

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX112919\
 Data File : VX013699.D
 Acq On : 29 Nov 2019 14:26
 Operator : JC/SP
 Sample : K6027-06
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-1A

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\82X112619W.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	1.070	26	30	42	rBV	2312726	2764365	39.32%	4.420%
2	1.393	79	83	88	rVB	137200	159538	2.27%	0.255%
3	1.783	142	147	158	rVB	286643	392604	5.58%	0.628%
4	1.997	177	182	191	rVB2	153323	326038	4.64%	0.521%
5	2.107	196	200	207	rVB	104324	149534	2.13%	0.239%
6	2.259	220	225	234	rVB	253909	402471	5.73%	0.643%
7	2.588	274	279	287	rVB	46534	85509	1.22%	0.137%
8	2.802	308	314	323	rBV3	80864	194036	2.76%	0.310%
9	2.881	323	327	330	rVV	59846	112384	1.60%	0.180%
10	2.930	330	335	344	rVB3	93855	221798	3.16%	0.355%
11	3.180	365	376	385	rBV3	169957	445010	6.33%	0.711%
12	3.807	473	479	489	rVB	65322	152232	2.17%	0.243%
13	3.917	489	497	504	rBV2	43890	109181	1.55%	0.175%
14	4.185	534	541	550	rVB2	59644	148671	2.11%	0.238%
15	4.393	564	575	587	rBV	165836	466242	6.63%	0.745%
16	4.521	587	596	604	rBV2	41335	109925	1.56%	0.176%
17	5.295	714	723	739	rVB	129079	405050	5.76%	0.648%
18	5.484	744	754	763	rVV	371254	1099906	15.65%	1.758%
19	5.575	763	769	773	rVV3	129524	374662	5.33%	0.599%
20	5.649	773	781	809	rVB	674783	2119185	30.15%	3.388%
21	5.929	819	827	837	rVB4	57413	181715	2.58%	0.291%
22	6.051	837	847	853	rBV	404240	1065032	15.15%	1.703%
23	6.136	853	861	873	rVB	2722059	7029896	100.00%	11.239%
24	6.289	873	886	892	rBV5	93003	331736	4.72%	0.530%
25	6.417	899	907	924	rVB	1043843	3051363	43.41%	4.878%
26	6.691	946	952	960	rBV6	32768	82850	1.18%	0.132%
27	6.850	969	978	987	rBV	1031305	2556688	36.37%	4.088%
28	6.929	988	991	1003	rVB4	151394	390163	5.55%	0.624%
29	7.063	1005	1013	1018	rBV2	38861	93145	1.32%	0.149%
30	7.252	1034	1044	1051	rBV2	190739	429809	6.11%	0.687%
31	7.453	1066	1077	1089	rBV4	332978	889065	12.65%	1.421%
32	7.728	1116	1122	1130	rVB2	59321	113108	1.61%	0.181%
33	7.929	1148	1155	1156	rBV2	74477	136047	1.94%	0.218%
34	7.953	1156	1159	1165	rVV3	87066	153178	2.18%	0.245%

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35	8.051	1165	1175	1186	rVB6	51518	184350	2.62%	0.295%
36	8.209	1186	1201	1206	rBV	439559	934487	13.29%	1.494%
37	8.258	1206	1209	1214	rVV	71934	128929	1.83%	0.206%
38	8.569	1254	1260	1268	rVB2	329297	645460	9.18%	1.032%
39	8.660	1268	1275	1277	rBV4	52589	91188	1.30%	0.146%
40	8.709	1277	1283	1290	rVV	2151894	3353637	47.71%	5.362%
41	8.782	1290	1295	1301	rVB	187444	293405	4.17%	0.469%
42	8.922	1313	1318	1323	rVV3	40152	107078	1.52%	0.171%
43	8.977	1323	1327	1331	rVV2	67396	125289	1.78%	0.200%
44	9.032	1331	1336	1341	rVV4	53752	127091	1.81%	0.203%
45	9.099	1342	1347	1351	rVV2	35768	76563	1.09%	0.122%
46	9.160	1353	1357	1362	rVV	52351	94329	1.34%	0.151%
47	9.221	1362	1367	1376	rVV	176578	287759	4.09%	0.460%
48	9.331	1382	1385	1389	rVB	51800	70797	1.01%	0.113%
49	9.514	1406	1415	1419	rVV2	90458	185027	2.63%	0.296%
50	9.556	1419	1422	1428	rVV4	48658	97287	1.38%	0.156%
51	9.812	1458	1464	1468	rBV4	39730	74922	1.07%	0.120%
52	10.105	1502	1512	1517	rBV	2013946	2829257	40.25%	4.523%
53	10.245	1531	1535	1546	rVB	1856034	2439325	34.70%	3.900%
54	10.355	1546	1553	1559	rVB	2086823	2747381	39.08%	4.392%
55	10.696	1603	1609	1614	rBV	689815	925767	13.17%	1.480%
56	10.946	1644	1650	1656	rVB2	34947	72275	1.03%	0.116%
57	11.013	1656	1661	1666	rBV	66311	91776	1.31%	0.147%
58	11.135	1675	1681	1686	rBV	1622580	2127479	30.26%	3.401%
59	11.355	1714	1717	1721	rVB	64656	77502	1.10%	0.124%
60	11.428	1721	1729	1731	rBV	335753	466406	6.63%	0.746%
61	11.446	1731	1732	1737	rVV	348020	354101	5.04%	0.566%
62	11.501	1737	1741	1746	rVB	728673	849463	12.08%	1.358%
63	11.666	1761	1768	1776	rBV	1326692	1727392	24.57%	2.762%
64	11.800	1786	1790	1796	rVV	1357316	1709613	24.32%	2.733%
65	12.074	1830	1835	1840	rVV	2106996	2573168	36.60%	4.114%
66	12.135	1840	1845	1850	rVB	1005292	1254702	17.85%	2.006%
67	12.294	1864	1871	1879	rBV	1483686	1962752	27.92%	3.138%
68	12.367	1879	1883	1890	rVB2	405734	511594	7.28%	0.818%
69	12.458	1894	1898	1903	rVV3	182972	255509	3.63%	0.409%
70	12.513	1903	1907	1913	rVB2	159483	225365	3.21%	0.360%
71	12.580	1913	1918	1920	rBV	222215	273040	3.88%	0.437%

