

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX052219\
 Data File : VX009834.D
 Acq On : 23 May 2019 04:01
 Operator : JC/SP
 Sample : K2976-15 50X
 Misc : 5.39g/5mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 EP1-SB-MW-IR(14-17)

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\82X042919W.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	5.501	702	713	726	rBV2	102693	298378	10.31%	0.880%
2	5.665	728	740	752	rVV	184742	521368	18.01%	1.537%
3	6.061	792	805	821	rBV	132142	347597	12.01%	1.025%
4	6.860	924	936	951	rBV	303034	723448	25.00%	2.133%
5	8.713	1234	1240	1255	rVB	662680	1027286	35.49%	3.029%
6	9.043	1287	1294	1300	rVB2	22688	40321	1.39%	0.119%
7	9.445	1352	1360	1369	rBV	66663	108515	3.75%	0.320%
8	9.561	1373	1379	1385	rBV2	49517	109952	3.80%	0.324%
9	9.622	1385	1389	1393	rVB	87524	120742	4.17%	0.356%
10	9.677	1393	1398	1402	rBV	146700	220291	7.61%	0.650%
11	9.817	1415	1421	1425	rBV2	25537	40566	1.40%	0.120%
12	9.878	1425	1431	1437	rVV5	98594	245510	8.48%	0.724%
13	9.957	1440	1444	1448	rVB	83583	115333	3.98%	0.340%
14	10.061	1455	1461	1465	rBV	177182	255426	8.83%	0.753%
15	10.110	1465	1469	1475	rVB	635866	835618	28.87%	2.464%
16	10.177	1475	1480	1485	rVB5	27357	61864	2.14%	0.182%
17	10.244	1485	1491	1497	rBV4	69651	139485	4.82%	0.411%
18	10.329	1497	1505	1507	rBV2	125866	215567	7.45%	0.636%
19	10.372	1507	1512	1516	rVB2	171208	302631	10.46%	0.892%
20	10.408	1516	1518	1530	rVB	121595	208842	7.22%	0.616%
21	10.512	1530	1535	1540	rBV	80324	121566	4.20%	0.358%
22	10.585	1540	1547	1549	rBV2	30506	41777	1.44%	0.123%
23	10.640	1549	1556	1564	rBV2	318168	551318	19.05%	1.626%
24	10.725	1564	1570	1573	rBV2	180717	277952	9.60%	0.820%
25	10.768	1574	1577	1580	rVB2	98521	123093	4.25%	0.363%
26	10.835	1580	1588	1592	rBV	917821	1331158	45.99%	3.925%
27	10.908	1592	1600	1604	rBV	593865	1020446	35.26%	3.009%
28	10.945	1604	1606	1611	rVB	366463	467130	16.14%	1.377%
29	11.012	1611	1617	1621	rBV4	140148	247904	8.57%	0.731%
30	11.055	1621	1624	1626	rBV2	131573	188266	6.50%	0.555%
31	11.079	1626	1628	1632	rVB	241577	288623	9.97%	0.851%
32	11.134	1632	1637	1642	rVB	1081155	1424021	49.20%	4.199%
33	11.183	1642	1645	1648	rVB	89762	96094	3.32%	0.283%
34	11.219	1648	1651	1653	rBV	73645	98078	3.39%	0.289%

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35	11.274	1653	1660	1669	rVB3	527569	907439	31.35%	2.676%
36	11.359	1669	1674	1676	rBV	200600	253988	8.78%	0.749%
37	11.390	1676	1679	1682	rBV2	105059	137355	4.75%	0.405%
38	11.432	1682	1686	1692	rBV2	458039	970746	33.54%	2.862%
39	11.487	1692	1695	1696	rBV	404710	482583	16.67%	1.423%
40	11.506	1696	1698	1701	rVB	423717	395317	13.66%	1.166%
41	11.579	1707	1710	1715	rVB4	67363	90572	3.13%	0.267%
42	11.628	1715	1718	1720	rBV	69313	79741	2.76%	0.235%
43	11.682	1720	1727	1732	rVV5	293477	529831	18.31%	1.562%
44	11.737	1732	1736	1738	rVV2	270311	371260	12.83%	1.095%
45	11.768	1738	1741	1744	rVV	1363996	1587230	54.84%	4.680%
46	11.804	1744	1747	1757	rVB	2062804	2894335	100.00%	8.534%
47	11.890	1757	1761	1763	rBV	227802	309909	10.71%	0.914%
48	11.945	1763	1770	1774	rVV3	521759	953508	32.94%	2.811%
49	11.999	1774	1779	1781	rVV	566085	774587	26.76%	2.284%
50	12.024	1781	1783	1787	rVV4	379061	493924	17.07%	1.456%
51	12.079	1787	1792	1797	rVV3	807353	1414217	48.86%	4.170%
52	12.140	1797	1802	1808	rVB2	568327	852430	29.45%	2.513%
53	12.213	1809	1814	1819	rVB5	100497	181124	6.26%	0.534%
54	12.323	1823	1832	1835	rBV2	476162	930952	32.16%	2.745%
55	12.371	1835	1840	1846	rVB3	787219	1338141	46.23%	3.946%
56	12.475	1846	1857	1861	rBV3	215297	612577	21.16%	1.806%
57	12.518	1861	1864	1870	rVB4	290415	522365	18.05%	1.540%
58	12.585	1870	1875	1876	rBV	313206	362174	12.51%	1.068%
59	12.609	1876	1879	1882	rVV2	369555	525006	18.14%	1.548%
60	12.664	1886	1888	1891	rVB	228616	221728	7.66%	0.654%
61	12.737	1896	1900	1902	rBV2	189913	268416	9.27%	0.791%
62	12.774	1903	1906	1911	rVB4	203692	363343	12.55%	1.071%
63	12.816	1911	1913	1915	rBV2	34902	42604	1.47%	0.126%
64	12.847	1915	1918	1920	rBV3	73294	88772	3.07%	0.262%
65	12.871	1920	1922	1925	rVB	59224	51693	1.79%	0.152%
66	12.938	1930	1933	1937	rBV3	217770	355030	12.27%	1.047%
67	12.975	1937	1939	1942	rVB	150910	149890	5.18%	0.442%
68	13.018	1942	1946	1948	rBV	224851	268131	9.26%	0.791%
69	13.079	1953	1956	1961	rVV3	96942	222955	7.70%	0.657%
70	13.121	1961	1963	1965	rVV	73351	87680	3.03%	0.259%
71	13.146	1965	1967	1972	rVV	65975	100647	3.48%	0.297%

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Integration Parameters: RTEINT.P

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72	13.213	1972	1978	1982	rVV4	86763	189125	6.53%	0.558%
73	13.261	1983	1986	1989	rVV3	56058	71961	2.49%	0.212%
74	13.304	1989	1993	1996	rVV2	76581	113806	3.93%	0.336%
75	13.341	1996	1999	2006	rVV2	285844	448391	15.49%	1.322%
76	13.432	2010	2014	2017	rVV2	44883	78254	2.70%	0.231%
77	13.475	2017	2021	2030	rVB	207737	331421	11.45%	0.977%
78	13.560	2030	2035	2038	rBV2	27742	42939	1.48%	0.127%
79	13.639	2044	2048	2053	rBV2	38784	68973	2.38%	0.203%
80	13.719	2059	2061	2066	rVB2	27395	35951	1.24%	0.106%
81	13.834	2071	2080	2086	rVB	69558	126295	4.36%	0.372%

