,4-Dichlorobenzene-d4	12.06

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	93
3		Mesitylene	120	C9H12	000108-67-8	91
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



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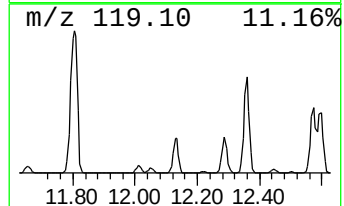
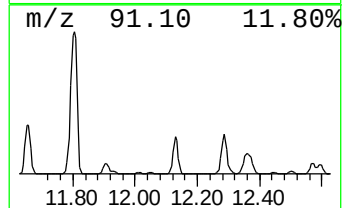
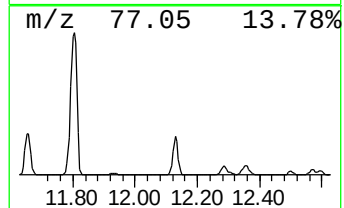
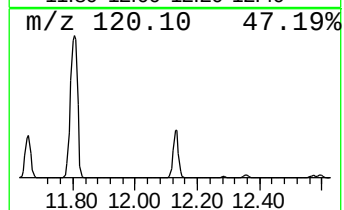
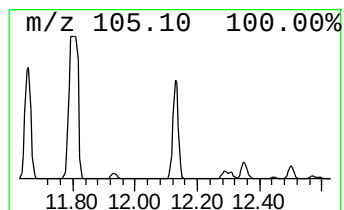
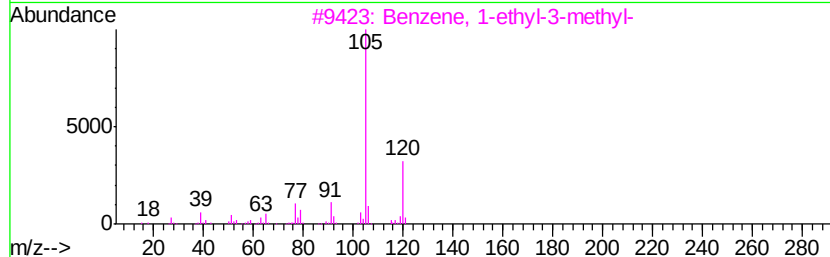
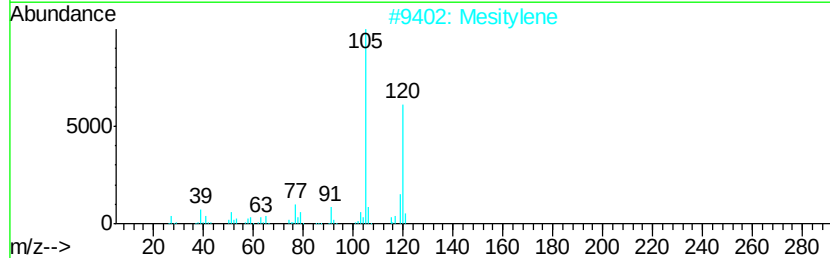
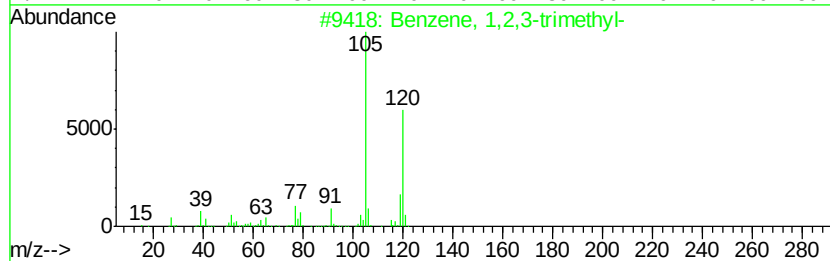
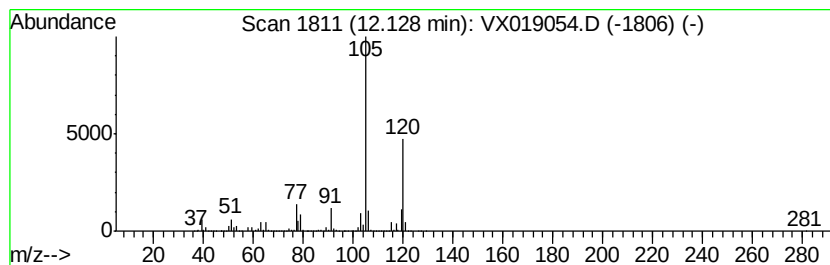
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.13	560.74 ug/l	21222600	1,4-Dichlorobenzene-d4	12.06

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
2		Mesitylene	120	C9H12	000108-67-8	94
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX102220\
Data File : VX019054.D