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72	12.605	1920	1922	1927	rVB	190326	198025	2.82%	0.317%
73	12.666	1927	1932	1935	rBV	518443	634740	9.03%	1.015%
74	12.751	1941	1946	1953	rVB2	535663	782012	11.12%	1.250%
75	12.897	1965	1970	1975	rBV2	147185	182968	2.60%	0.293%
76	12.970	1977	1982	1986	rBV	388777	485241	6.90%	0.776%
77	13.013	1986	1989	1995	rVB	496066	576457	8.20%	0.922%
78	13.092	1995	2002	2005	rBV3	60839	119302	1.70%	0.191%
79	13.220	2020	2023	2028	rVV	153948	192873	2.74%	0.308%
80	13.342	2039	2043	2049	rVB2	1010668	1303601	18.54%	2.084%
81	13.476	2061	2065	2071	rVB2	115784	151212	2.15%	0.242%
82	13.598	2081	2085	2088	rVV	76192	111622	1.59%	0.178%
83	13.635	2088	2091	2097	rVV3	112068	204669	2.91%	0.327%
84	13.696	2097	2101	2102	rVV	70126	92705	1.32%	0.148%
85	13.714	2102	2104	2110	rVB2	107165	148270	2.11%	0.237%
86	13.830	2116	2123	2131	rVB	132962	197851	2.81%	0.316%
87	14.269	2191	2195	2201	rBV	45255	72862	1.04%	0.116%

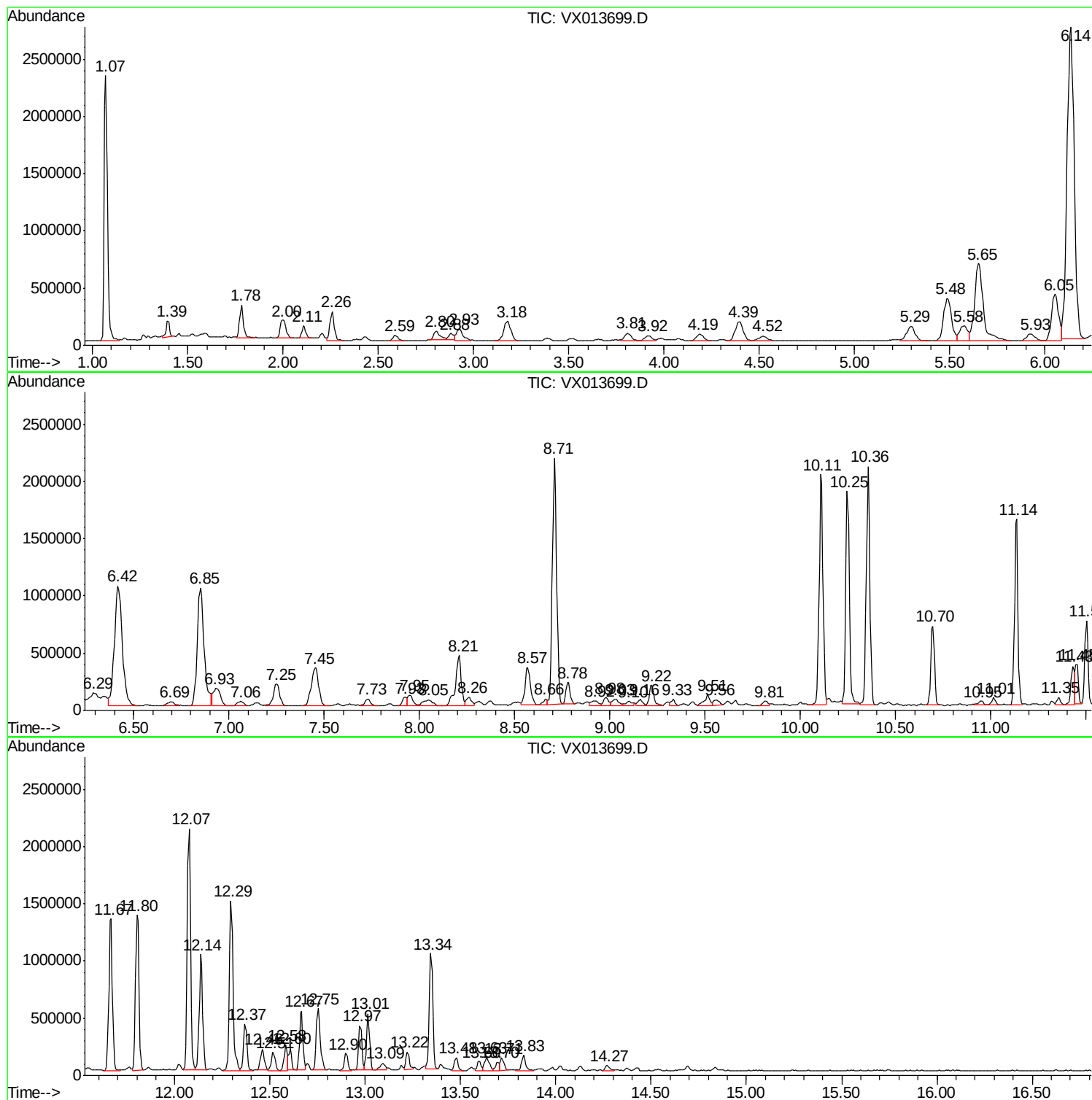
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.M
Quant Title : SW846 8260