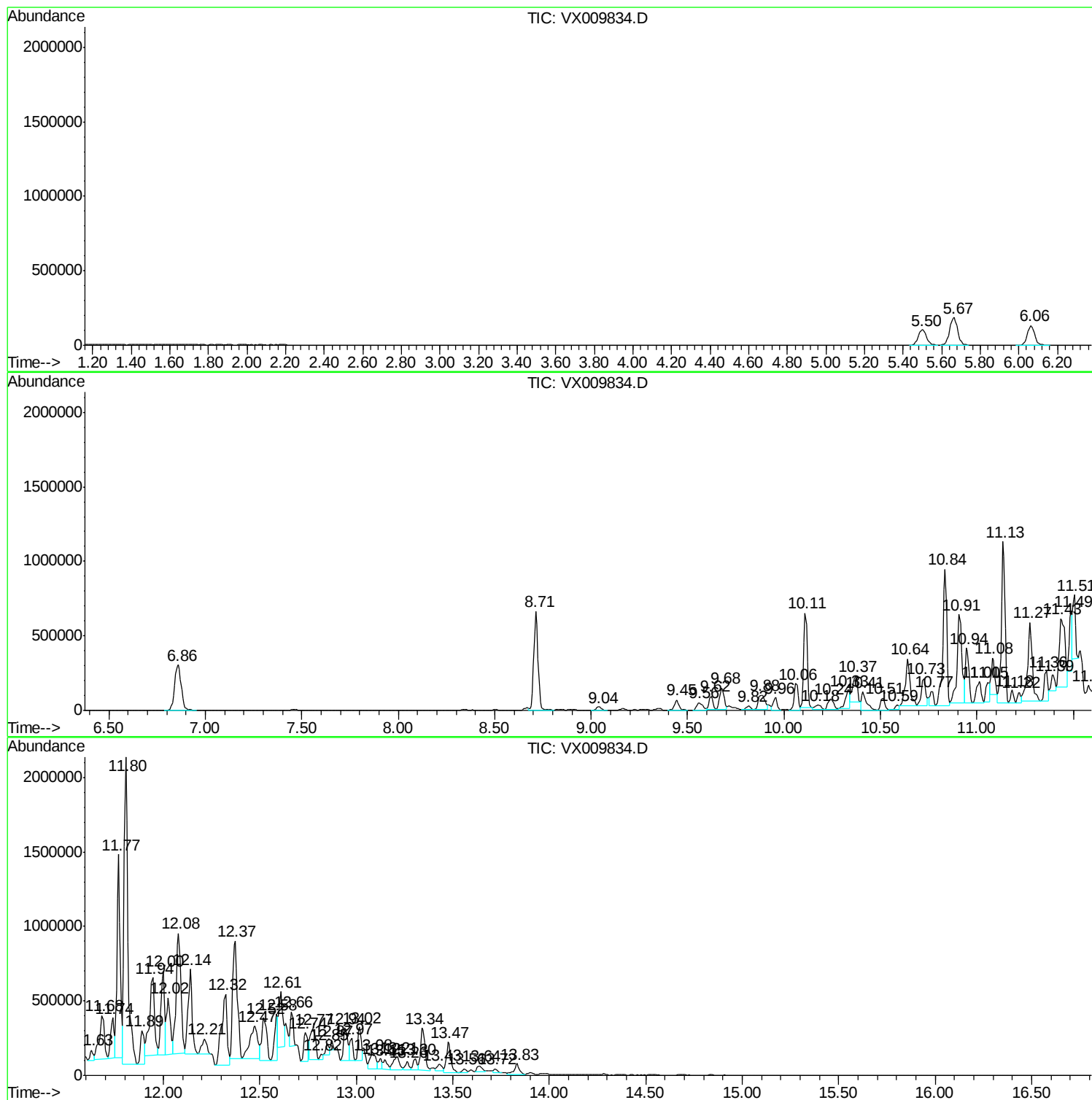
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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



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ALS Vial : 46 Sample Multiplier: 1

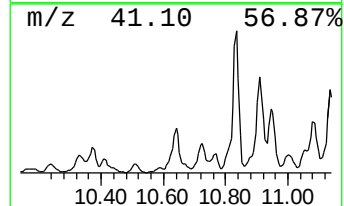
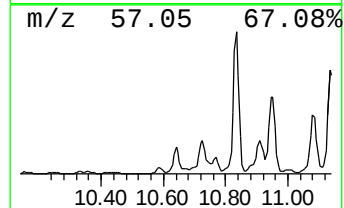
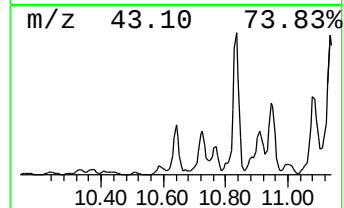
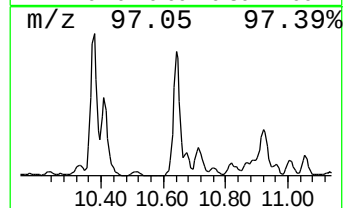
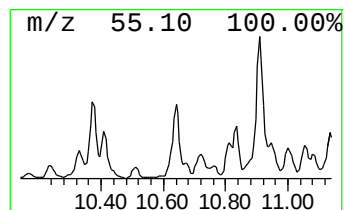
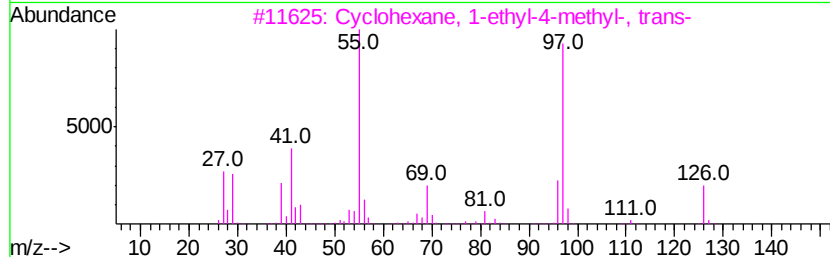
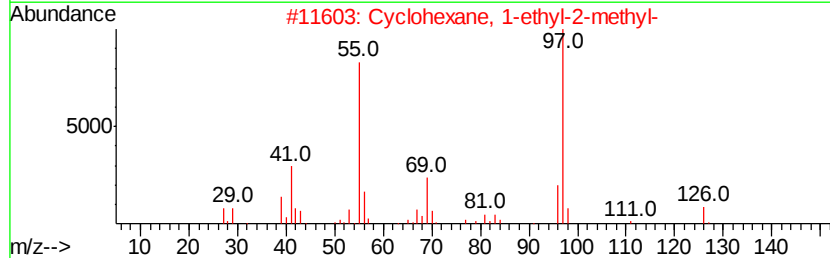
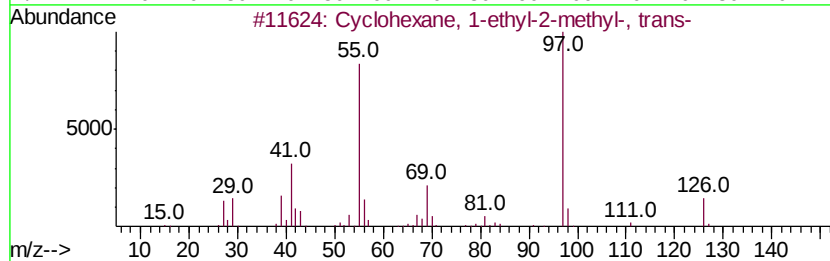
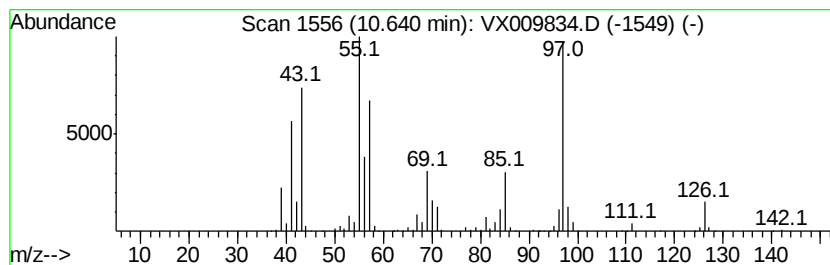
Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclohexane, 1-ethyl-2-meth... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.64	32.99 ug/l	551318	Chlorobenzene-d5	10.11		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	70
2		Cyclohexane, 1-ethyl-2-methyl-	126	C9H18	003728-54-9	62
3		Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	58
4		Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	58
5		1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	52



