

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX052219\
 Data File : VX009832.D
 Acq On : 23 May 2019 03:14
 Operator : JC/SP
 Sample : K2976-13 50X
 Misc : 6.02g/5mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 EP1-SB-MW-HR(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	700	713	726	rBV	102454	293662	27.75%	1.775%
2	5.659	728	739	754	rVV	179673	515505	48.71%	3.115%
3	6.062	794	805	825	rBV	130793	352231	33.28%	2.129%
4	6.860	923	936	952	rBV	302645	717192	67.77%	4.334%
5	8.714	1228	1240	1257	rBV	628110	1005978	95.06%	6.079%
6	9.445	1353	1360	1371	rBV2	6435	11149	1.05%	0.067%
7	9.561	1374	1379	1385	rBV4	6735	13871	1.31%	0.084%
8	9.677	1392	1398	1403	rBV	18920	31806	3.01%	0.192%
9	9.890	1426	1433	1440	rBV4	18974	50400	4.76%	0.305%
10	9.957	1440	1444	1449	rVB	14013	20511	1.94%	0.124%
11	10.061	1455	1461	1464	rBV	32228	46183	4.36%	0.279%
12	10.110	1464	1469	1477	rVV	610069	829233	78.36%	5.011%
13	10.238	1485	1490	1496	rVB3	15977	28814	2.72%	0.174%
14	10.335	1496	1506	1508	rBV2	33762	65539	6.19%	0.396%
15	10.372	1508	1512	1516	rVV2	50855	75747	7.16%	0.458%
16	10.408	1516	1518	1529	rVB2	33715	62185	5.88%	0.376%
17	10.512	1529	1535	1541	rBV2	31486	49510	4.68%	0.299%
18	10.585	1543	1547	1549	rBV2	16064	20563	1.94%	0.124%
19	10.640	1549	1556	1564	rBV4	131362	231085	21.84%	1.396%
20	10.725	1564	1570	1574	rBV3	80115	130591	12.34%	0.789%
21	10.762	1574	1576	1580	rVB	42680	51125	4.83%	0.309%
22	10.835	1580	1588	1592	rBV	379918	547632	51.75%	3.309%
23	10.908	1592	1600	1603	rBV3	207986	364876	34.48%	2.205%
24	10.945	1603	1606	1611	rVB	200377	277631	26.23%	1.678%
25	11.000	1611	1615	1620	rVB3	55823	91516	8.65%	0.553%
26	11.055	1620	1624	1625	rBV	77215	88597	8.37%	0.535%
27	11.079	1625	1628	1633	rVV2	139947	203027	19.18%	1.227%
28	11.134	1633	1637	1641	rVB	766209	930964	87.97%	5.626%
29	11.176	1641	1644	1648	rVB2	66310	89183	8.43%	0.539%
30	11.219	1648	1651	1653	rBV	42714	53445	5.05%	0.323%
31	11.274	1653	1660	1670	rVB3	271722	444646	42.02%	2.687%
32	11.353	1670	1673	1675	rBV3	19728	21247	2.01%	0.128%
33	11.390	1675	1679	1684	rBV4	80908	145690	13.77%	0.880%
34	11.445	1684	1688	1691	rVV3	102226	149210	14.10%	0.902%