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



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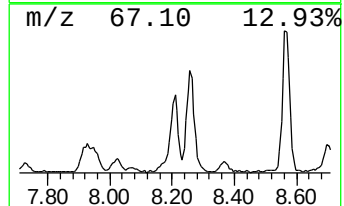
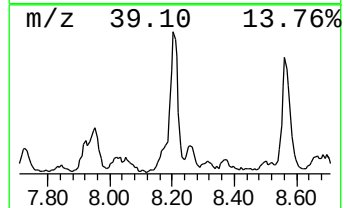
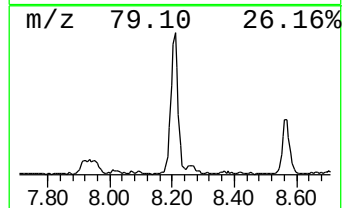
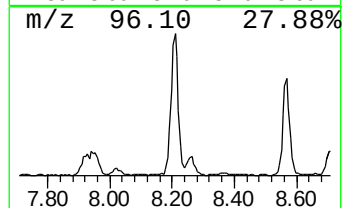
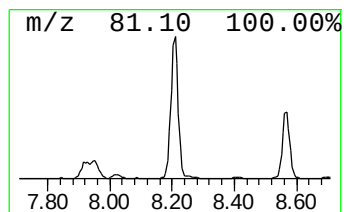
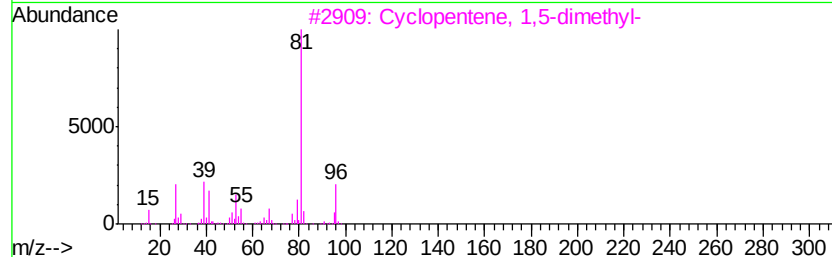
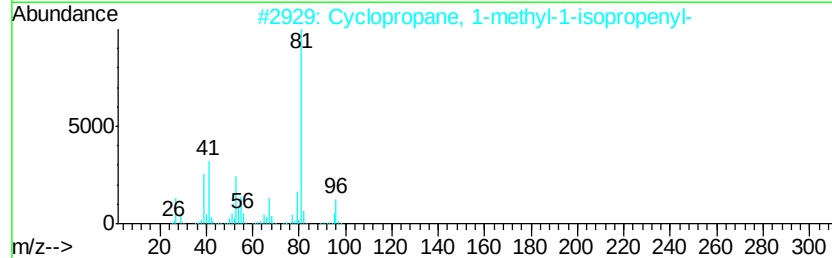
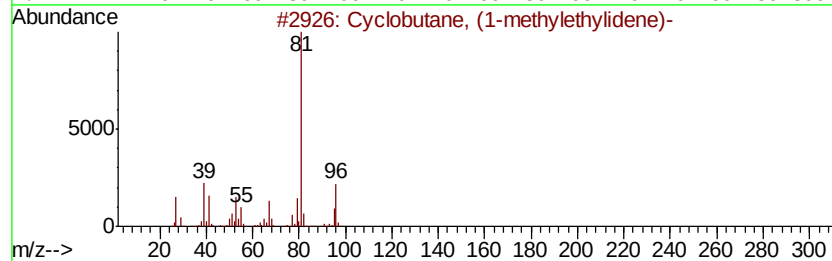
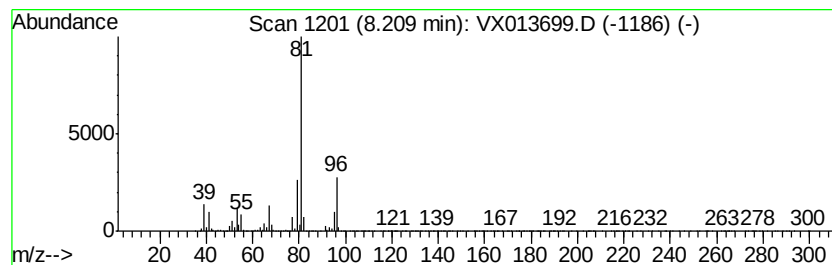
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.M
Quant Title : SW846 8260

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

Peak Number 2 Cyclobutane, (1-methylethyl... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.21	18.28 ug/l	934487	1,4-Difluorobenzene	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclobutane, (1-methylethylidene)-	96	C7H12	001528-22-9	90
2		Cyclopropane, 1-methyl-1-isopropyl-	96	C7H12	003422-07-9	86
3		Cyclopentene, 1,5-dimethyl-	96	C7H12	016491-15-9	83
4		3,5-Dimethylcyclopentene	96	C7H12	007459-71-4	72
5		1,4-Pentadiene, 3,3-dimethyl-	96	C7H12	001112-35-2	72



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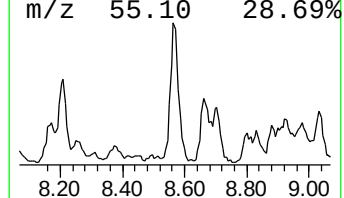
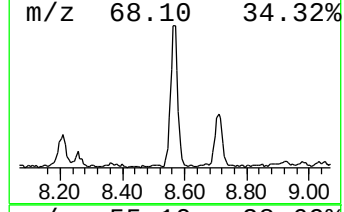
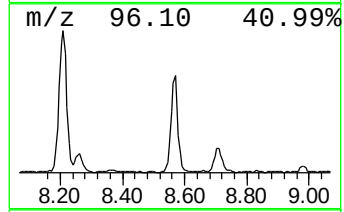
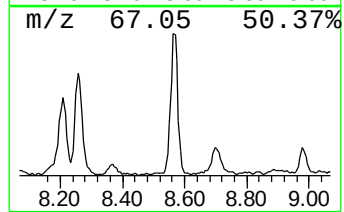
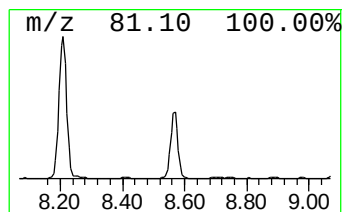
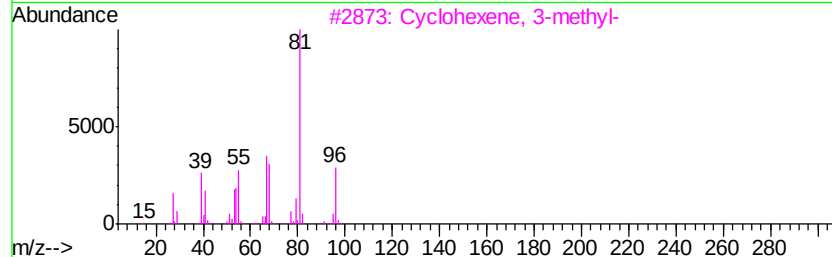
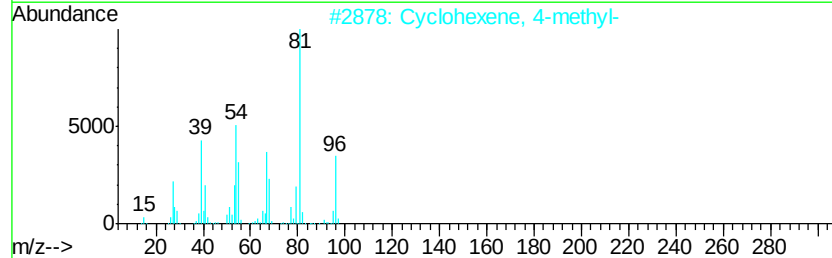
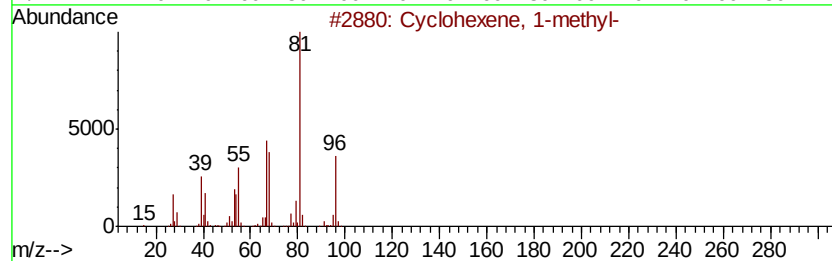
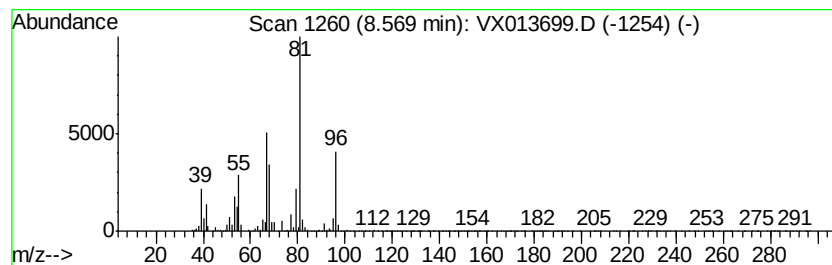
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Cyclohexene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.57	11.41 ug/l	645460	Chlorobenzene-d5	10.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexene, 1-methyl-	96	C7H12	000591-49-1	94
2		Cyclohexene, 4-methyl-	96	C7H12	000591-47-9	87
3		Cyclohexene, 3-methyl-	96	C7H12	000591-48-0	81
4		2,4-Hexadiene, 2-methyl-	96	C7H12	028823-41-8	76
5		2,4-Hexadiene, 3-methyl-	96	C7H12	028823-42-9	64



