

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX060618\
 Data File : VX002360.D
 Acq On : 06 Jun 2018 18:36
 Operator : JC/MD
 Sample : J3312-03
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 W-16

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\82X060418W.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	1.075	76	80	94	rBV	1096210	1379497	35.87%	3.526%
2	1.270	108	112	115	rBV	36657	48617	1.26%	0.124%
3	1.337	118	123	128	rVV2	45369	138444	3.60%	0.354%
4	1.404	131	134	140	rVV	59663	97812	2.54%	0.250%
5	1.459	140	143	152	rVB	86358	111931	2.91%	0.286%
6	2.270	270	276	283	rVB	163617	249051	6.48%	0.637%
7	2.812	357	365	377	rBV	615840	1166913	30.34%	2.983%
8	2.934	379	385	397	rVB	33562	73297	1.91%	0.187%
9	5.513	795	808	815	rBV	152377	436748	11.36%	1.116%
10	5.665	825	833	852	rVB	236788	668336	17.38%	1.708%
11	6.074	890	900	911	rBV	173561	452744	11.77%	1.157%
12	6.306	929	938	946	rBV3	18913	49798	1.29%	0.127%
13	6.866	1016	1030	1053	rBV	349933	827484	21.52%	2.115%
14	7.464	1115	1128	1139	rBV	43209	99656	2.59%	0.255%
15	8.720	1328	1334	1341	rVV	846278	1318876	34.29%	3.371%
16	8.787	1341	1345	1350	rVB	88816	135388	3.52%	0.346%
17	9.037	1380	1386	1393	rVB4	20959	40010	1.04%	0.102%
18	10.116	1555	1563	1571	rBV	716607	1016800	26.44%	2.599%
19	10.341	1594	1600	1609	rBV4	24538	51476	1.34%	0.132%
20	10.658	1642	1652	1657	rBV6	19080	53962	1.40%	0.138%
21	10.841	1669	1682	1689	rBV3	34911	71574	1.86%	0.183%
22	10.914	1689	1694	1701	rBV4	23850	40408	1.05%	0.103%
23	11.024	1706	1712	1721	rBV	707875	935174	24.32%	2.391%
24	11.140	1726	1731	1736	rBV	750793	935469	24.32%	2.391%
25	11.366	1762	1768	1774	rBV	527361	663362	17.25%	1.696%
26	11.677	1807	1819	1826	rBV3	35219	68472	1.78%	0.175%
27	11.774	1826	1835	1839	rBV	117944	154870	4.03%	0.396%
28	11.926	1854	1860	1861	rBV	161634	205010	5.33%	0.524%
29	11.951	1861	1864	1870	rVB	364338	436792	11.36%	1.117%
30	12.085	1880	1886	1891	rBV	844966	1089603	28.33%	2.785%
31	12.237	1906	1911	1916	rVB	123962	152348	3.96%	0.389%
32	12.304	1916	1922	1928	rBV	1180433	1456590	37.88%	3.723%
33	12.372	1928	1933	1941	rVV2	229045	460204	11.97%	1.176%
34	12.463	1943	1948	1956	rBV	359746	437555	11.38%	1.119%

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Integration Parameters: RTEINT.P

Integrator: RTE
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 Sampling : 1
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 Max Peaks: 100
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 Title : SW846 8260

35	12.670	1974	1982	1985	rBV	1509309	1882106	48.94%	4.811%
36	12.707	1985	1988	1992	rVV	581331	730947	19.01%	1.869%
37	12.762	1992	1997	2004	rVV2	1725772	2314934	60.20%	5.918%
38	12.823	2004	2007	2012	rVV2	61966	83580	2.17%	0.214%
39	12.902	2012	2020	2025	rVB	294602	434323	11.29%	1.110%
40	12.981	2025	2033	2038	rBV	1362258	1750716	45.52%	4.475%
41	13.030	2038	2041	2046	rVB	63284	77466	2.01%	0.198%
42	13.091	2046	2051	2057	rVB3	143085	218947	5.69%	0.560%
43	13.158	2057	2062	2065	rBV	169583	234616	6.10%	0.600%
44	13.201	2065	2069	2071	rVV2	217529	279861	7.28%	0.715%
45	13.231	2071	2074	2083	rVB2	455319	646814	16.82%	1.653%
46	13.316	2083	2088	2089	rBV2	107444	126508	3.29%	0.323%
47	13.353	2089	2094	2100	rVV2	2676687	3845716	100.00%	9.831%
48	13.408	2100	2103	2105	rVV2	88661	123205	3.20%	0.315%
49	13.432	2105	2107	2109	rVV	117079	116050	3.02%	0.297%
50	13.487	2113	2116	2121	rVB2	84600	109425	2.85%	0.280%
51	13.573	2122	2130	2132	rBV2	106810	168790	4.39%	0.431%
52	13.603	2132	2135	2138	rVV	277950	362661	9.43%	0.927%
53	13.646	2138	2142	2146	rVV2	418953	597542	15.54%	1.527%
54	13.707	2146	2152	2153	rVV3	299285	534752	13.91%	1.367%
55	13.725	2153	2155	2163	rVB2	541009	739047	19.22%	1.889%
56	13.810	2163	2169	2173	rBV	130791	180279	4.69%	0.461%
57	13.859	2173	2177	2182	rVB2	146698	198853	5.17%	0.508%
58	13.914	2182	2186	2191	rVB	187359	244871	6.37%	0.626%
59	13.987	2191	2198	2202	rBV2	299127	483012	12.56%	1.235%
60	14.036	2202	2206	2210	rVV	137401	170133	4.42%	0.435%
61	14.133	2218	2222	2225	rBV2	142990	189145	4.92%	0.484%
62	14.170	2225	2228	2232	rVB	89410	100591	2.62%	0.257%
63	14.280	2241	2246	2252	rBV2	285087	493767	12.84%	1.262%
64	14.383	2259	2263	2267	rVB	122320	134461	3.50%	0.344%
65	14.438	2267	2272	2276	rVB	177039	228820	5.95%	0.585%
66	14.481	2276	2279	2284	rVB	61007	79058	2.06%	0.202%
67	14.542	2284	2289	2294	rBV2	393754	557503	14.50%	1.425%
68	14.676	2303	2311	2317	rBV4	69081	160107	4.16%	0.409%
69	14.737	2317	2321	2325	rVV	113079	147049	3.82%	0.376%
70	14.786	2326	2329	2333	rVB2	45013	54571	1.42%	0.139%
71	14.847	2333	2339	2347	rBV	2042185	2709407	70.45%	6.926%

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Integration Parameters: RTEINT.P