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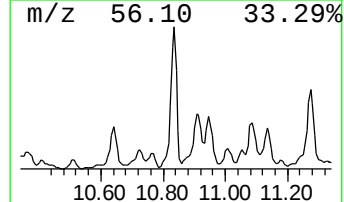
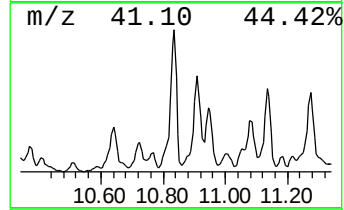
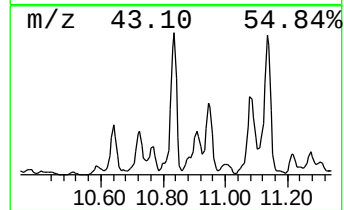
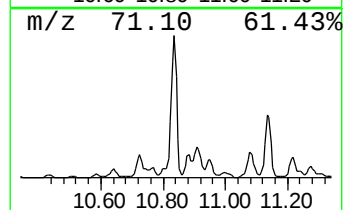
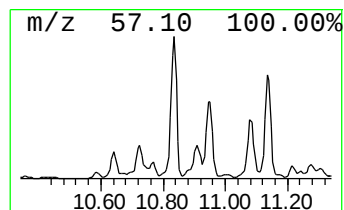
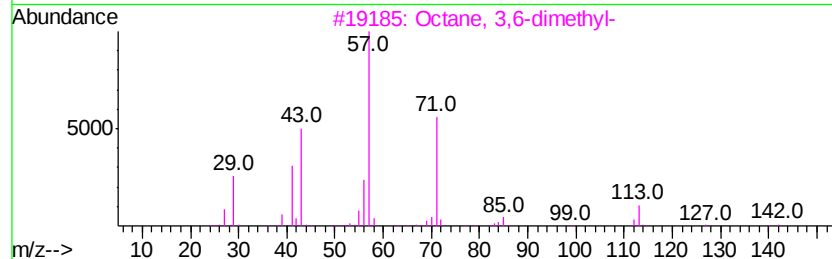
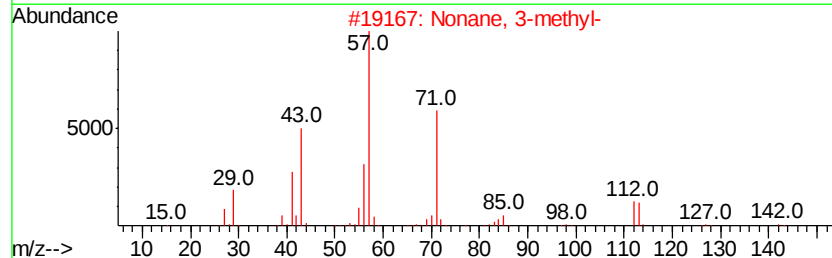
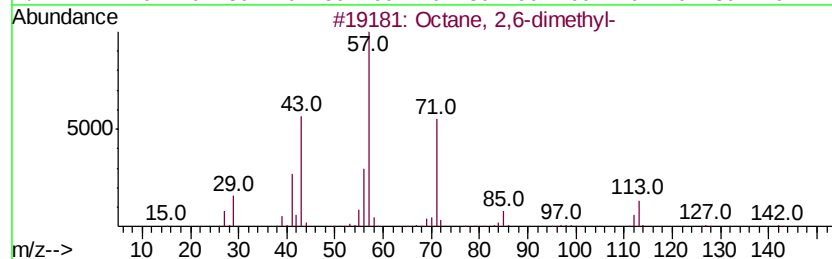
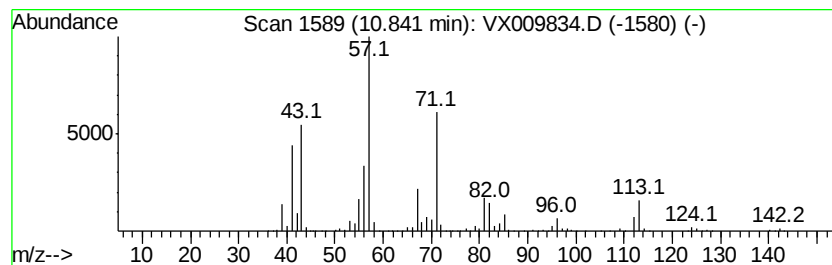
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Octane, 2,6-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.84	79.65 ug/l	1331160	Chlorobenzene-d5	10.11		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	94
2		Nonane, 3-methyl-	142	C10H22	005911-04-6	70
3		Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	70
4		Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	49
5		Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	43



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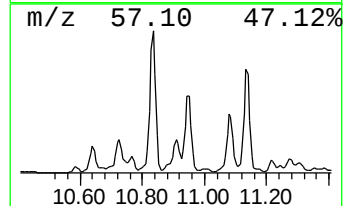
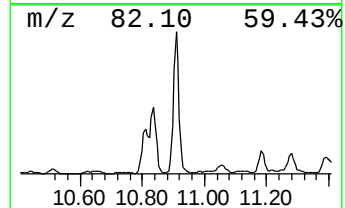
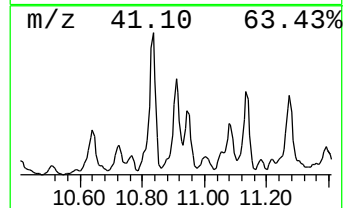
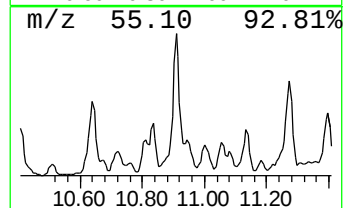
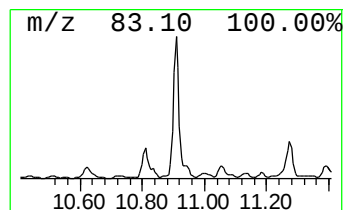
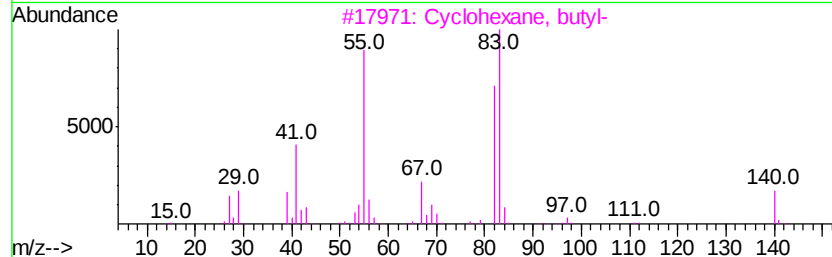
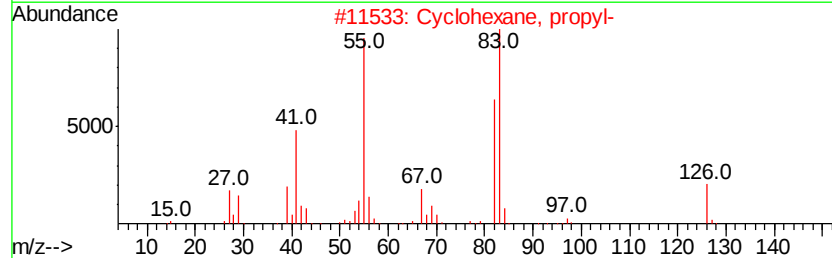
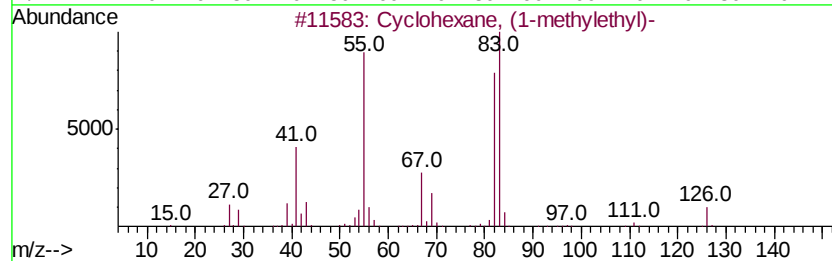
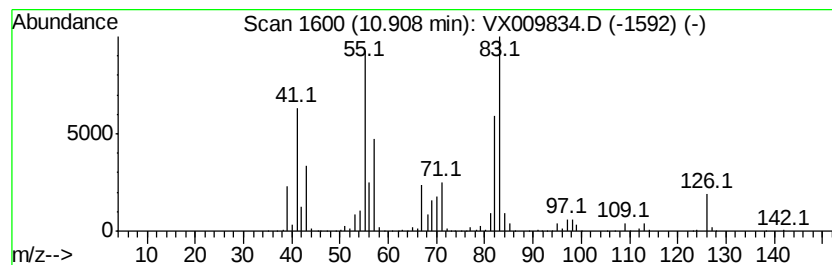
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Quant Title : SW846 8260