Acq On : 23 Oct 2020 02:27
Operator : JC/SP
Sample : L4417-17
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 45 Sample Multiplier: 1

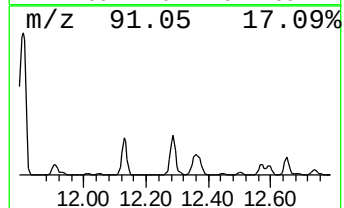
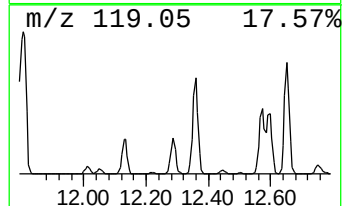
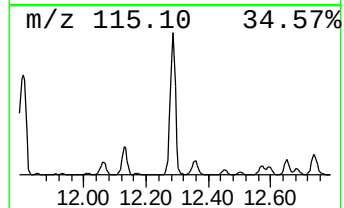
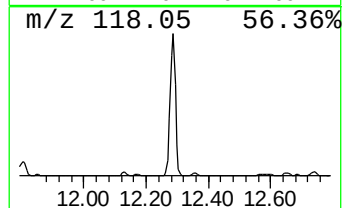
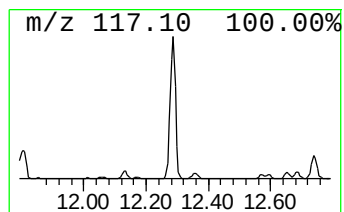
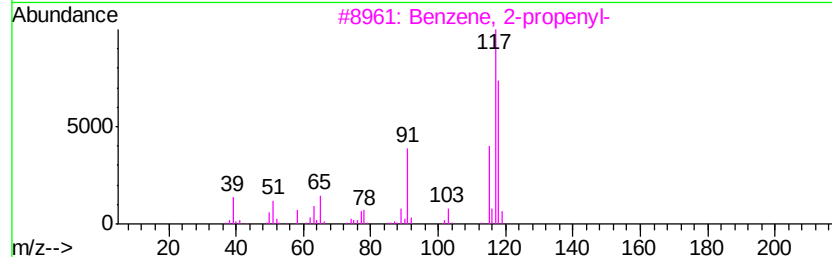
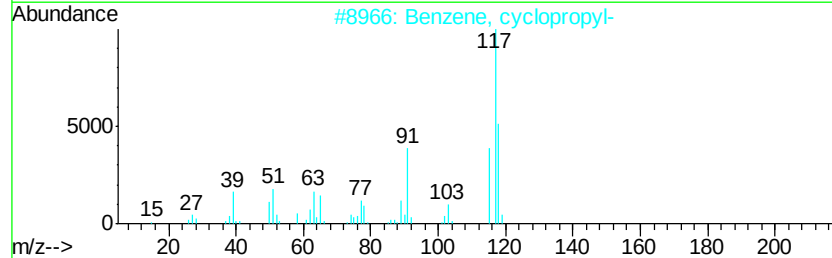
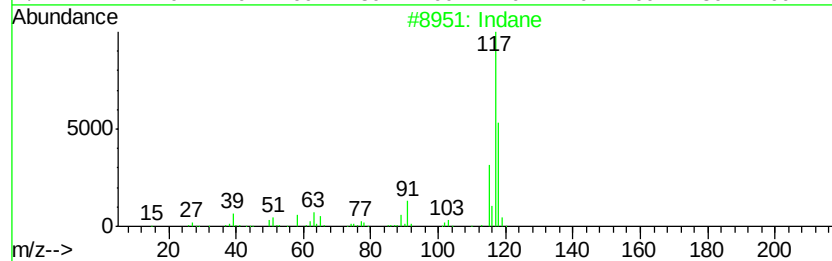
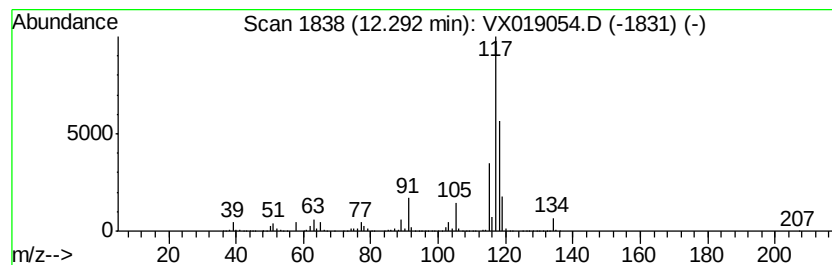
Instrument :
MSVOA_X
ClientSampleId :
MW-10

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X102120W.M
Quant Title : SW846 8260

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

Peak Number 6 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.29	455.65 ug/l	17245000	1,4-Dichlorobenzene-d4	12.06	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	76