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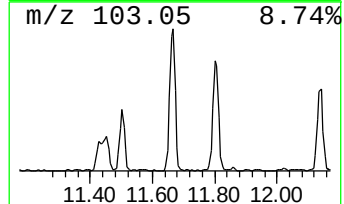
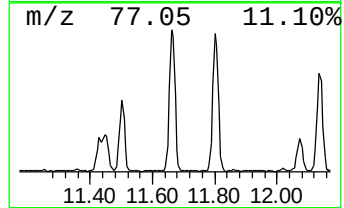
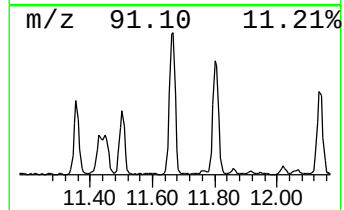
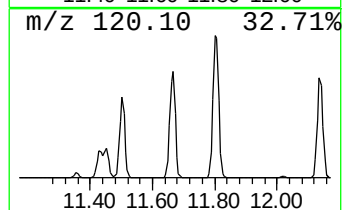
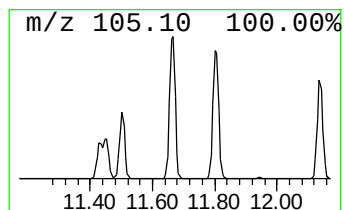
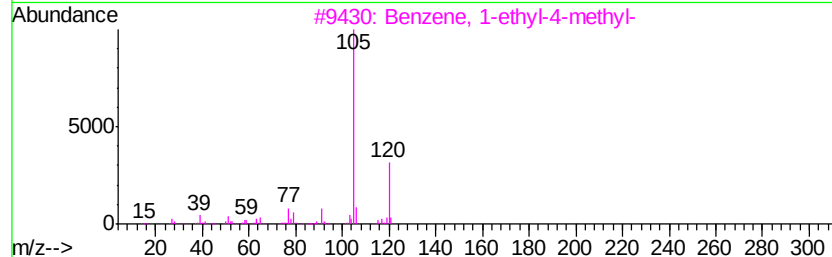
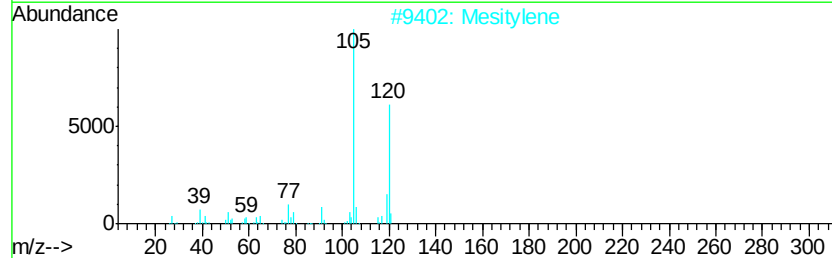
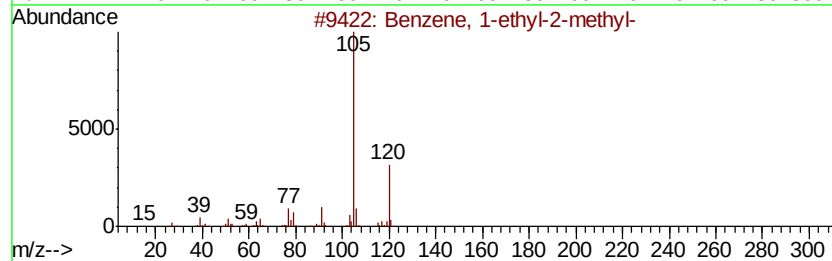
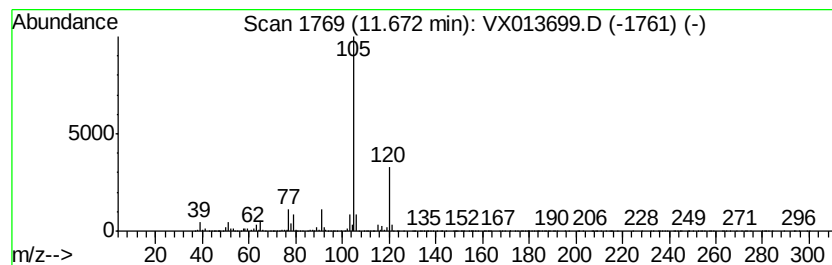
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.67	33.57 ug/l	1727390	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
2		Mesitylene	120	C9H12	000108-67-8	90
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90