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

Peak Number 3 Cyclohexane, (1-methylethyl)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.91	61.06 ug/l	1020450	Chlorobenzene-d5	10.11
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclohexane, (1-methylethyl)-	126 C9H18	000696-29-7 81
2		Cyclohexane, propyl-	126 C9H18	001678-92-8 81
3		Cyclohexane, butyl-	140 C10H20	001678-93-9 59
4		Cyclohexane, octyl-	196 C14H28	001795-15-9 59
5		Cyclohexane, ethyl-	112 C8H16	001678-91-7 53



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EP1-SB-MW-IR(14-17)

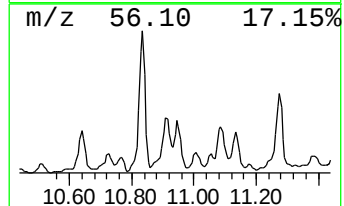
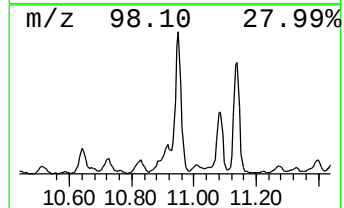
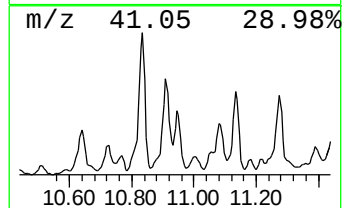
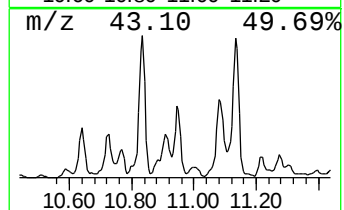
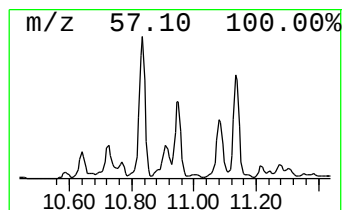
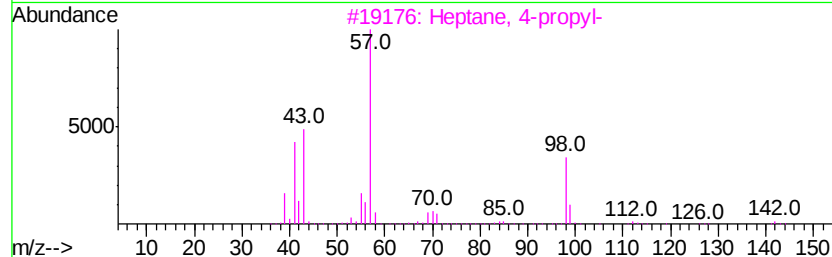
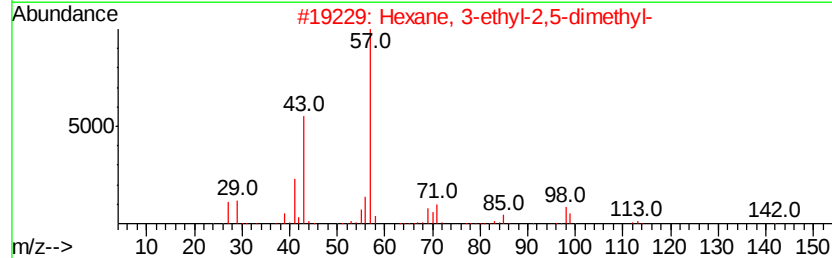
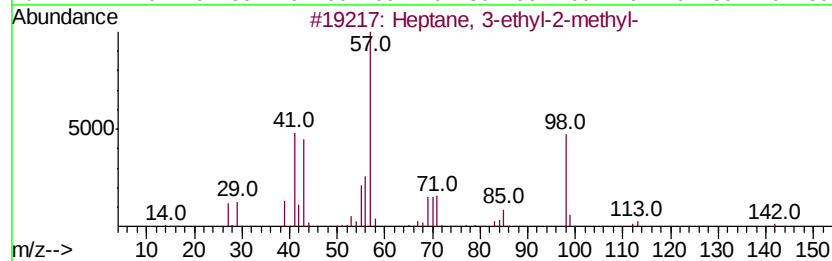
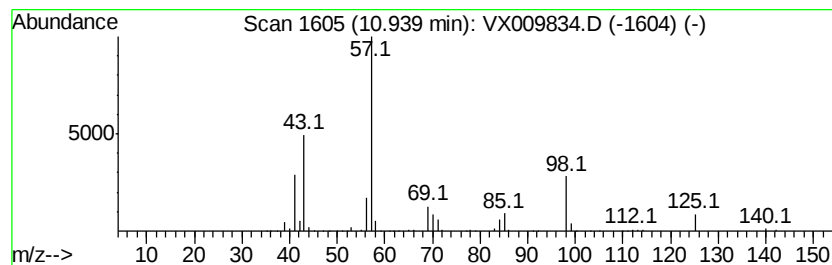
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Quant Title : SW846 8260

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

Peak Number 4 Heptane, 3-ethyl-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.94	27.95 ug/l	467130	Chlorobenzene-d5	10.11

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	80
2		Hexane, 3-ethyl-2,5-dimethyl-	142	C10H22	052897-04-8	59
3		Heptane, 4-propyl-	142	C10H22	003178-29-8	56
4		Undecane, 6-methyl-	170	C12H26	017302-33-9	56
5		Ether, tert-butyl isopropylidene...	154	C10H18O	024524-56-9	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009834.D
Acq On : 23 May 2019 04:01
Operator : JC/SP
Sample : K2976-15 50X
Misc : 5.39g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 46 Sample Multiplier: 1