2	Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	68
4	Deltacyclene	118	C9H10	007785-10-6	58
5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	49



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX102220\
Data File : VX019054.D
Acq On : 23 Oct 2020 02:27
Operator : JC/SP
Sample : L4417-17
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
MW-10

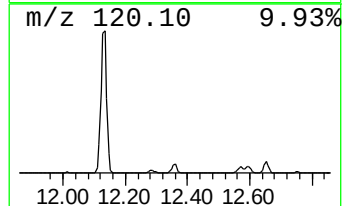
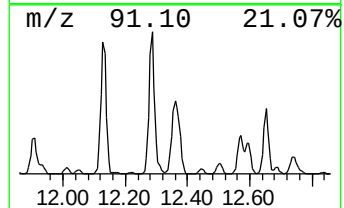
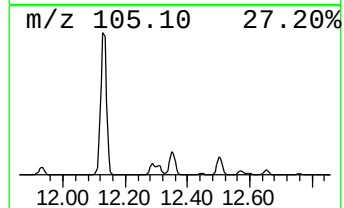
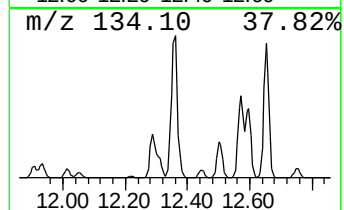
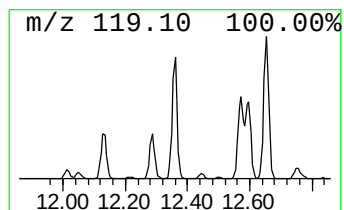
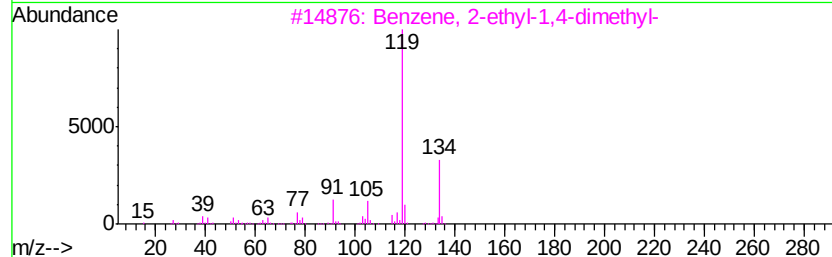
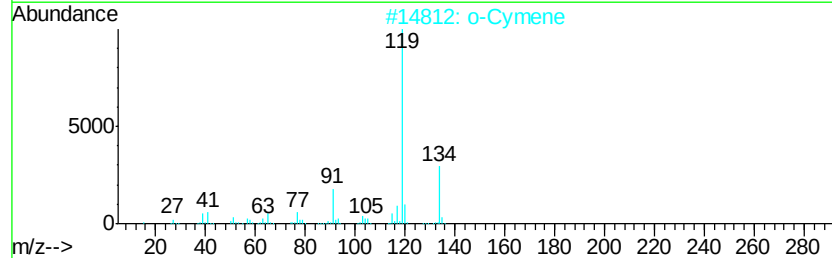
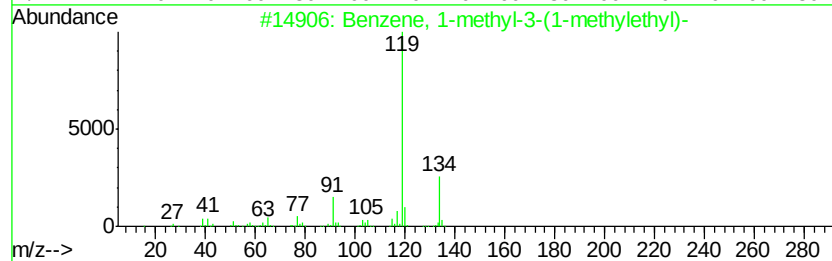
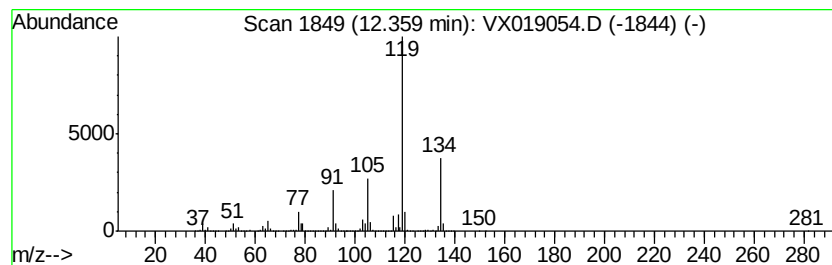
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-methyl-3-(1-meth... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.36	218.31 ug/l	8262600	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