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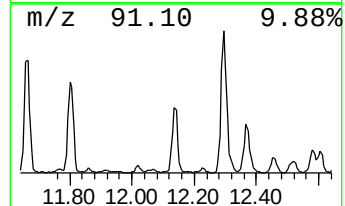
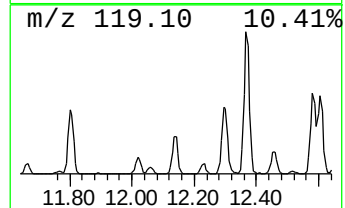
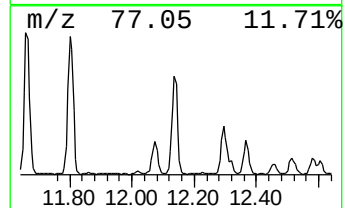
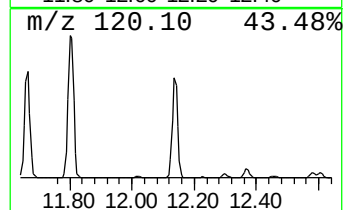
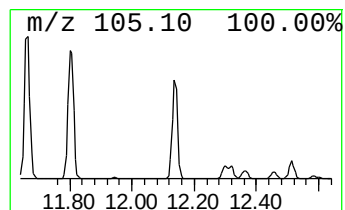
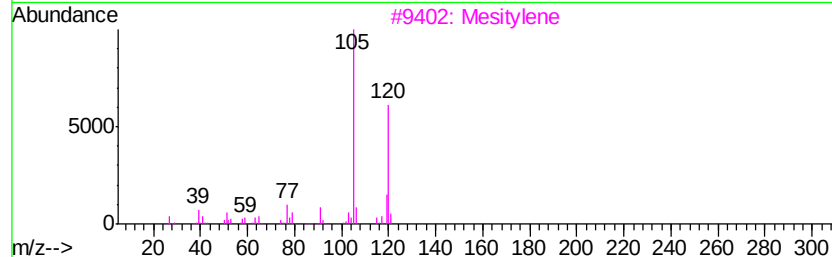
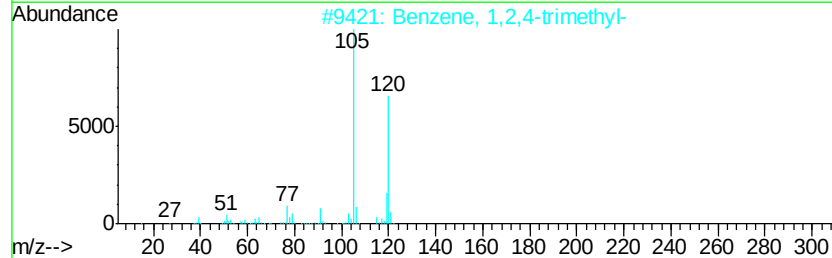
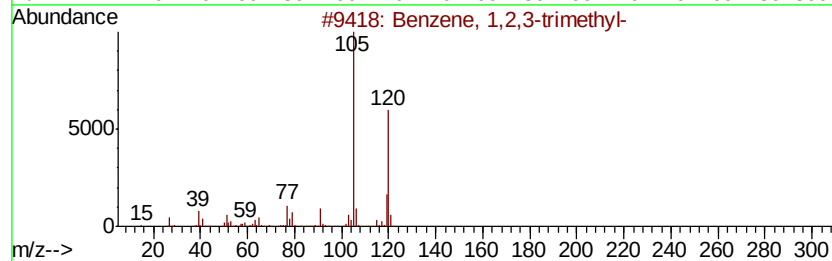
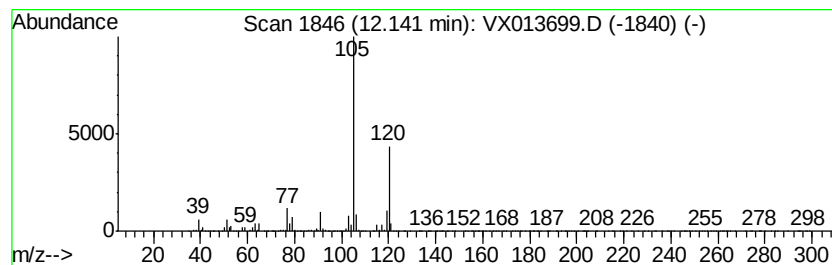
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	24.38 ug/l	1254700	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3		Mesitylene	120	C9H12	000108-67-8	93
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX112919\
Data File : VX013699.D
Acq On : 29 Nov 2019 14:26
Operator : JC/SP
Sample : K6027-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-1A

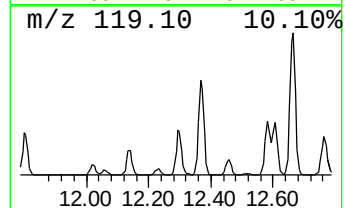
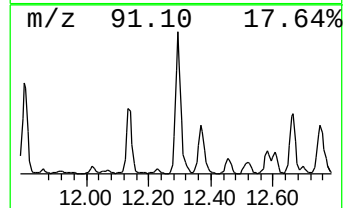
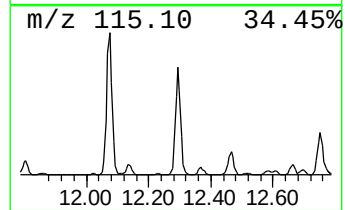
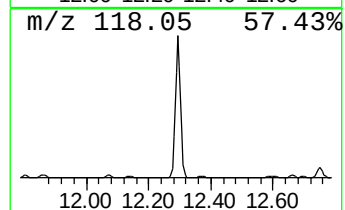
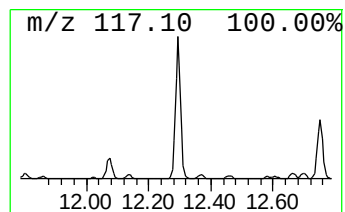
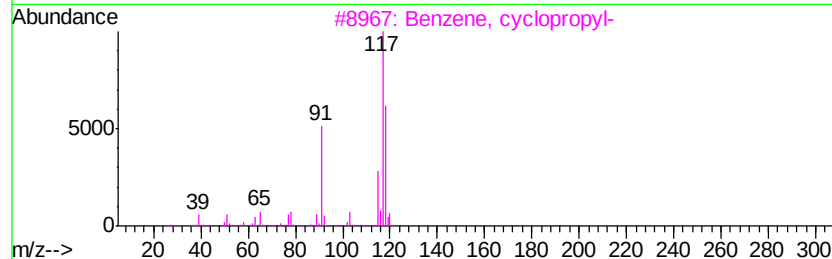
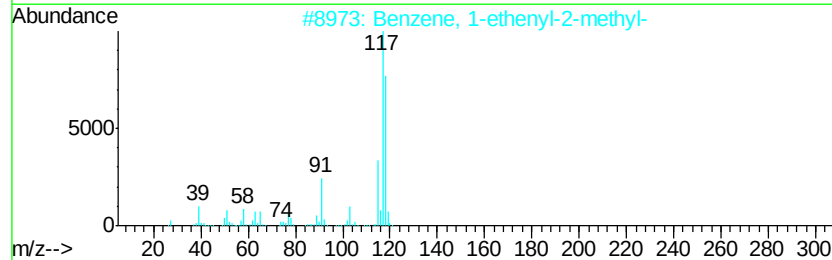
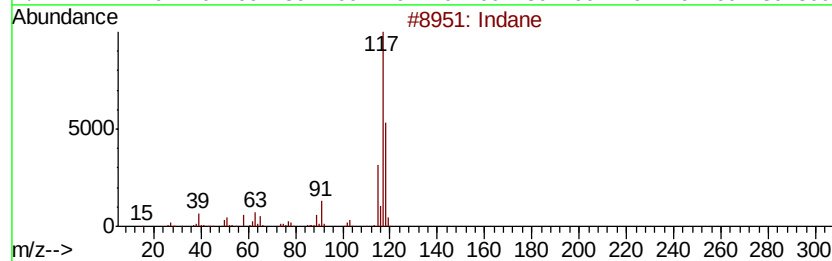
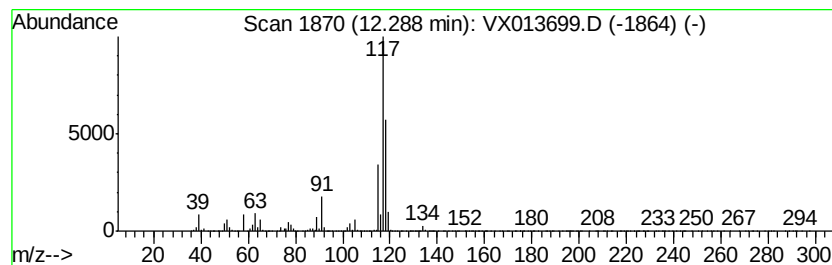
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Quant Title : SW846 8260