Integrator: RTE
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 Sampling : 1 Min Area: 3 % of largest Peak
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72	14.914	2347	2350	2353	rVB4	40126	52673	1.37%	0.135%
73	14.981	2353	2361	2367	rVB6	26197	65239	1.70%	0.167%
74	15.255	2401	2406	2413	rVB3	75657	133489	3.47%	0.341%
75	15.328	2413	2418	2425	rVB	33539	52086	1.35%	0.133%
76	15.450	2431	2438	2440	rBV2	47710	76008	1.98%	0.194%
77	15.487	2440	2444	2448	rVV	90994	134470	3.50%	0.344%
78	15.554	2449	2455	2460	rVV	194388	318896	8.29%	0.815%
79	15.603	2460	2463	2468	rVV5	27852	47965	1.25%	0.123%
80	15.694	2469	2478	2482	rVV	372615	631596	16.42%	1.615%
81	15.731	2482	2484	2492	rVB	99218	138658	3.61%	0.354%
82	15.816	2492	2498	2504	rBV2	43795	85985	2.24%	0.220%
83	15.926	2513	2516	2526	rVB	110471	207957	5.41%	0.532%
84	16.078	2535	2541	2550	rBV	81597	140217	3.65%	0.358%

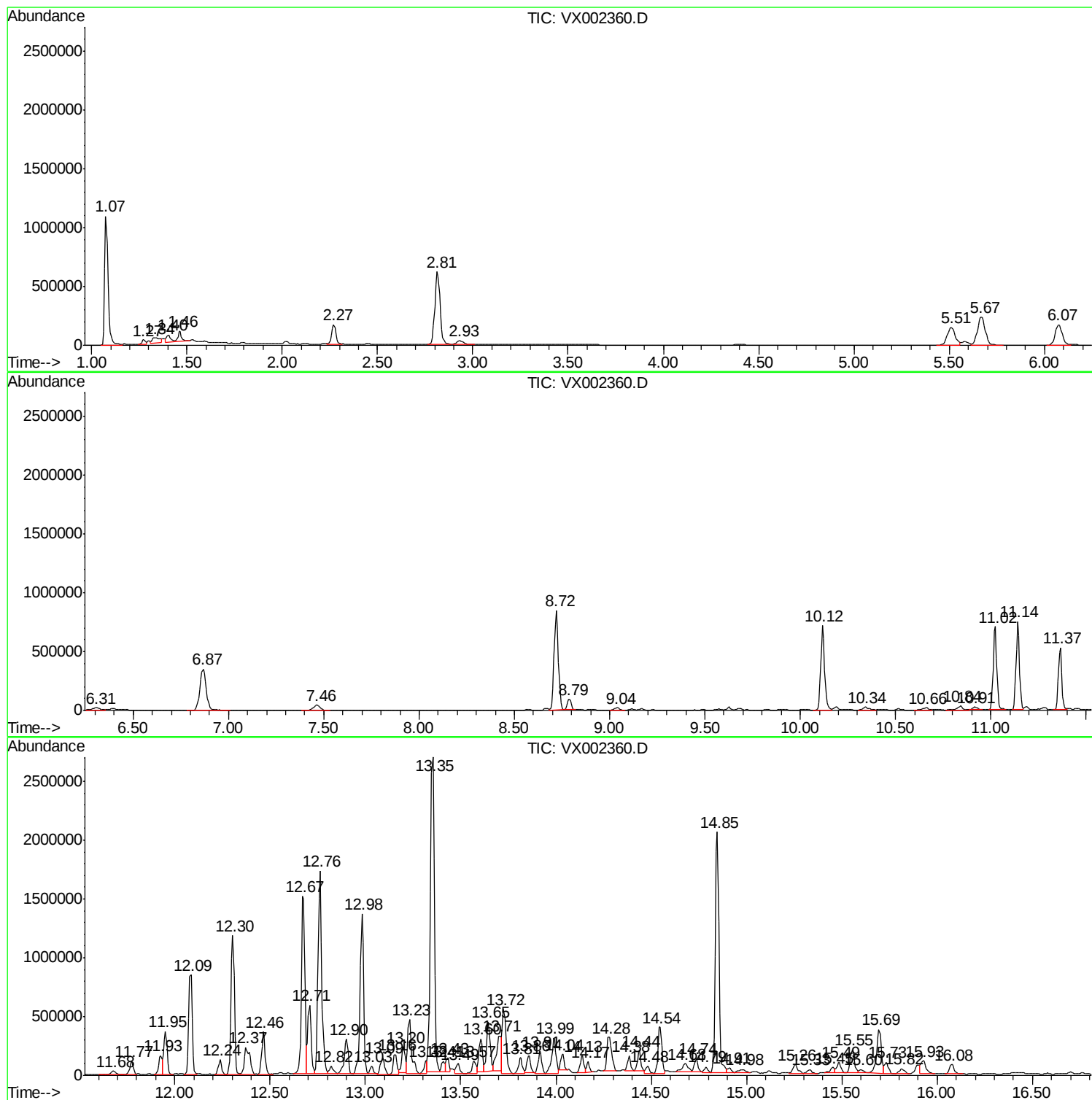
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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