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35	11.481	1691	1694	1698	rVV	233183	307282	29.04%	1.857%
36	11.536	1698	1703	1707	rVB5	136993	224937	21.26%	1.359%
37	11.579	1707	1710	1715	rVB5	42125	66053	6.24%	0.399%
38	11.628	1715	1718	1720	rBV	32589	38672	3.65%	0.234%
39	11.682	1720	1727	1731	rBV4	163379	264789	25.02%	1.600%
40	11.737	1731	1736	1737	rBV	151722	184691	17.45%	1.116%
41	11.768	1737	1741	1744	rVV	705500	852561	80.56%	5.152%
42	11.804	1744	1747	1757	rVB	612410	1043339	98.59%	6.305%
43	11.890	1757	1761	1763	rBV	147748	203140	19.20%	1.228%
44	11.939	1763	1769	1774	rVV4	238617	430997	40.73%	2.605%
45	11.993	1774	1778	1782	rVV2	241537	353665	33.42%	2.137%
46	12.030	1782	1784	1786	rVV3	70167	74981	7.09%	0.453%
47	12.079	1786	1792	1797	rVB3	686323	1058262	100.00%	6.395%
48	12.140	1797	1802	1805	rBV	130241	181428	17.14%	1.096%
49	12.298	1823	1828	1833	rBV4	137928	228959	21.64%	1.384%
50	12.365	1834	1839	1846	rVV2	304194	535082	50.56%	3.234%
51	12.469	1846	1856	1861	rVV3	104833	299930	28.34%	1.813%
52	12.524	1861	1865	1870	rVB4	110018	213085	20.14%	1.288%
53	12.585	1870	1875	1876	rBV	165068	194813	18.41%	1.177%
54	12.609	1876	1879	1882	rVV2	187067	282201	26.67%	1.705%
55	12.664	1885	1888	1891	rVB	124165	134589	12.72%	0.813%
56	12.737	1896	1900	1911	rVB7	91680	244342	23.09%	1.477%
57	12.816	1911	1913	1914	rBV2	19205	16375	1.55%	0.099%
58	12.871	1919	1922	1925	rBV3	52894	60191	5.69%	0.364%
59	12.938	1930	1933	1936	rBV3	79861	114841	10.85%	0.694%
60	12.975	1936	1939	1942	rVB2	83374	99082	9.36%	0.599%
61	13.018	1942	1946	1948	rBV	109044	137551	13.00%	0.831%
62	13.079	1953	1956	1963	rVB4	49094	88031	8.32%	0.532%
63	13.146	1963	1967	1972	rBV2	29900	46592	4.40%	0.282%
64	13.194	1972	1975	1978	rVB2	24122	25234	2.38%	0.152%
65	13.261	1983	1986	1989	rVV3	25715	30579	2.89%	0.185%
66	13.304	1989	1993	1995	rVV2	35184	43722	4.13%	0.264%
67	13.341	1995	1999	2006	rVB2	121198	203898	19.27%	1.232%
68	13.432	2010	2014	2017	rVV2	28142	46883	4.43%	0.283%
69	13.475	2017	2021	2030	rVB	65999	113316	10.71%	0.685%
70	13.560	2031	2035	2038	rBV3	11184	16674	1.58%	0.101%
71	13.633	2043	2047	2052	rBV3	19032	33362	3.15%	0.202%

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Integration Parameters: RTEINT.P

Integrator: RTE
Smoothing : ON Filtering: 5
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Stop Thrs : 0 Peak Location: TOP

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72	13.719	2059	2061	2066	rVB2	11327	15527	1.47%	0.094%
73	13.804	2070	2075	2077	rBV4	6667	10734	1.01%	0.065%
74	13.896	2086	2090	2097	rVB6	10790	16633	1.57%	0.101%

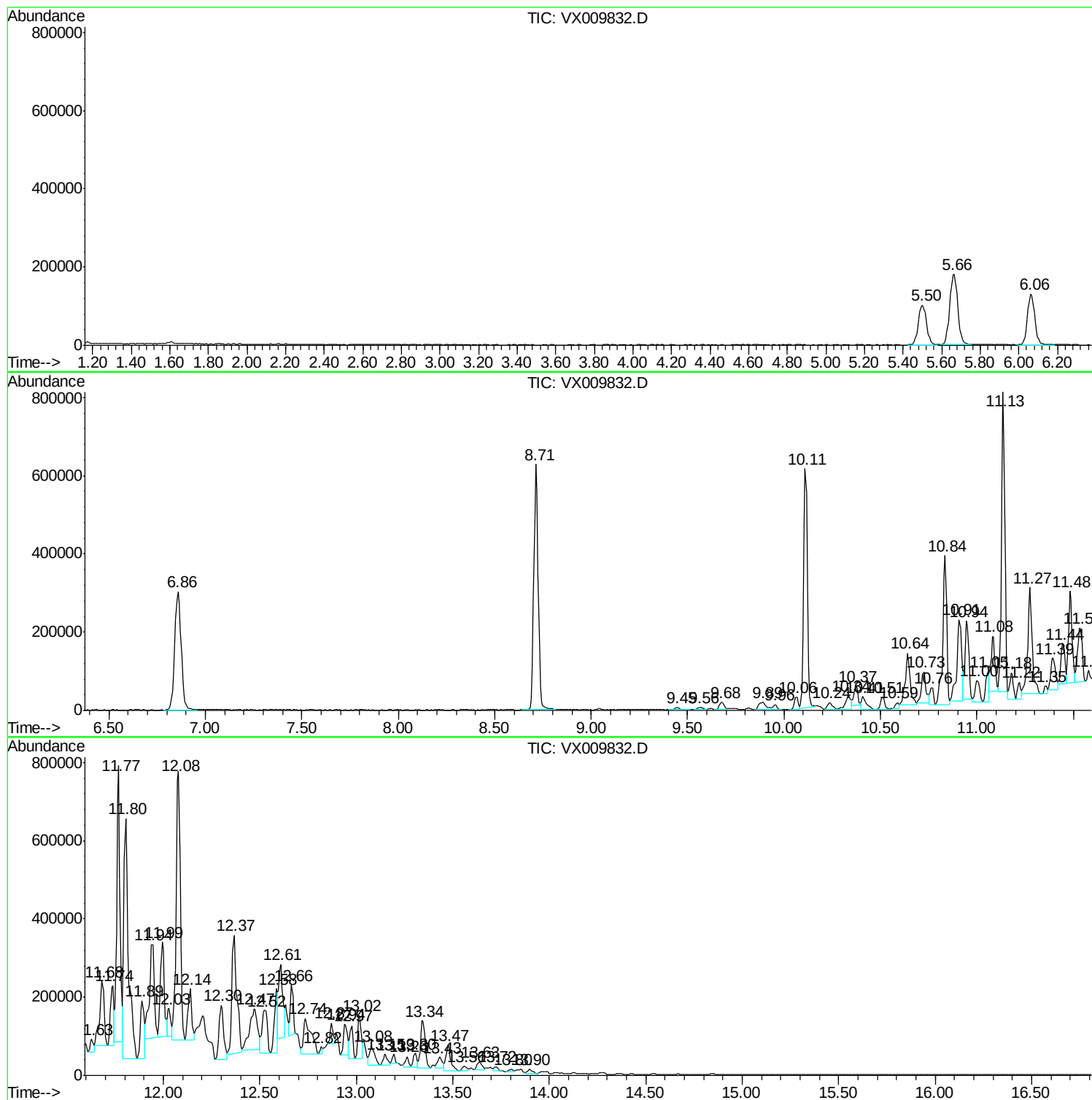
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Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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