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

Peak Number 6 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.29	38.14 ug/l	1962750	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indane	118	C9H10	000496-11-7	76
2		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	76
3		Benzene, cyclopropyl-	118	C9H10	000873-49-4	68
4		Deltacyclene	118	C9H10	007785-10-6	58
5		Benzene, 2-propenyl-	118	C9H10	000300-57-2	55



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX112919\
Data File : VX013699.D
Acq On : 29 Nov 2019 14:26
Operator : JC/SP
Sample : K6027-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-1A

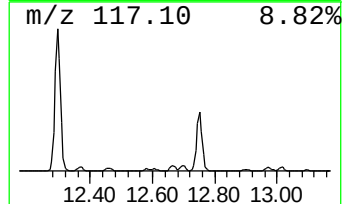
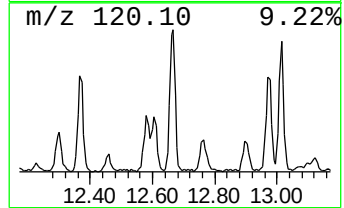
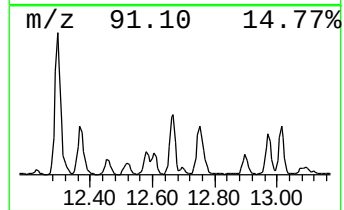
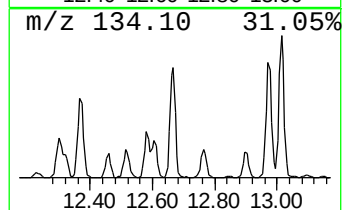
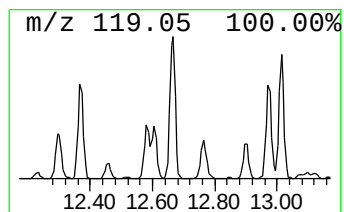
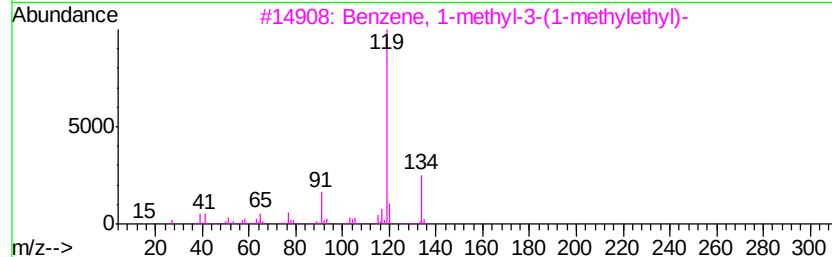
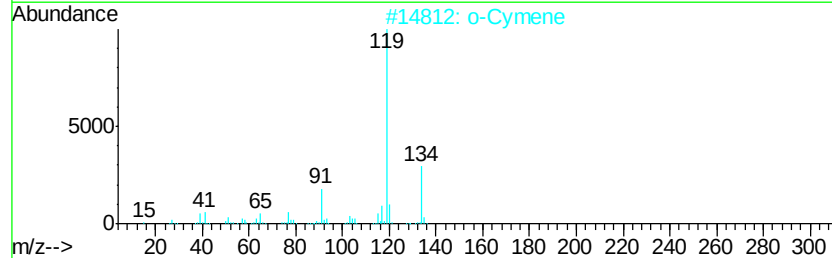
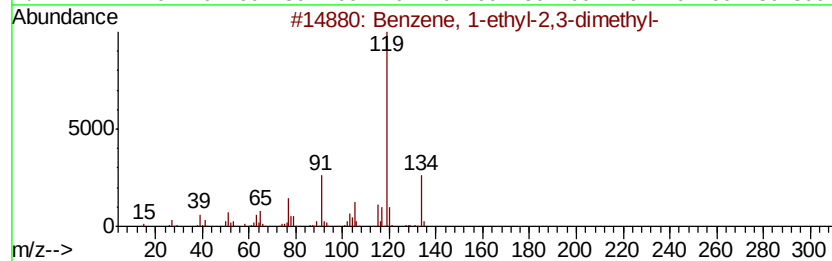
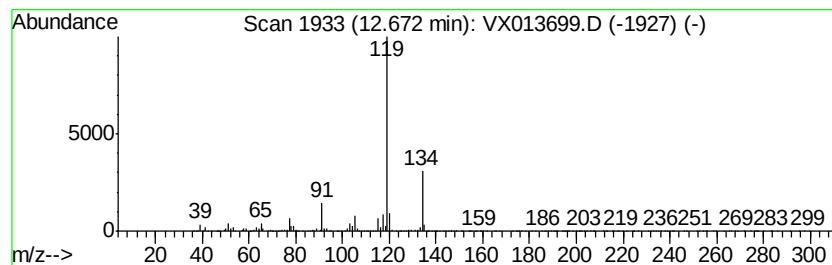
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	12.33 ug/l	634740	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
2		o-Cymene	134	C10H14	000527-84-4	95
3		Benzene, 1-methyl-3-(1-methylethyl-)	134	C10H14	000535-77-3	95
4		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX112919\
Data File : VX013699.D
Acq On : 29 Nov 2019 14:26
Operator : JC/SP
Sample : K6027-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-1A