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Instrument :
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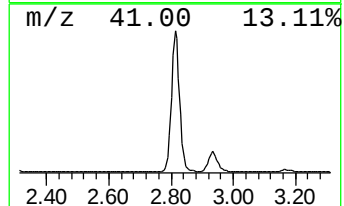
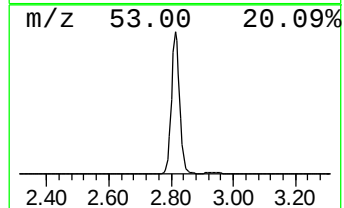
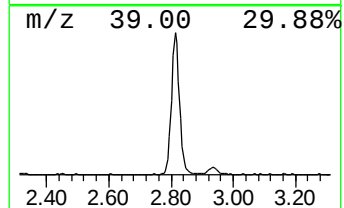
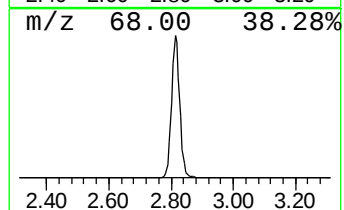
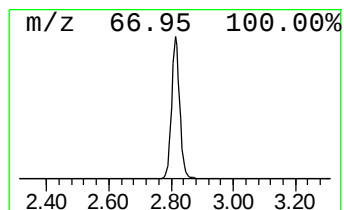
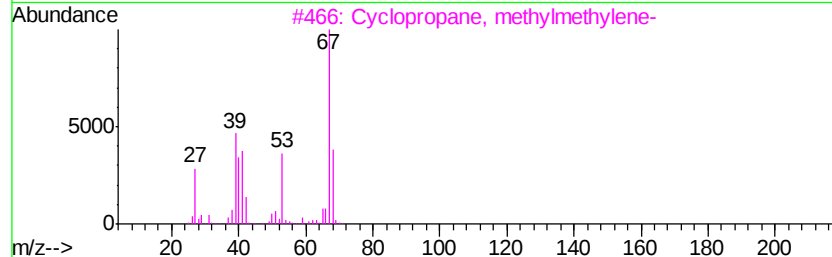
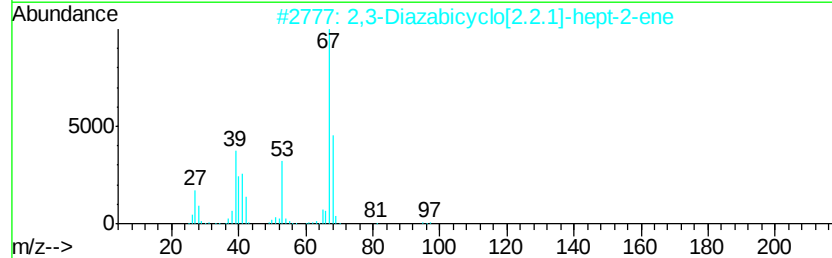
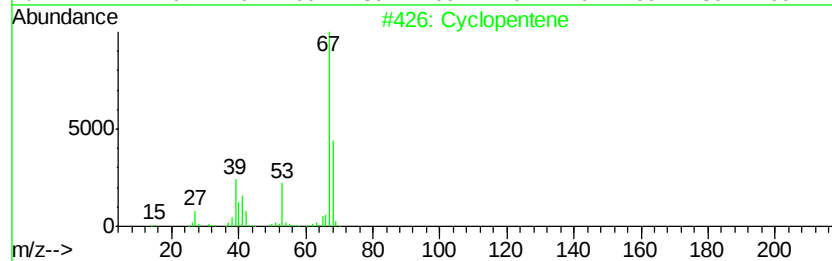
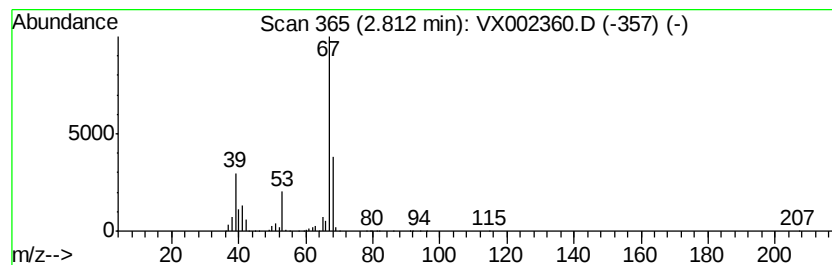
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

Peak Number 2 Cyclopentene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.81	87.30 ug/l	1166910	Pentafluorobenzene	5.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentene	68	C5H8	000142-29-0	91
2		2,3-Diazabicyclo[2.2.1]-hept-2-ene	96	C5H8N2	002721-32-6	78
3		Cyclopropane, methvlmethvlene-	68	C5H8	018631-84-0	59
4		1,4-Pentadiene	68	C5H8	000591-93-5	58
5		1,3-Pentadiene	68	C5H8	000504-60-9	53



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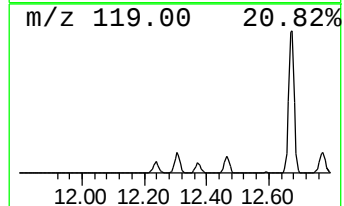
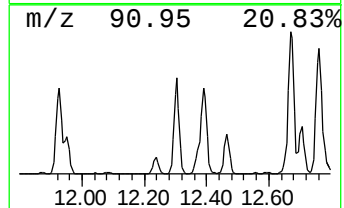
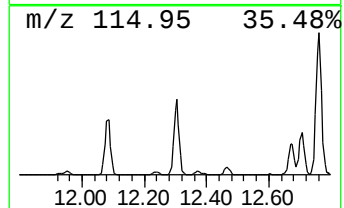
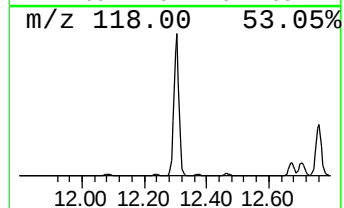
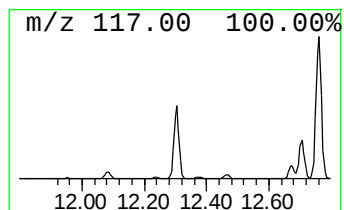
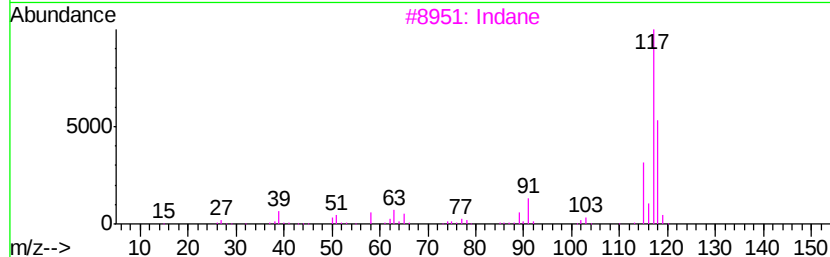
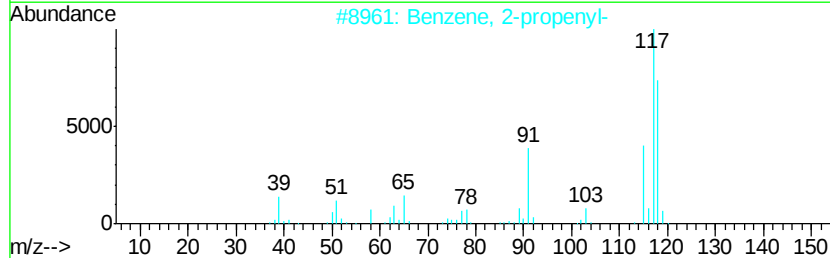
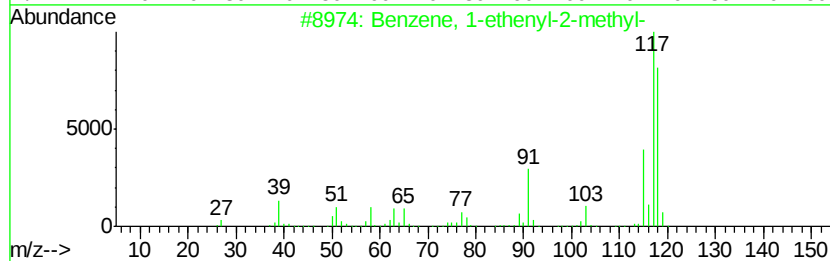
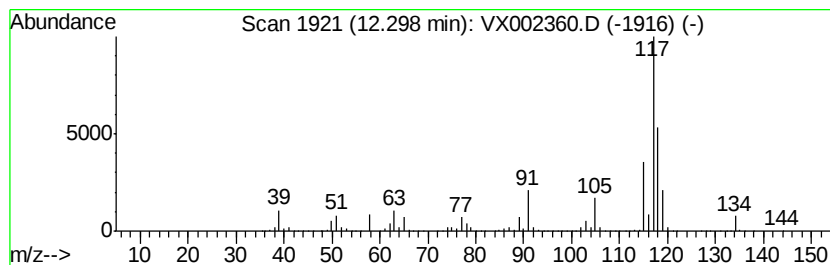
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Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1-ethenyl-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.30	66.84 ug/l	1456590	1,4-Dichlorobenzene-d4	12.09		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	86
2		Benzene, 2-propenyl-	118	C9H10	000300-57-2	70
3		Indane	118	C9H10	000496-11-7	64
4		Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
5		Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	50