2		o-Cymene	134	C10H14	000527-84-4	95
3		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	93
5		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX102220\
Data File : VX019054.D
Acq On : 23 Oct 2020 02:27
Operator : JC/SP
Sample : L4417-17
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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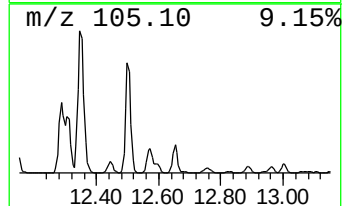
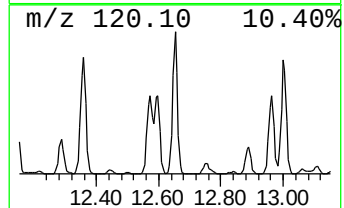
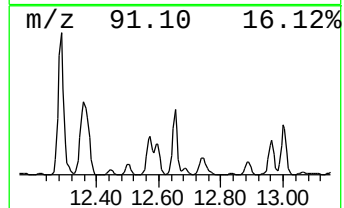
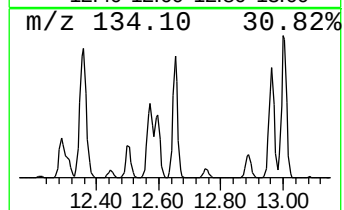
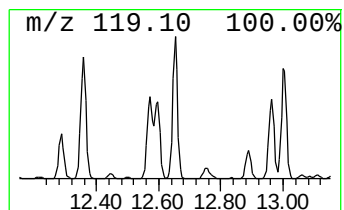
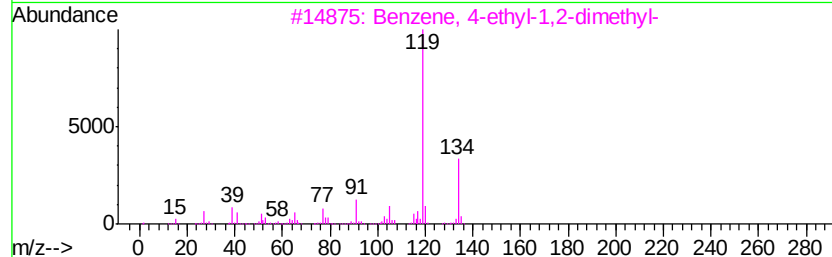
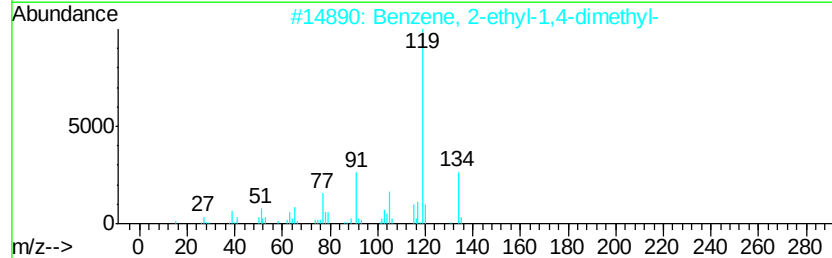
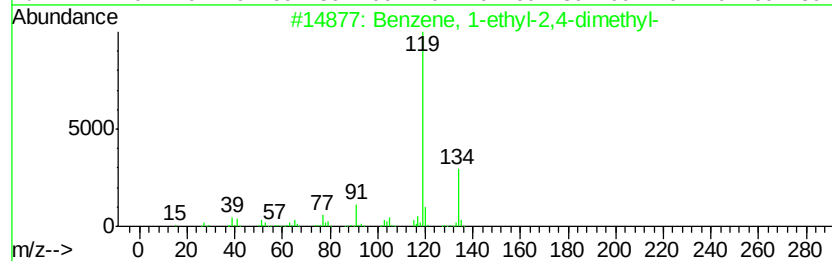
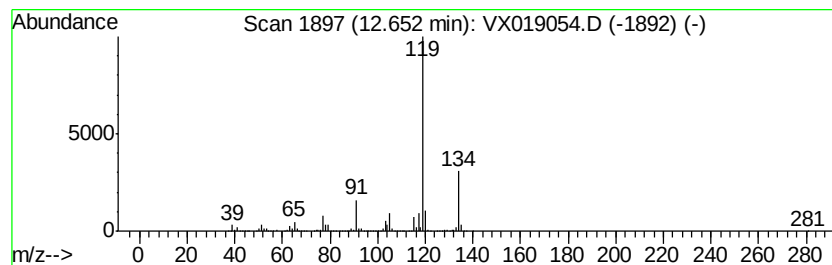
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	164.02 ug/l	6207550	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
4		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