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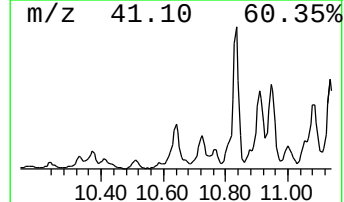
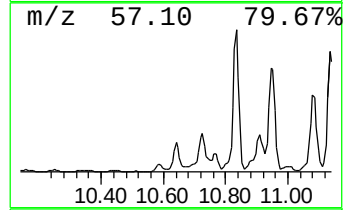
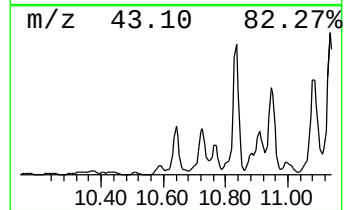
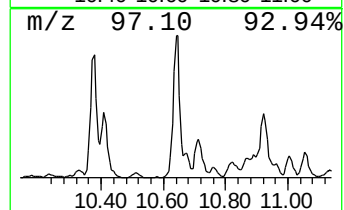
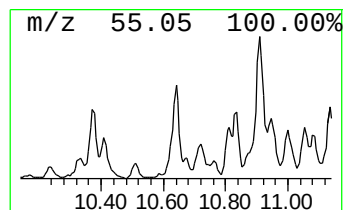
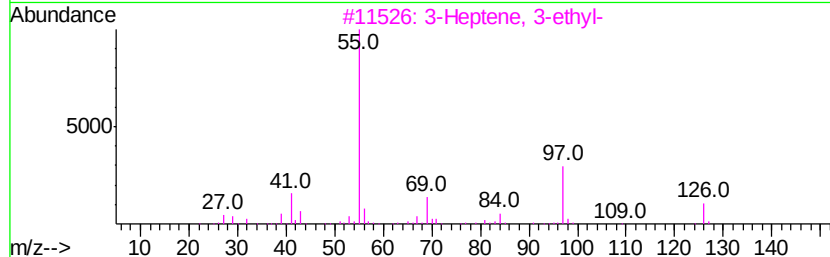
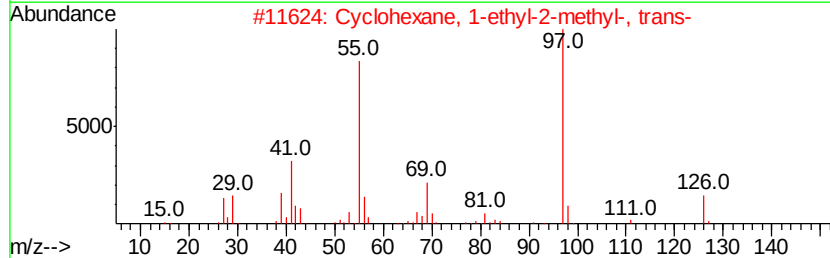
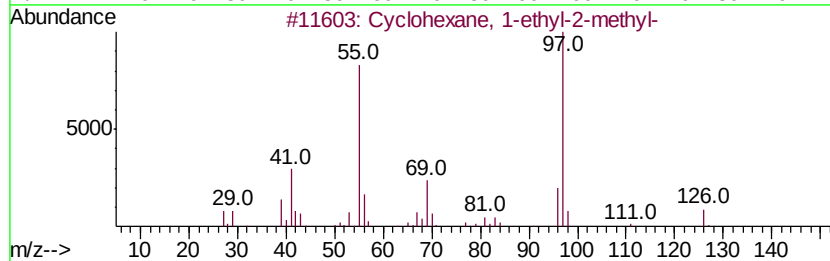
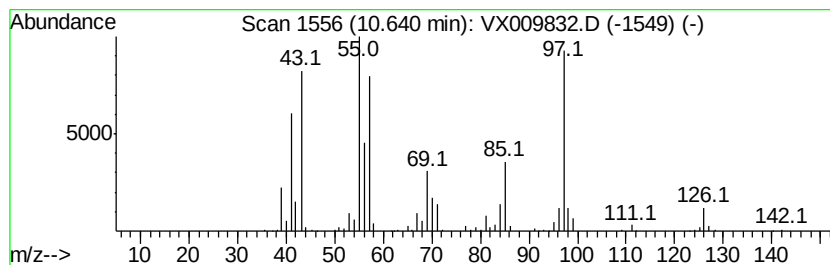
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