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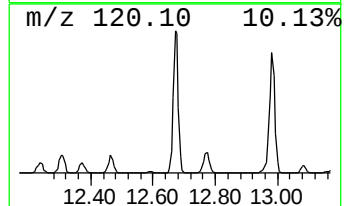
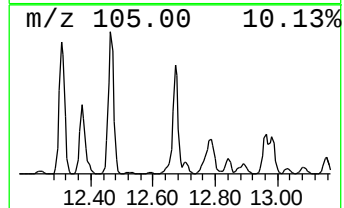
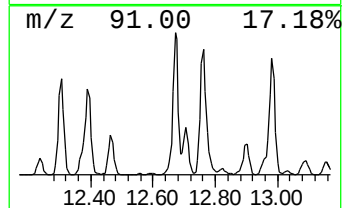
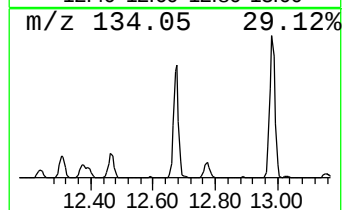
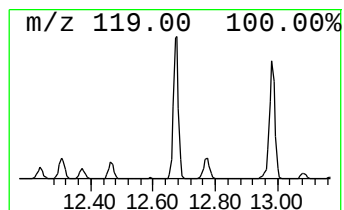
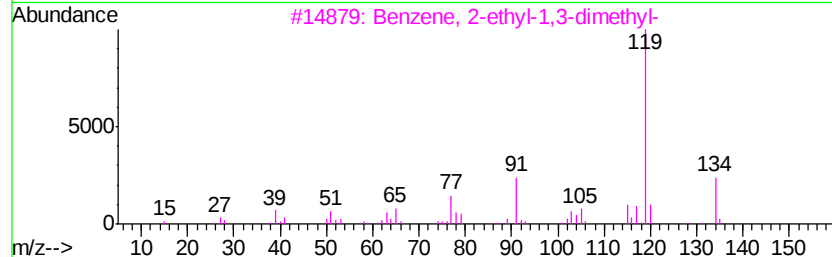
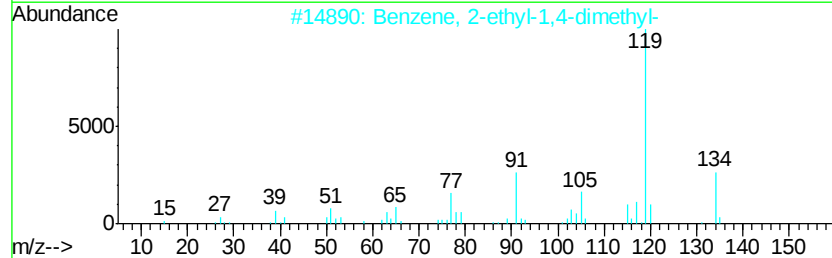
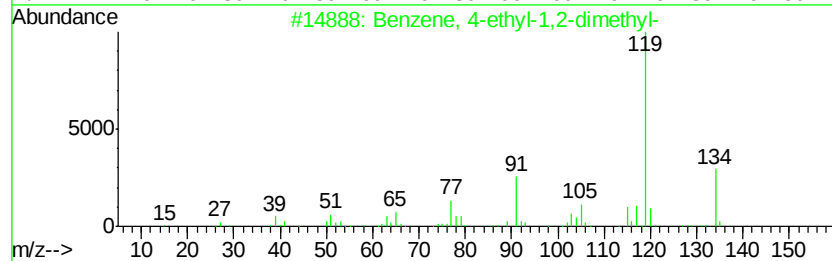
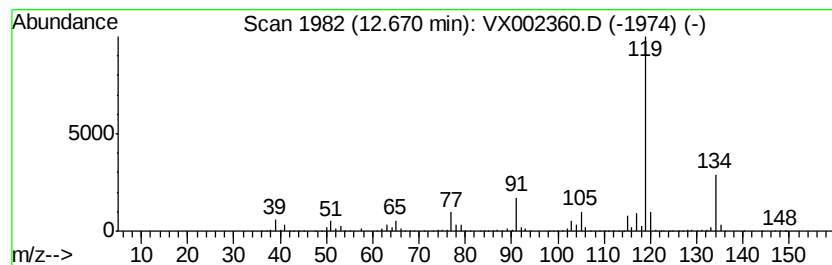
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	86.37 ug/l	1882110	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	97
2		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4		o-Cymene	134	C10H14	000527-84-4	95
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95