5		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX102220\
Data File : VX019054.D
Acq On : 23 Oct 2020 02:27
Operator : JC/SP
Sample : L4417-17
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10

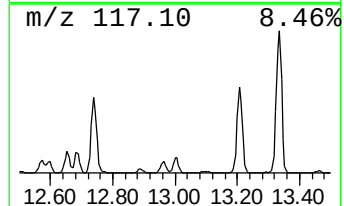
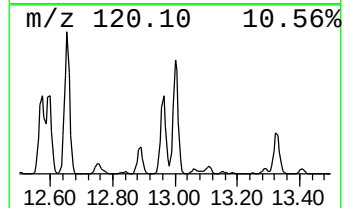
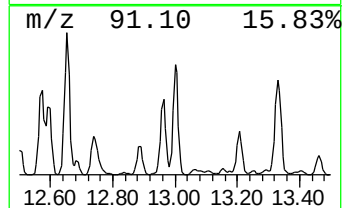
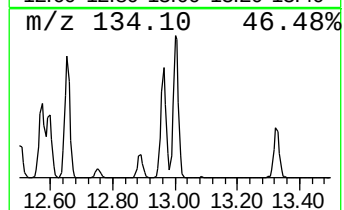
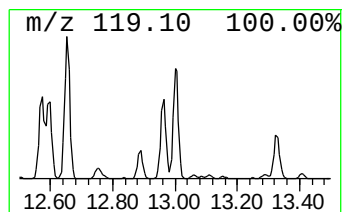
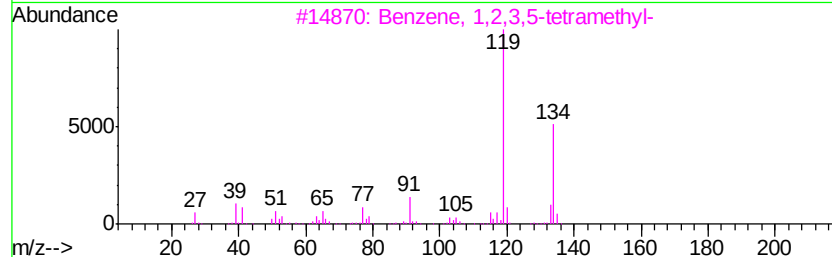
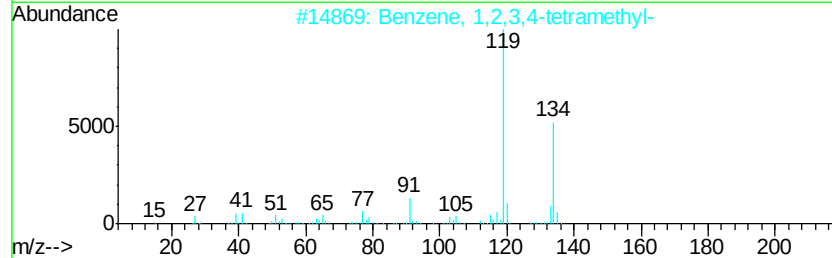
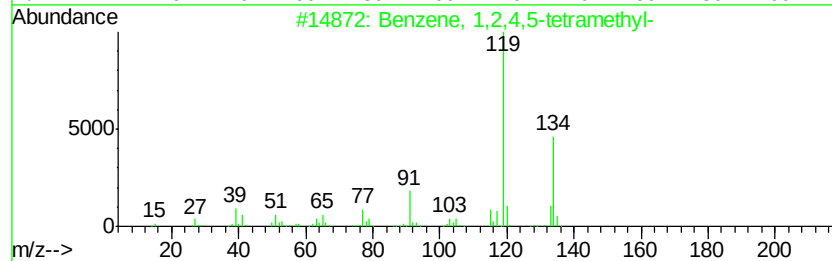
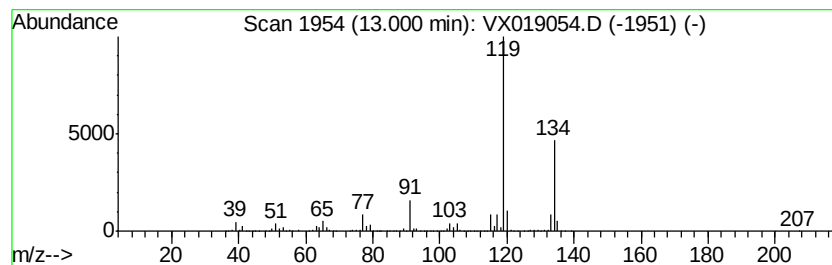
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X102120W.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.00	145.19 ug/l	5495120	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
5		o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX102220\
Data File : VX019054.D