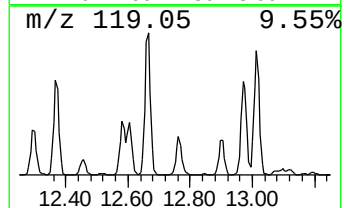
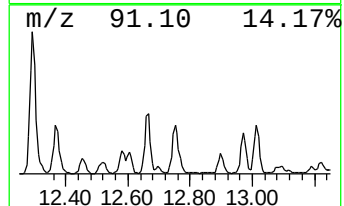
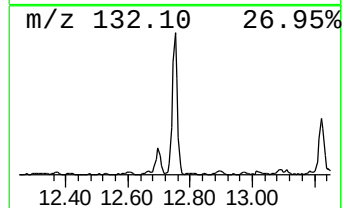
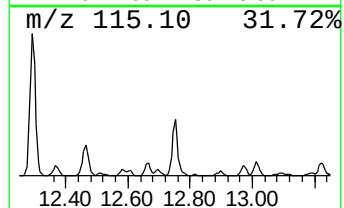
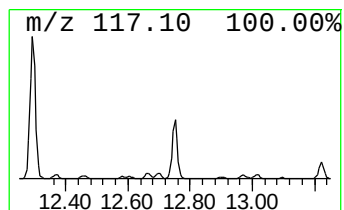
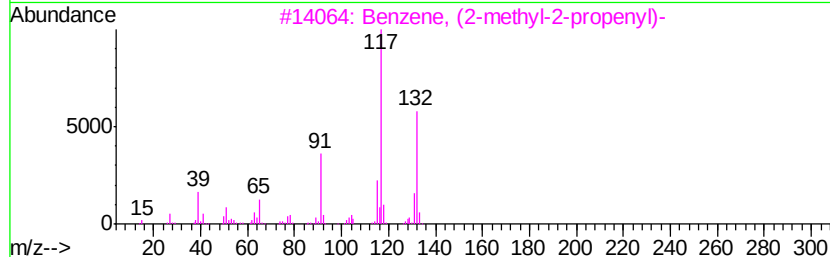
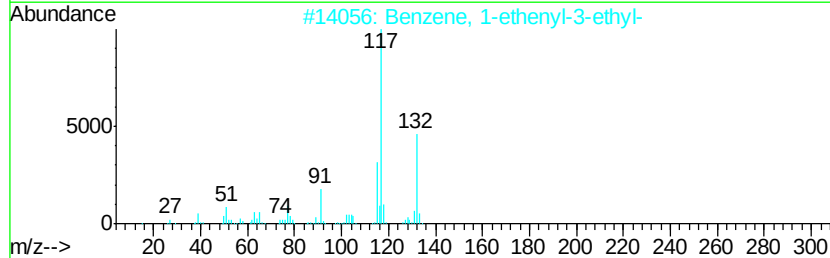
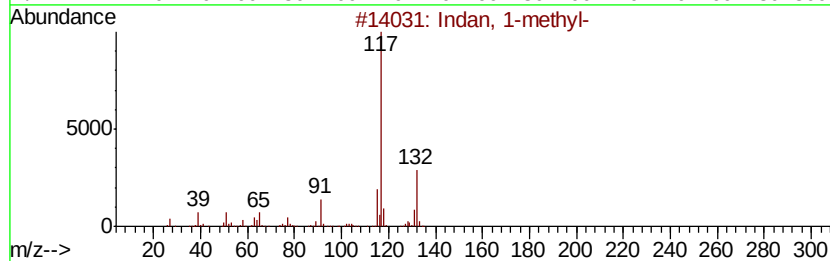
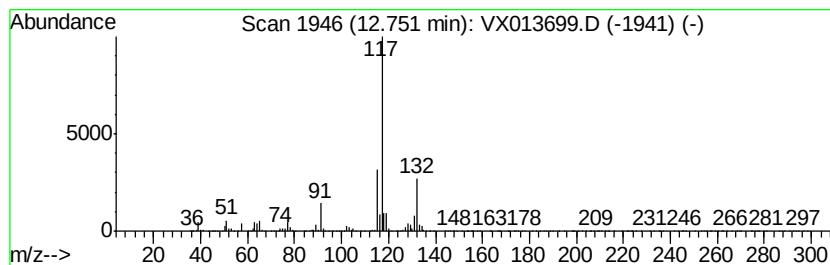
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Quant Title : SW846 8260

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

Peak Number 8 Indan, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.75	15.20 ug/l	782012	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indan, 1-methyl-	132	C10H12	000767-58-8	87
2		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
3		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	83
4		1-Phenyl-1-butene	132	C10H12	000824-90-8	83
5		Benzene, 2-butenyl-	132	C10H12	001560-06-1	83



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX112919\
Data File : VX013699.D
Acq On : 29 Nov 2019 14:26
Operator : JC/SP
Sample : K6027-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-1A

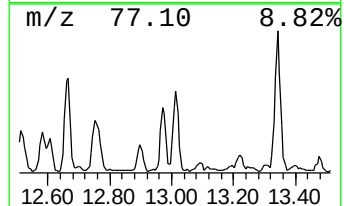
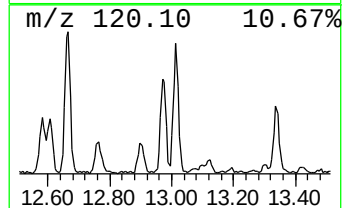
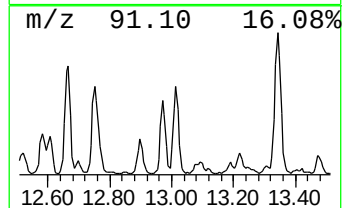
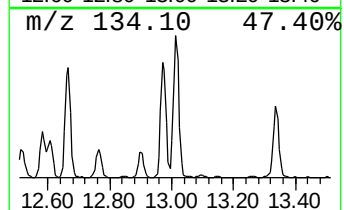
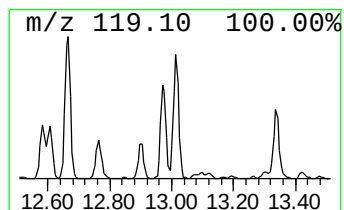
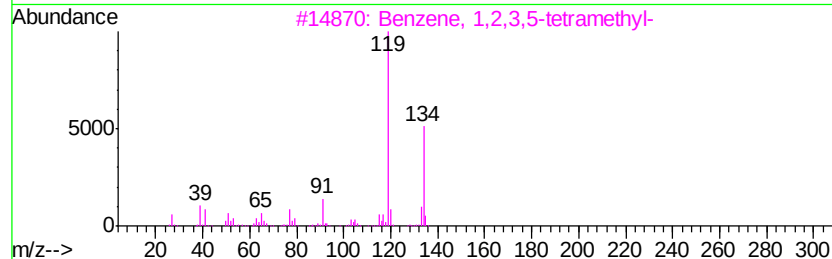
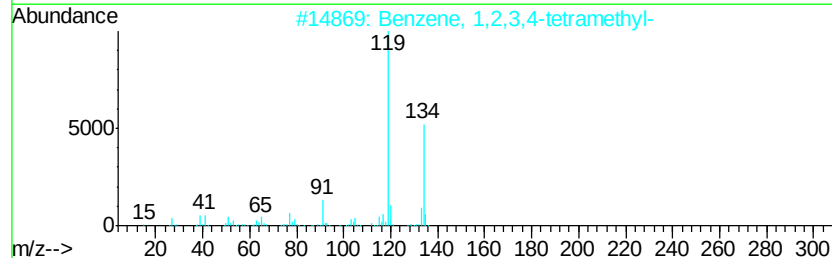
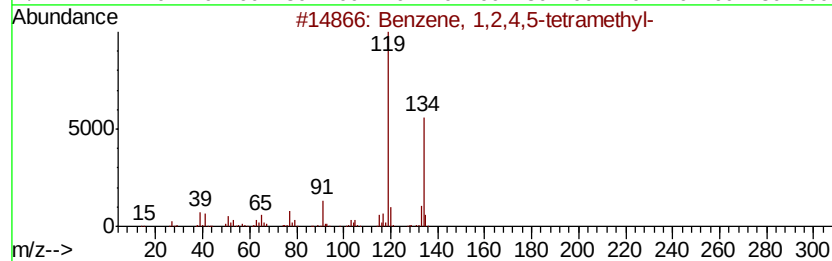
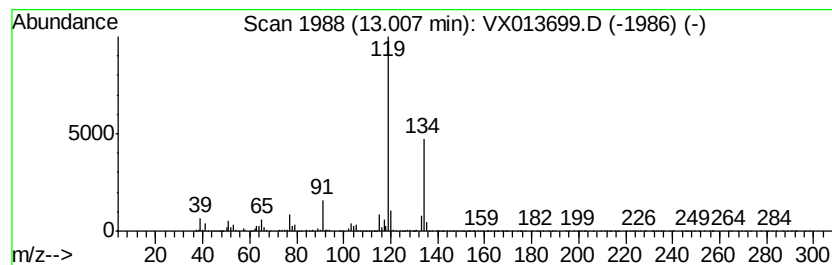
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.01	11.20 ug/l	576457	1,4-Dichlorobenzene-d4	12.07