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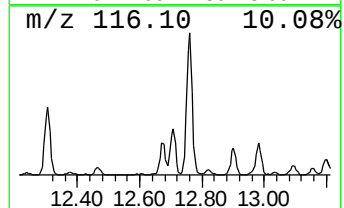
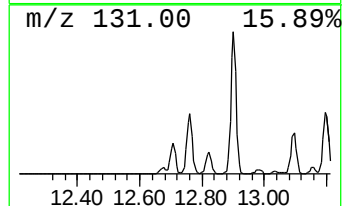
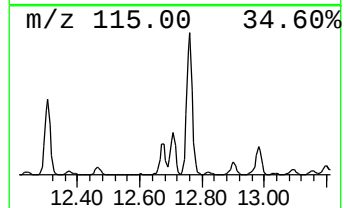
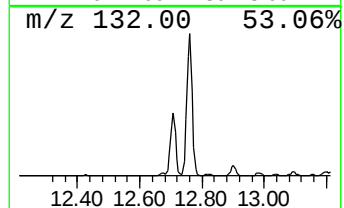
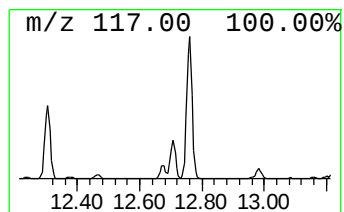
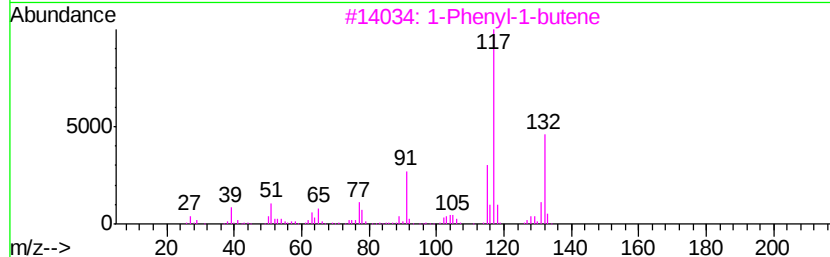
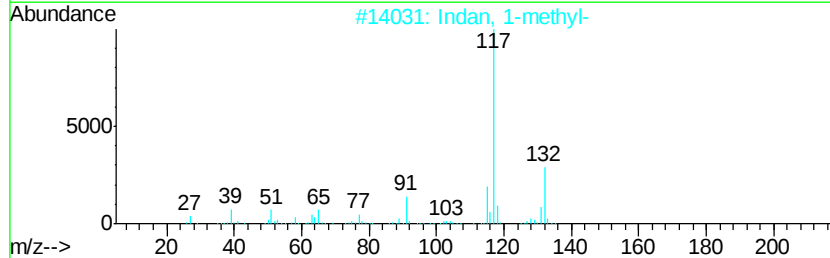
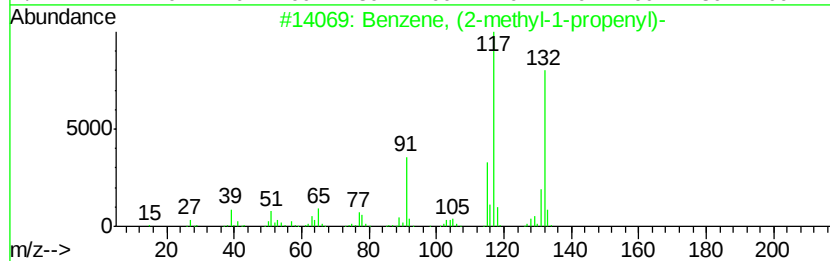
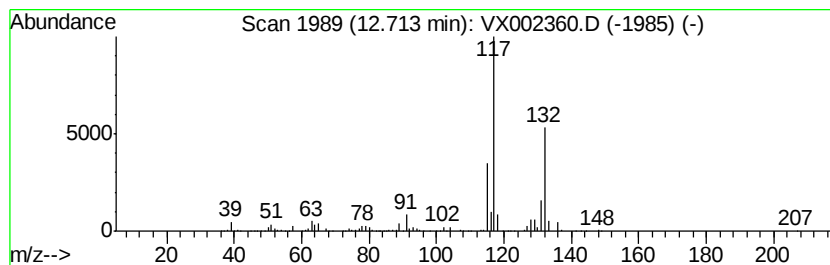
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, (2-methyl-1-propen... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.71	33.54 ug/l	730947	1,4-Dichlorobenzene-d4	12.09

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	91
2		Indan, 1-methyl-	132	C10H12	000767-58-8	90
3		1-Phenyl-1-butene	132	C10H12	000824-90-8	90
4		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	90
5		1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX060618\
Data File : VX002360.D
Acq On : 06 Jun 2018 18:36
Operator : JC/MD
Sample : J3312-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
W-16

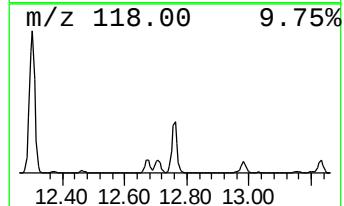
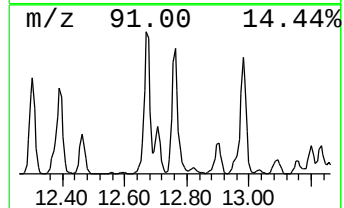
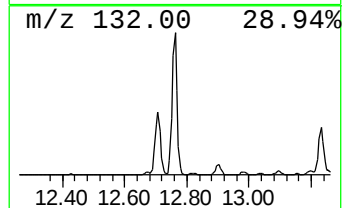
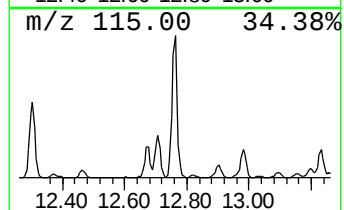
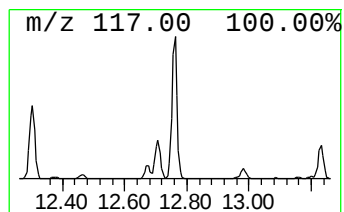
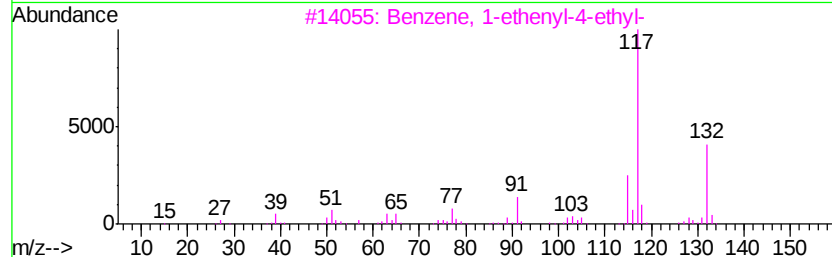
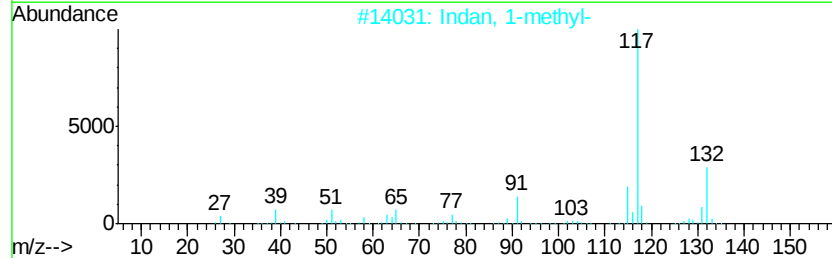
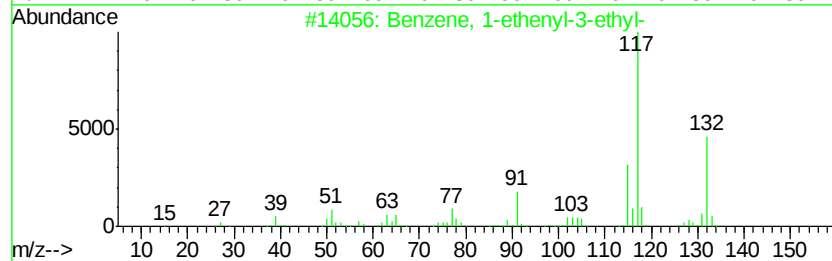
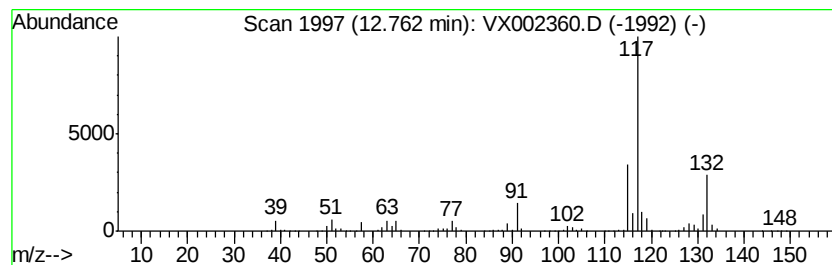
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	106.23 ug/l	2314930	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
2		Indan, 1-methyl-	132	C10H12	000767-58-8	87
3		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	86
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	83
5		Benzene, (1-methyl-1-propenyl)-, ...	132	C10H12	000768-00-3	83