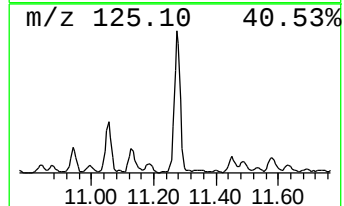
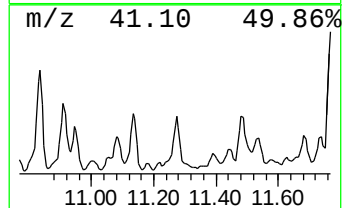
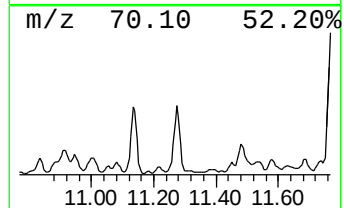
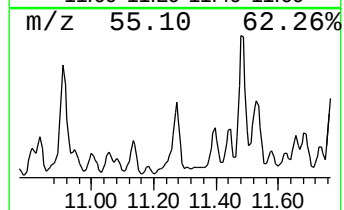
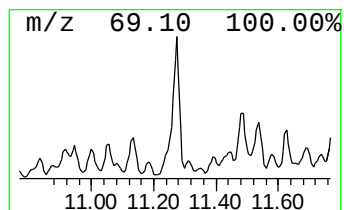
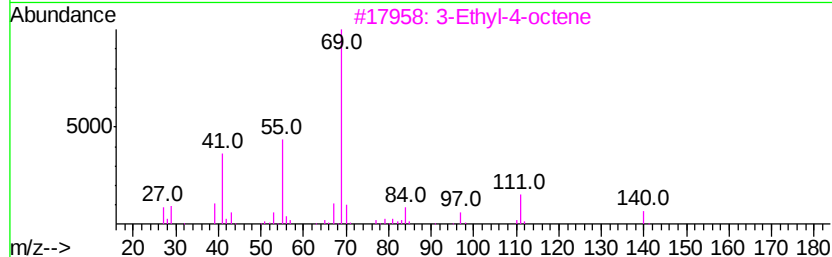
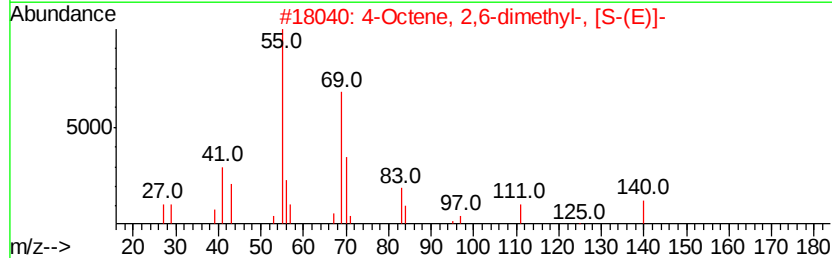
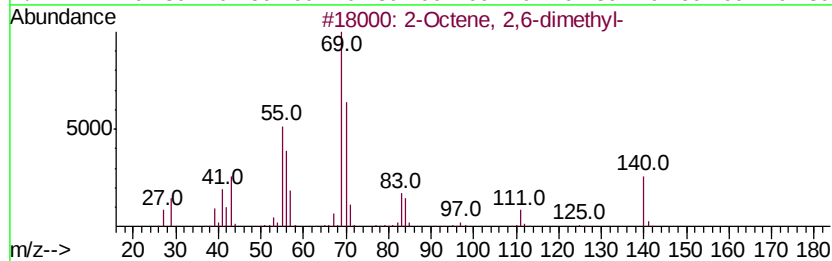
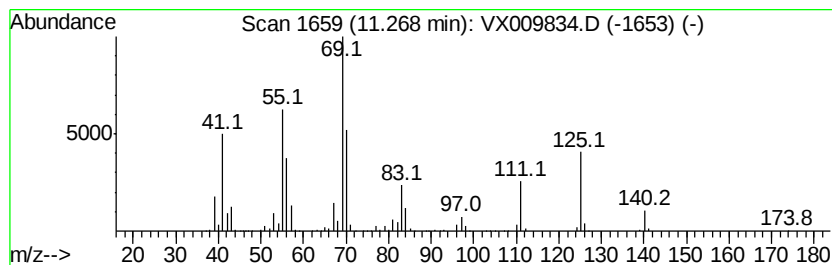
Instrument :
MSVOA_X
ClientSampleId :
EP1-SB-MW-IR(14-17)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

Peak Number 5 2-Octene, 2,6-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.27	32.08 ug/l	907439	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2-Octene, 2,6-dimethyl-	140 C10H20	004057-42-5	70
2	4-Octene, 2,6-dimethyl-, [S-(E)]-	140 C10H20	062960-76-3	59
3	3-Ethyl-4-octene	140 C10H20	019781-32-9	49
4	2-Octene, 2,7-dimethyl-	140 C10H20	002050-81-9	45
5	2-Methyl-2-octene	126 C9H18	016993-86-5	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009834.D
Acq On : 23 May 2019 04:01
Operator : JC/SP
Sample : K2976-15 50X
Misc : 5.39g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleID :
EP1-SB-MW-IR(14-17)

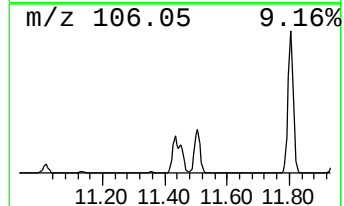
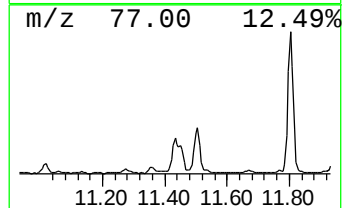
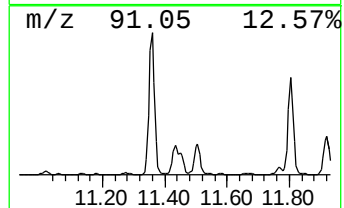
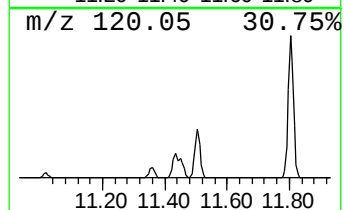
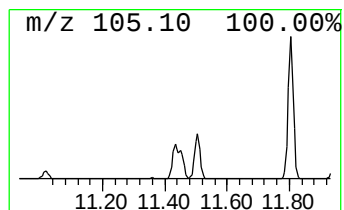
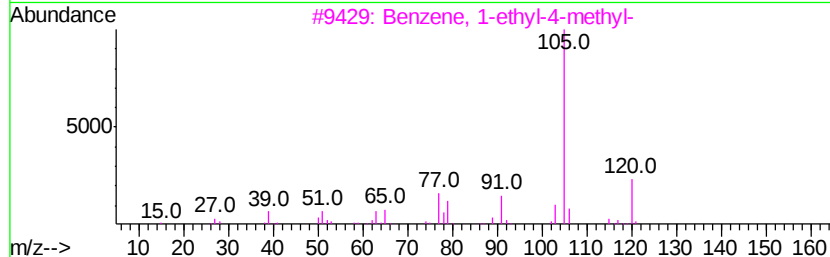
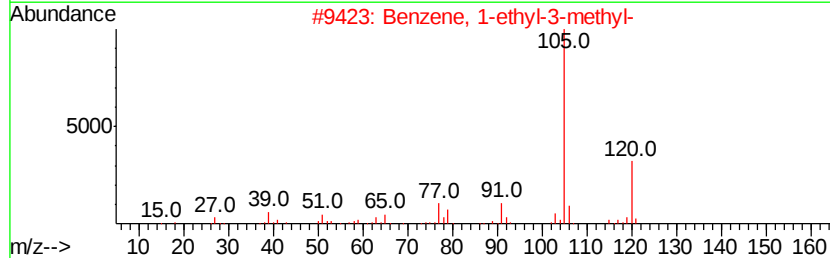
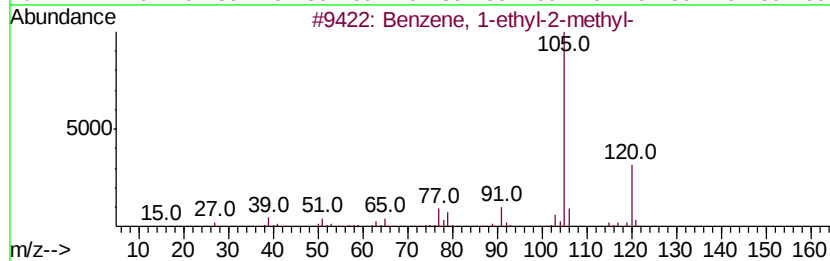
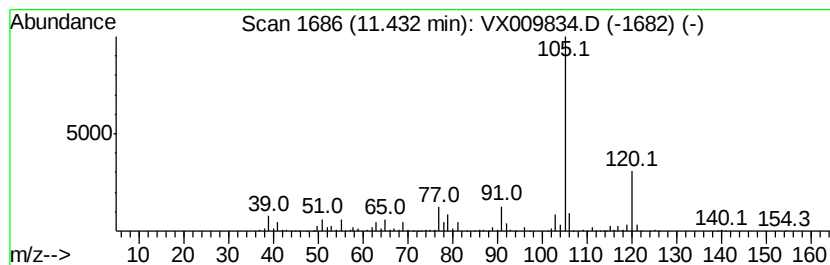
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.43	34.32 ug/l	970746	1,4-Dichlorobenzene-d4	12.08