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		o-Cymene	134	C10H14	000527-84-4	94
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX112919\
Data File : VX013699.D
Acq On : 29 Nov 2019 14:26
Operator : JC/SP
Sample : K6027-06
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-1A

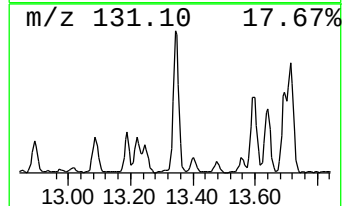
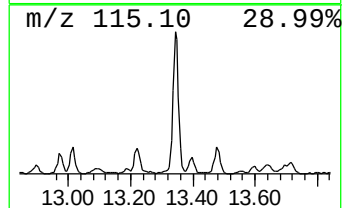
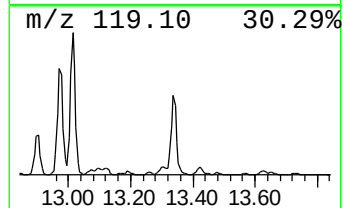
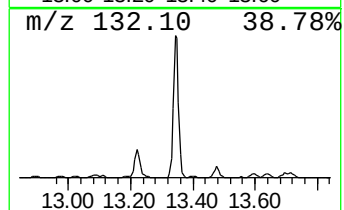
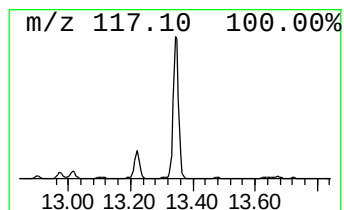
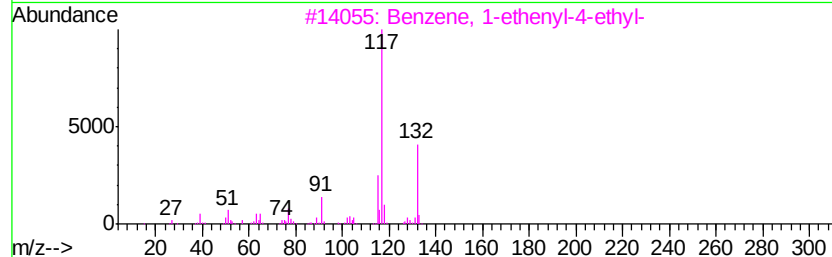
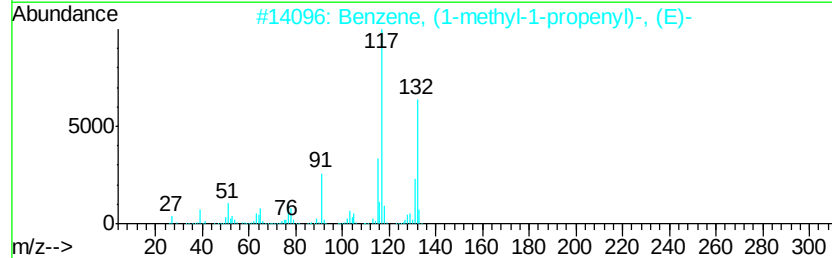
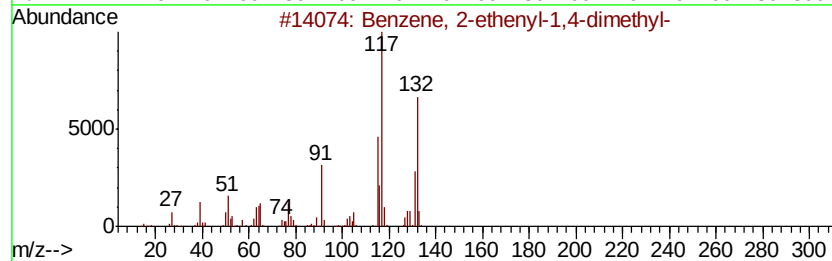
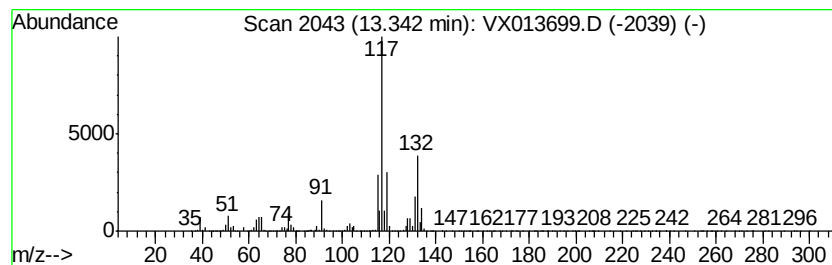
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.34	25.33 ug/l	1303600	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	91
3		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	90
4		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
5		3-Phenylbut-1-ene	132	C10H12	000934-10-1	87



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX112919\
Data File : VX013699.D
Acq On : 29 Nov 2019 14:26
Operator : JC/SP
Sample : K6027-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-1A

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112619W.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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Cyclohexene, 1-me...	8.57	11.4	ug/l	645460	3	10.11	2829260	50.0
Benzene, 1-ethyl-...	11.67	33.6	ug/l	1727390	4	12.07	2573170	50.0
Benzene, 1,2,3-tr...	12.14	24.4	ug/l	1254700	4	12.07	2573170	50.0
Indane	12.29	38.1	ug/l	1962750	4	12.07	2573170	50.0
Benzene, 1-ethyl-...	12.67	12.3	ug/l	634740	4	12.07	2573170	50.0
Indan, 1-methyl-	12.75	15.2	ug/l	782012	4	12.07	2573170	50.0
Benzene, 1,2,4,5-...	13.01	11.2	ug/l	576457	4	12.07	2573170	50.0
Benzene, 2-etheny...	13.34	25.3	ug/l	1303600	4	12.07	2573170	50.0