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX060618\
Data File : VX002360.D
Acq On : 06 Jun 2018 18:36
Operator : JC/MD
Sample : J3312-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
W-16

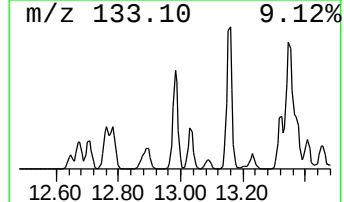
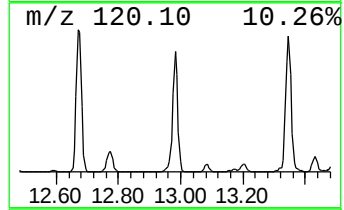
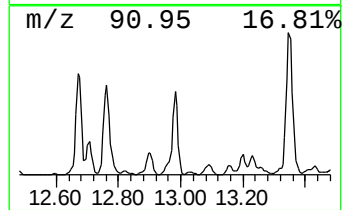
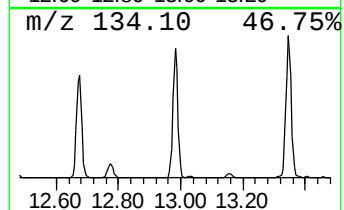
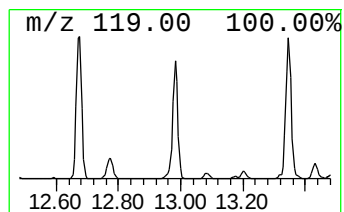
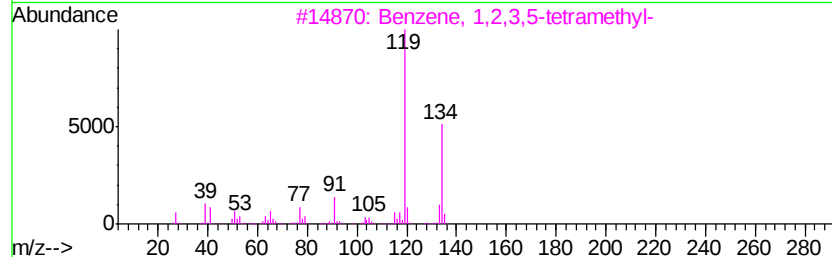
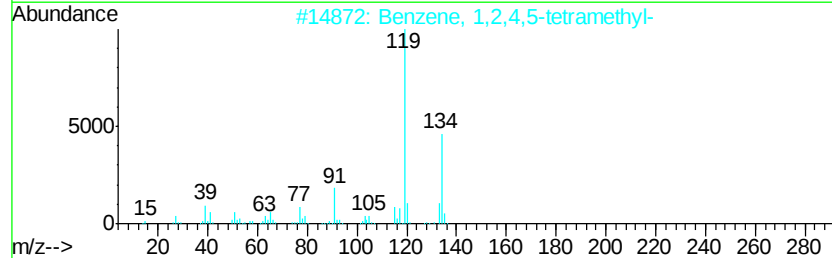
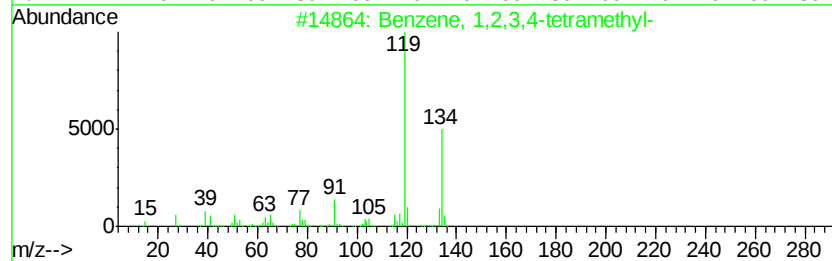
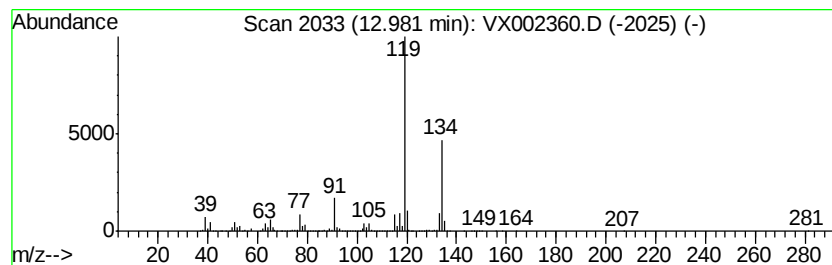
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.98	80.34 ug/l	1750720	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4		o-Cymene	134	C10H14	000527-84-4	95
5		1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX060618\
Data File : VX002360.D
Acq On : 06 Jun 2018 18:36
Operator : JC/MD
Sample : J3312-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
W-16