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009834.D
Acq On : 23 May 2019 04:01
Operator : JC/SP
Sample : K2976-15 50X
Misc : 5.39g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 46 Sample Multiplier: 1

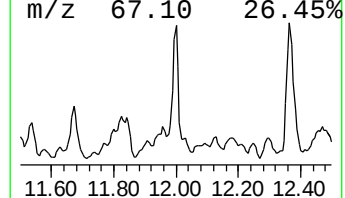
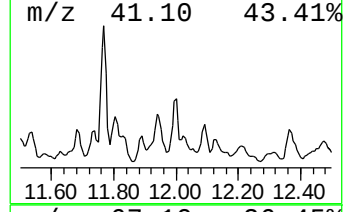
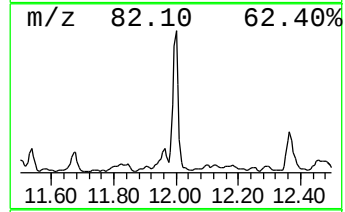
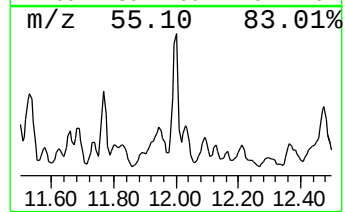
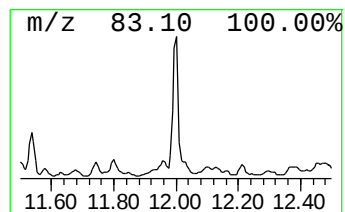
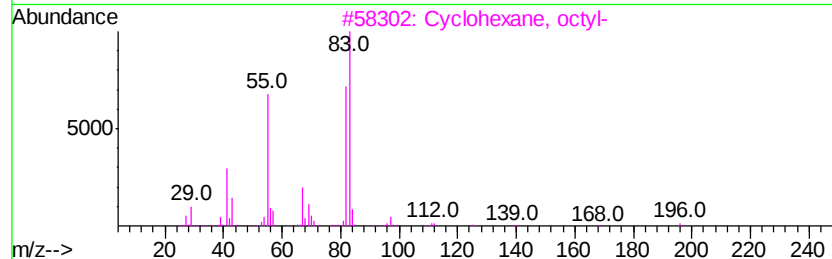
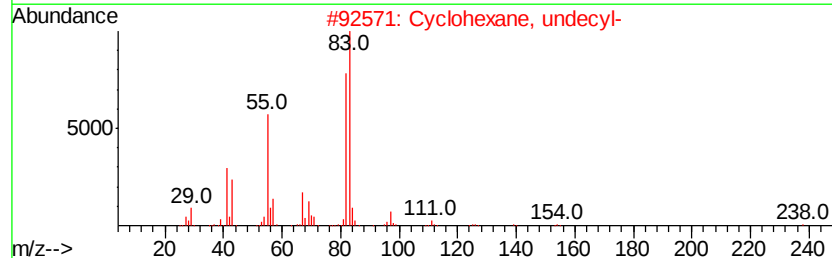
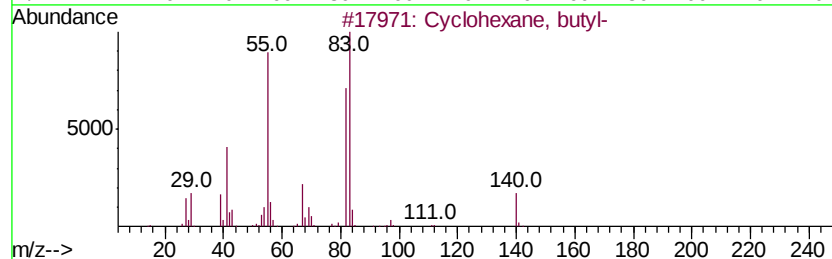
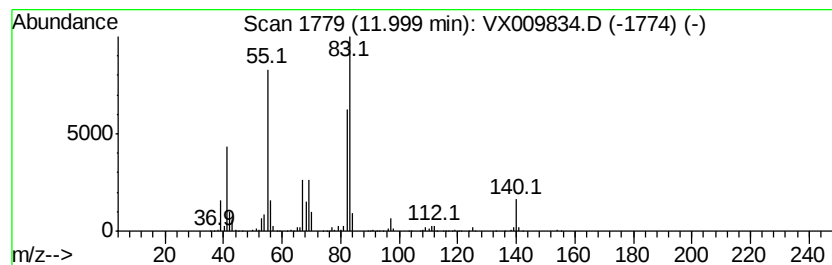
Instrument :
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ClientSampleId :
EP1-SB-MW-IR(14-17)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

Peak Number 7 Cyclohexane, butyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.00	27.39 ug/l	774587	1,4-Dichlorobenzene-d4	12.08	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cvclohexane, butyl-	140	C10H20	001678-93-9	91
2	Cyclohexane, undecyl-	238	C17H34	054105-66-7	78
3	Cvclohexane, octyl-	196	C14H28	001795-15-9	72
4	Cyclohexane, propyl-	126	C9H18	001678-92-8	72
5	Cyclohexane, pentyl-	154	C11H22	004292-92-6	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009834.D
Acq On : 23 May 2019 04:01
Operator : JC/SP
Sample : K2976-15 50X
Misc : 5.39g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 46 Sample Multiplier: 1