Acq On : 23 Oct 2020 02:27
Operator : JC/SP
Sample : L4417-17
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
MW-10

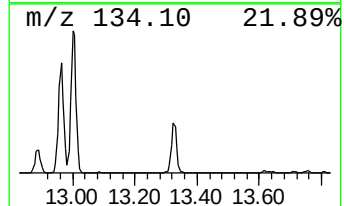
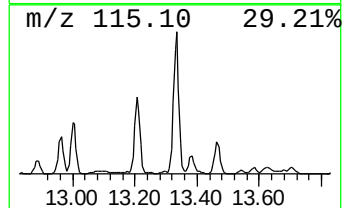
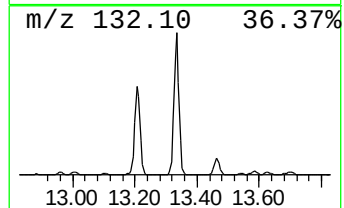
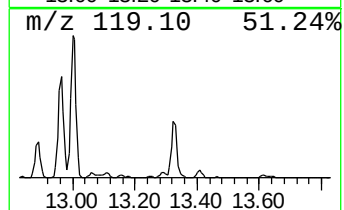
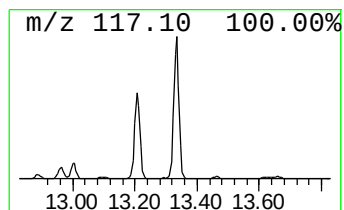
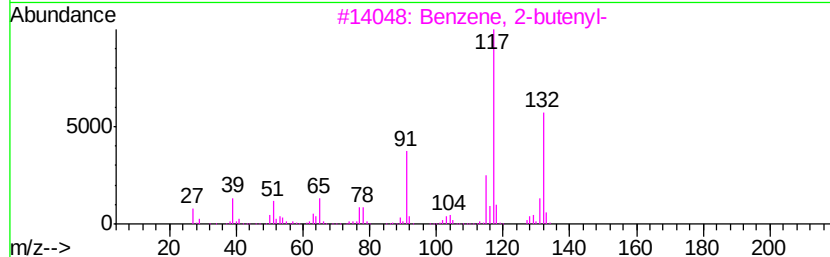
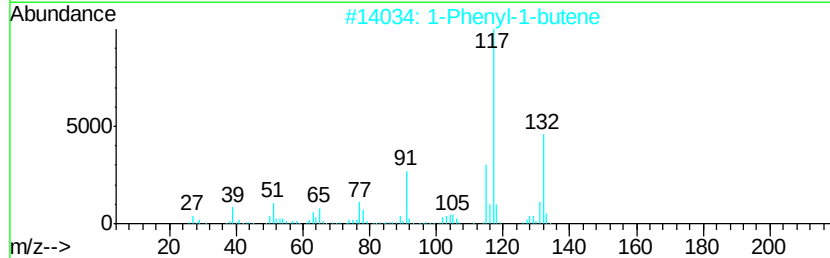
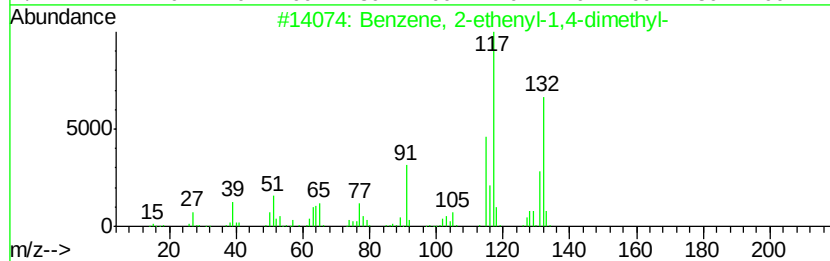
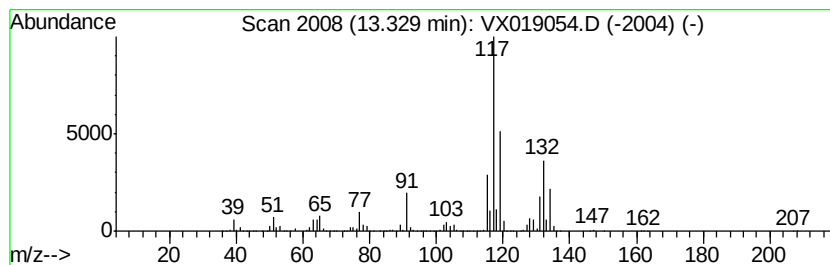
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X102120W.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.33	172.17 ug/l	6516130	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2		1-Phenyl-1-butene	132	C10H12	000824-90-8	70
3		Benzene, 2-butenyl-	132	C10H12	001560-06-1	70
4		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	70
5		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	64



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX102220\
Data File : VX019054.D
Acq On : 23 Oct 2020 02:27
Operator : JC/SP
Sample : L4417-17
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X102120W.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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Cyclopentane, met...	4.38	117.3	ug/l	5284510	1	5.64	2252230	50.0
Heptane, 2,2,3,5-...	6.32	123.3	ug/l	2943090	2	6.83	1193380	50.0
Benzene, 1-ethyl-...	11.66	603.1	ug/l	22825400	4	12.06	1892360	50.0
Benzene, 1,2,3-tr...	12.13	560.7	ug/l	21222600	4	12.06	1892360	50.0
Indane	12.29	455.6	ug/l	17245000	4	12.06	1892360	50.0
Benzene, 1-methyl...	12.36	218.3	ug/l	8262600	4	12.06	1892360	50.0
Benzene, 1-ethyl-...	12.65	164.0	ug/l	6207550	4	12.06	1892360	50.0
Benzene, 1,2,4,5-...	13.00	145.2	ug/l	5495120	4	12.06	1892360	50.0
Benzene, 2-etheny...	13.33	172.2	ug/l	6516130	4	12.06	1892360	50.0