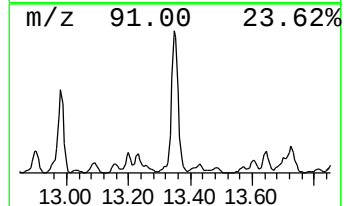
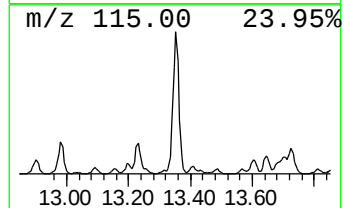
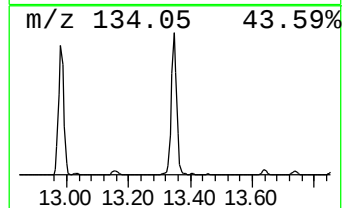
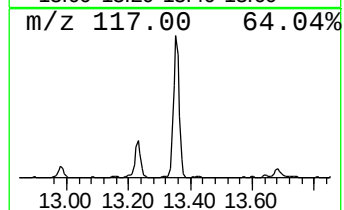
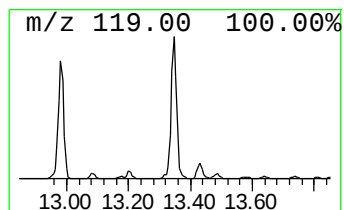
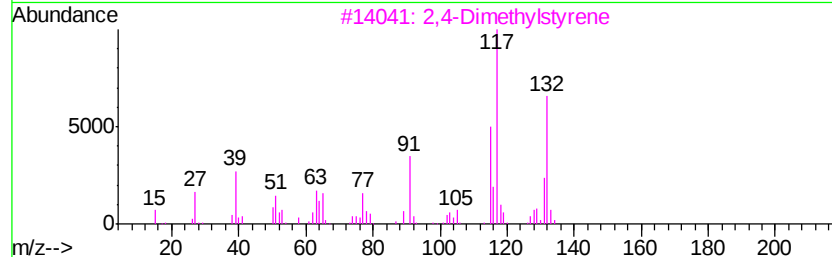
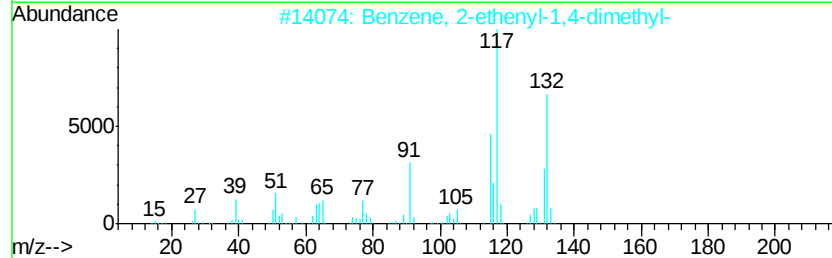
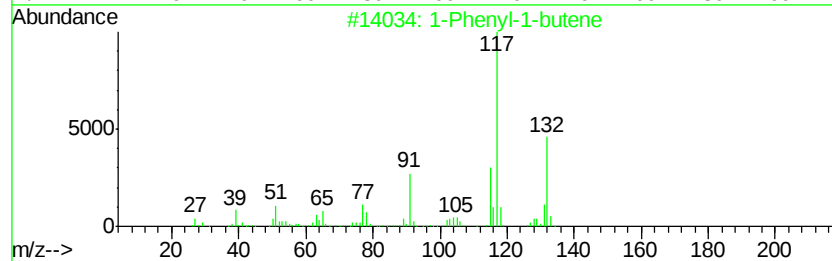
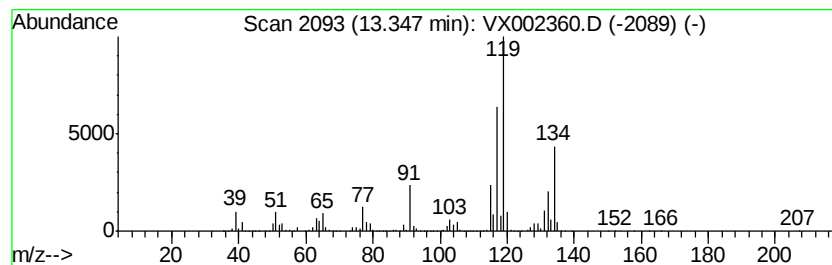
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Quant Title : SW846 8260

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

Peak Number 8 1-Phenyl-1-butene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	176.47 ug/l	3845720	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Phenyl-1-butene	132	C10H12	000824-90-8	96
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
3		2,4-Dimethylstyrene	132	C10H12	002234-20-0	95
4		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	92
5		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX060618\
Data File : VX002360.D
Acq On : 06 Jun 2018 18:36
Operator : JC/MD
Sample : J3312-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
W-16

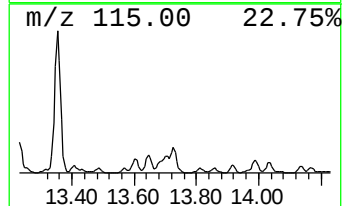
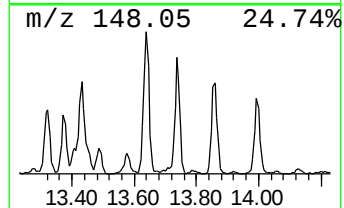
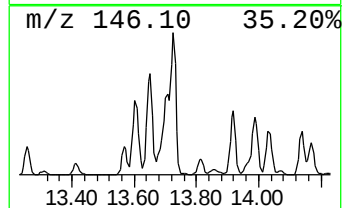
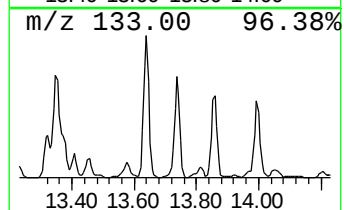
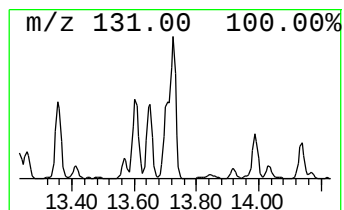
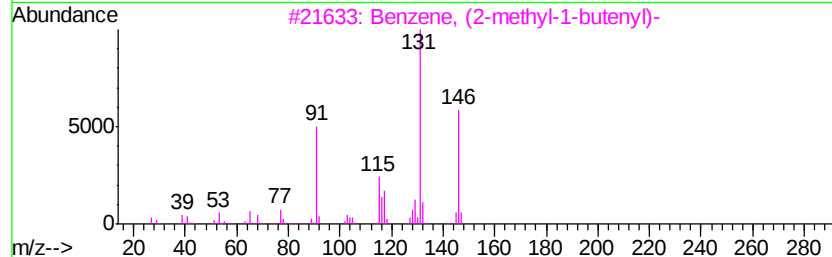
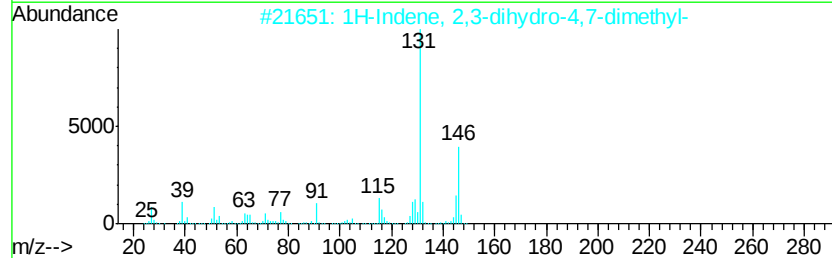
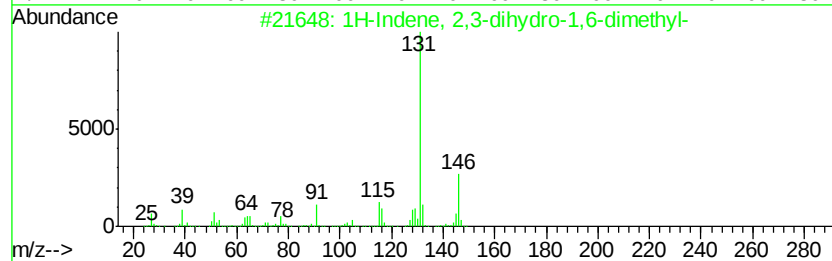
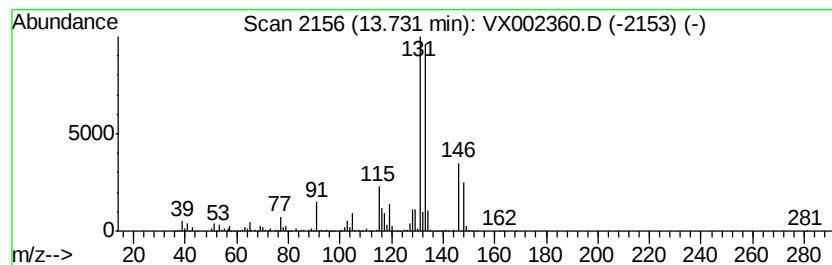
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