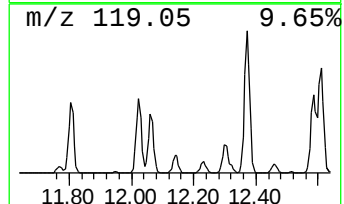
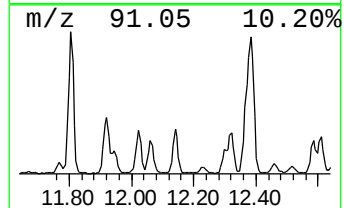
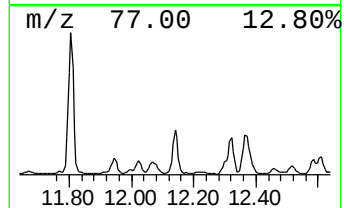
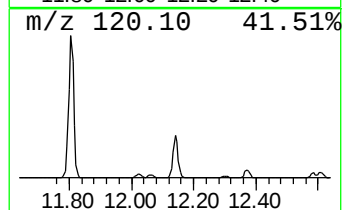
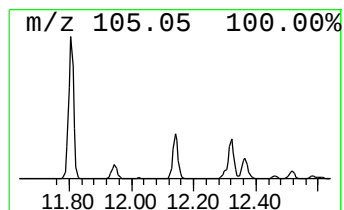
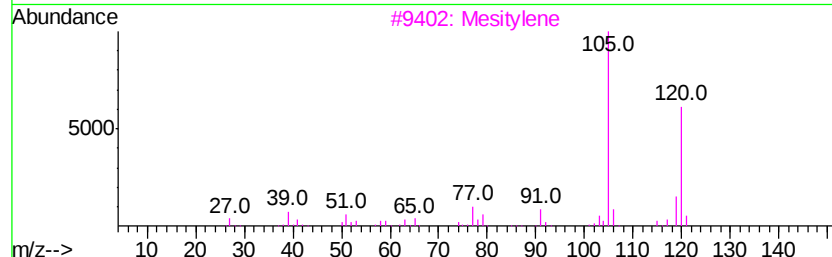
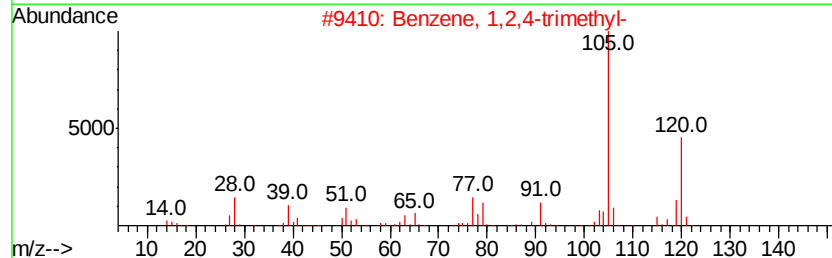
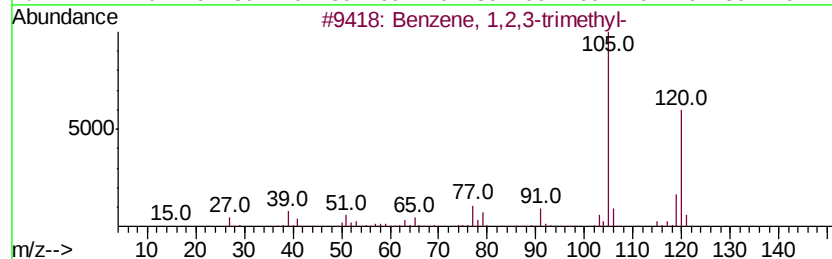
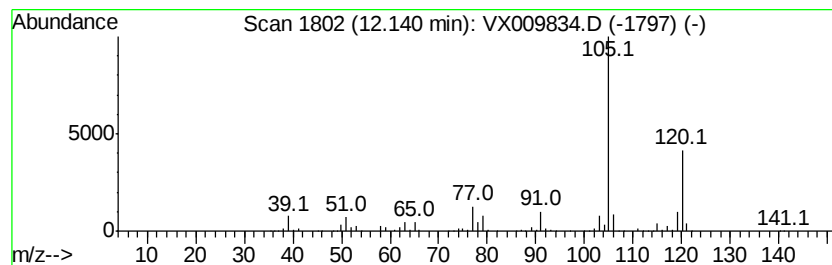
Instrument :
MSVOA_X
ClientSampleId :
EP1-SB-MW-IR(14-17)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	30.14 ug/l	852430	1,4-Dichlorobenzene-d4	12.08
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		Mesitylene	120 C9H12	000108-67-8 93
4		Benzene, 1-ethyl-3-methyl-	120 C9H12	000620-14-4 90
5		Benzene, 1-ethyl-2-methyl-	120 C9H12	000611-14-3 87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009834.D
Acq On : 23 May 2019 04:01
Operator : JC/SP
Sample : K2976-15 50X
Misc : 5.39g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 46 Sample Multiplier: 1

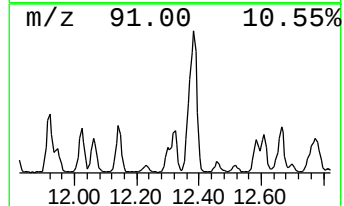
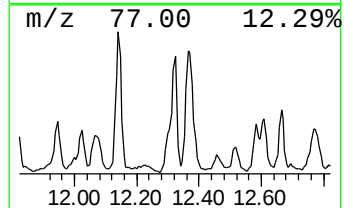
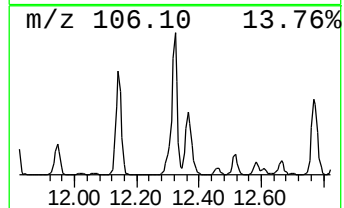
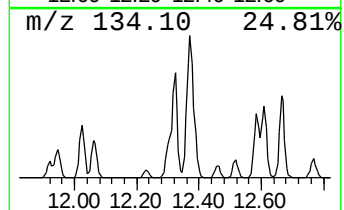
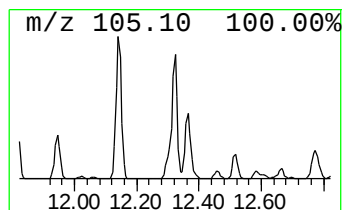
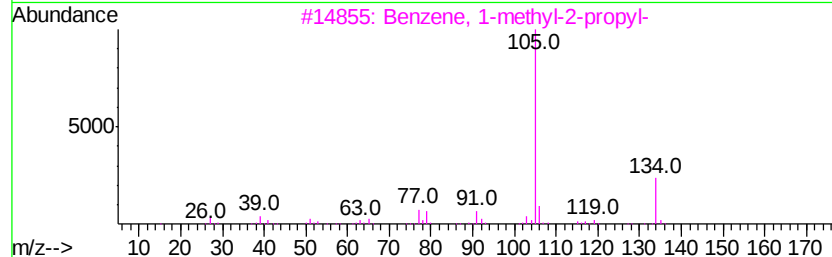
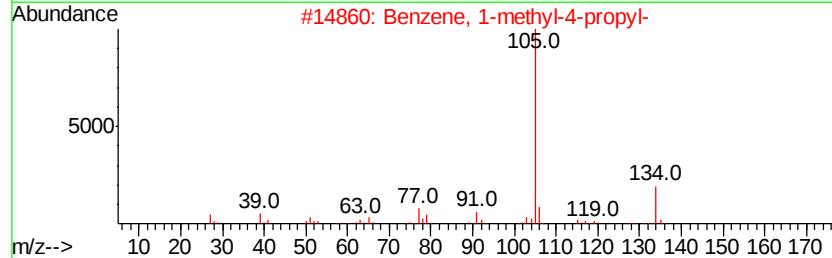
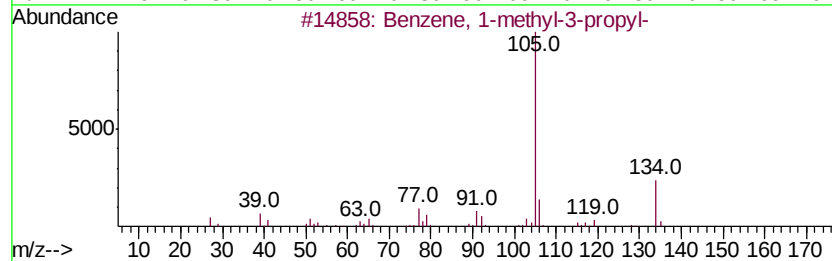
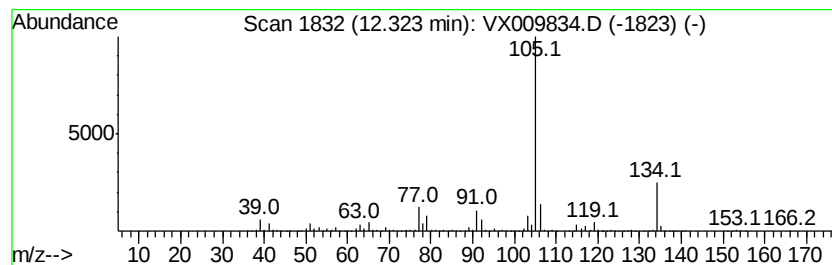
Instrument :
MSVOA_X
ClientSampleId :
EP1-SB-MW-IR(14-17)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-methyl-3-propyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.32	32.91 ug/l	930952	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-methyl-3-propyl-	134 C10H14	001074-43-7 95
2		Benzene, 1-methyl-4-propyl-	134 C10H14	001074-55-1 91
3		Benzene, 1-methyl-2-propyl-	134 C10H14	001074-17-5 91
4		Benzene, (1-methylpropyl)-	134 C10H14	000135-98-8 87
5		2-Tolyloxirane	134 C9H10O	002783-26-8 83



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009834.D
Acq On : 23 May 2019 04:01
Operator : JC/SP
Sample : K2976-15 50X
Misc : 5.39g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EP1-SB-MW-IR(14-17)

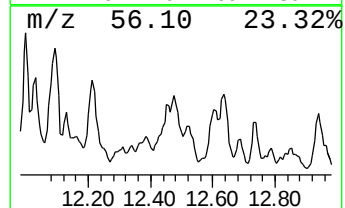
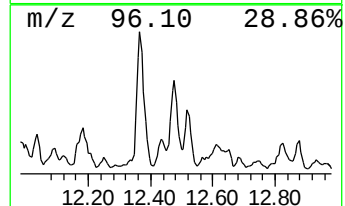
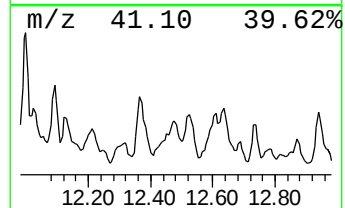
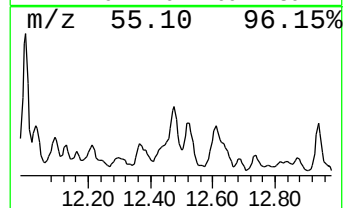
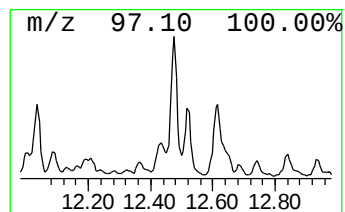
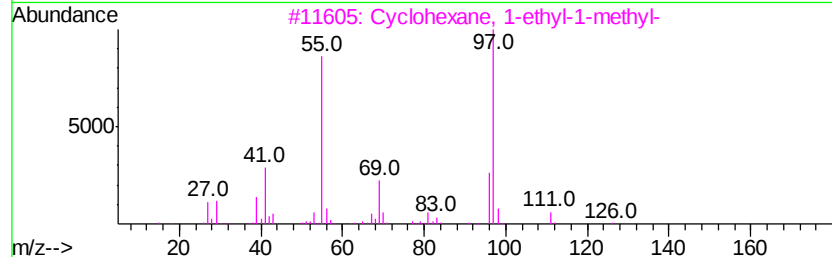
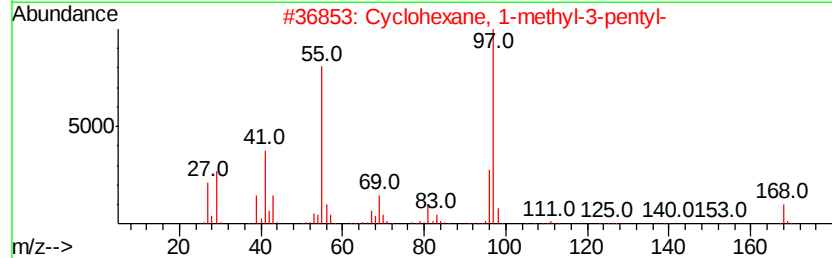
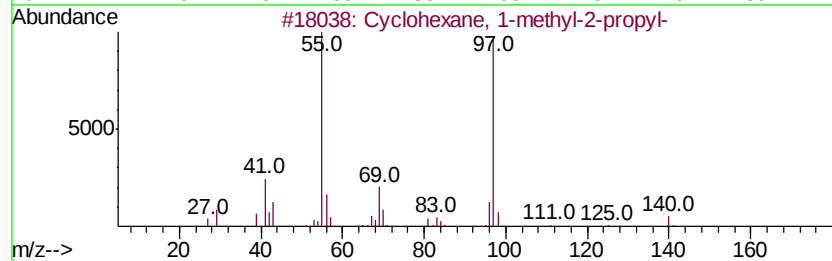
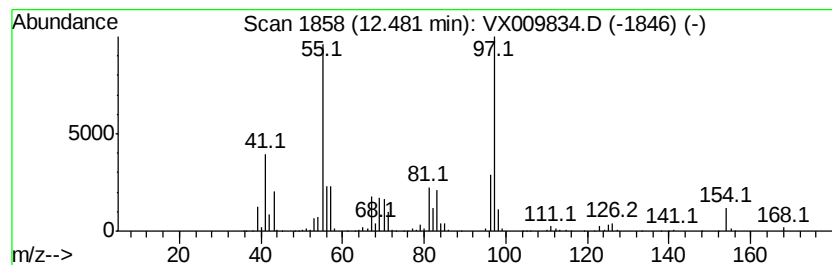
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

Peak Number 10 Cyclohexane, 1-methyl-2-pro... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.48	21.66 ug/l	612577	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1-methyl-2-propyl-	140	C10H20	004291-79-6	64
2		Cyclohexane, 1-methyl-3-pentyl-	168	C12H24	054411-02-8	64
3		Cyclohexane, 1-ethyl-1-methyl-	126	C9H18	004926-90-3	59
4		Cyclopentane, 1-hydroxymethyl-1,...	128	C8H16O	1000156-73-8	53
5		Cyclohexane, 1,3-dimethyl-	112	C8H16	000591-21-9	53



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\WX052219\
 Data File : VX009834.D
 Acq On : 23 May 2019 04:01
 Operator : JC/SP
 Sample : K2976-15 50X
 Misc : 5.39g/5mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 EP1-SB-MW-IR(14-17)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.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
Cyclohexane, 1-et...	10.64	33.0	ug/l	551318	3	10.11	835618	50.0
Octane, 2,6-dimet...	10.84	79.7	ug/l	1331160	3	10.11	835618	50.0
Cyclohexane, (1-m...	10.91	61.1	ug/l	1020450	3	10.11	835618	50.0
Heptane, 3-ethyl...	10.94	27.9	ug/l	467130	3	10.11	835618	50.0
2-Octene, 2,6-dim...	11.27	32.1	ug/l	907439	4	12.08	1414220	50.0
Benzene, 1-ethyl...	11.43	34.3	ug/l	970746	4	12.08	1414220	50.0
Cyclohexane, butyl-	12.00	27.4	ug/l	774587	4	12.08	1414220	50.0
Benzene, 1,2,3-tr...	12.14	30.1	ug/l	852430	4	12.08	1414220	50.0
Benzene, 1-methyl...	12.32	32.9	ug/l	930952	4	12.08	1414220	50.0
Cyclohexane, 1-me...	12.48	21.7	ug/l	612577	4	12.08	1414220	50.0