Peak Number 9 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.73	33.91 ug/l	739047	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	91
2		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91
3		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	91
4		1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	87
5		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX060618\
Data File : VX002360.D
Acq On : 06 Jun 2018 18:36
Operator : JC/MD
Sample : J3312-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
W-16

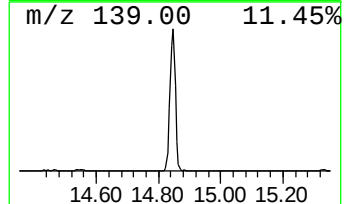
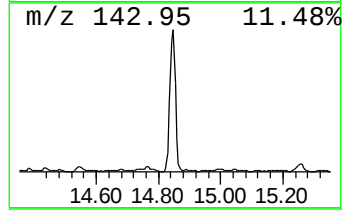
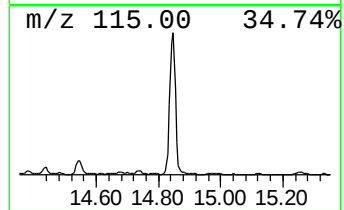
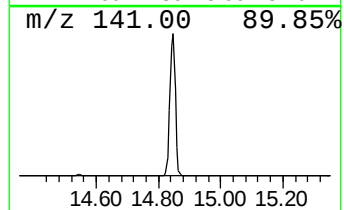
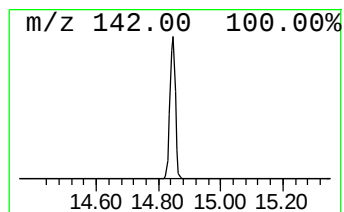
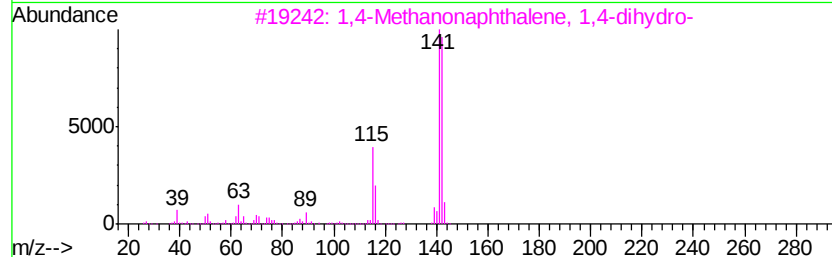
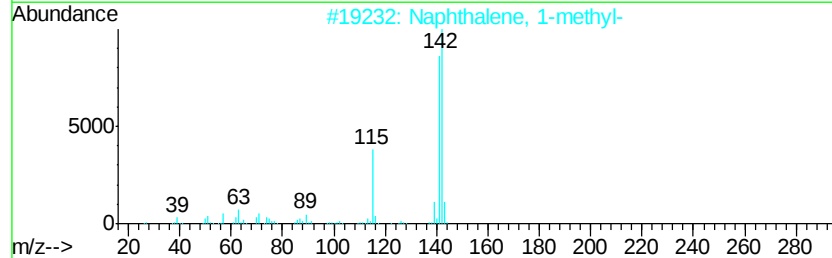
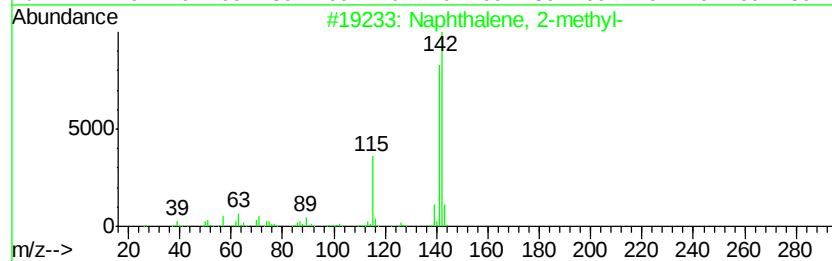
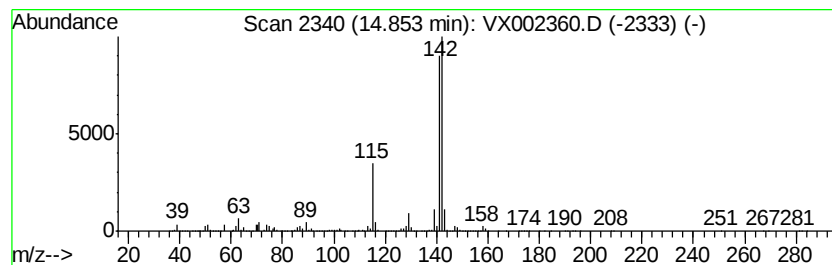
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.85	124.33 ug/l	2709410	1,4-Dichlorobenzene-d4	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3		1,4-Methanonaphthalene, 1,4-dihv...	142	C11H10	004453-90-1	93
4		Benzocycloheptatriene	142	C11H10	000264-09-5	90
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	59



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX060618\
Data File : VX002360.D
Acq On : 06 Jun 2018 18:36
Operator : JC/MD
Sample : J3312-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
W-16

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.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
Cyclopentene	2.81	87.3	ug/l	1166910	1	5.67	668336	50.0
Benzene, 1-etheny...	12.30	66.8	ug/l	1456590	4	12.09	1089600	50.0
Benzene, 4-ethyl-...	12.67	86.4	ug/l	1882110	4	12.09	1089600	50.0
Benzene, (2-methy...	12.71	33.5	ug/l	730947	4	12.09	1089600	50.0
Benzene, 1-etheny...	12.76	106.2	ug/l	2314930	4	12.09	1089600	50.0
Benzene, 1,2,3,4-...	12.98	80.3	ug/l	1750720	4	12.09	1089600	50.0
1-Phenyl-1-butene	13.35	176.5	ug/l	3845720	4	12.09	1089600	50.0
1H-Indene, 2,3-di...	13.73	33.9	ug/l	739047	4	12.09	1089600	50.0
Naphthalene, 2-me...	14.85	124.3	ug/l	2709410	4	12.09	1089600	50.0