Peak Number 1 unknown10.64 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.64	13.93 ug/l	231085	Chlorobenzene-d5	10.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1-ethyl-2-methyl-	126	C9H18	003728-54-9	49
2		Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	49
3		3-Heptene, 3-ethyl-	126	C9H18	074764-46-8	49
4		Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	46
5		Thiophene, 2-propyl-	126	C7H10S	001551-27-5	38



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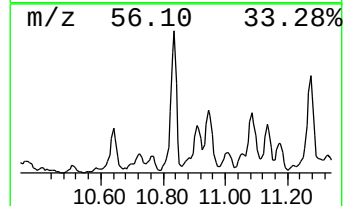
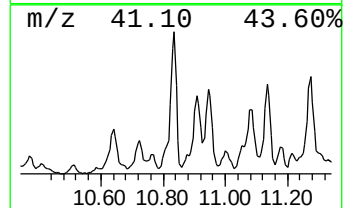
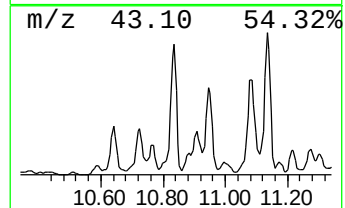
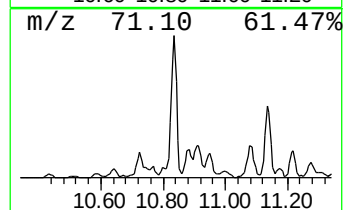
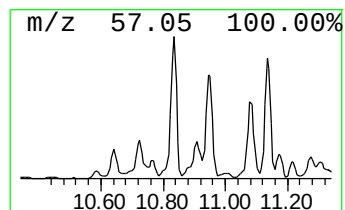
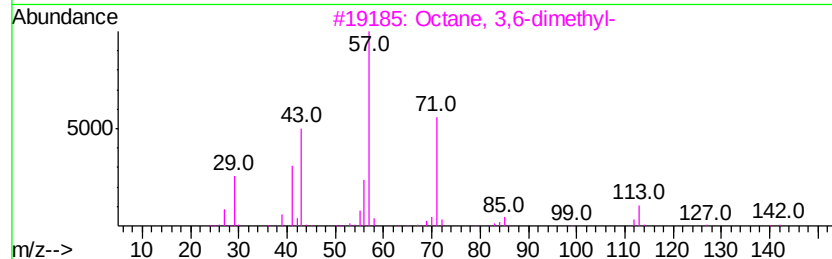
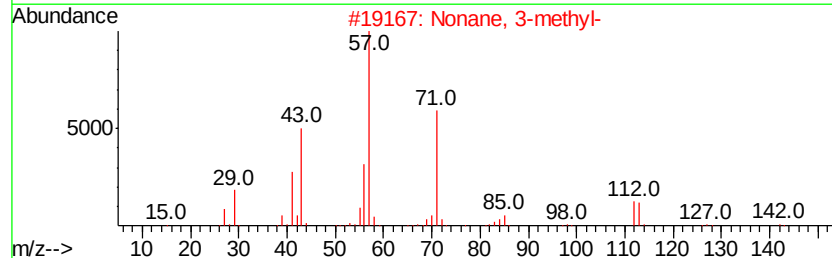
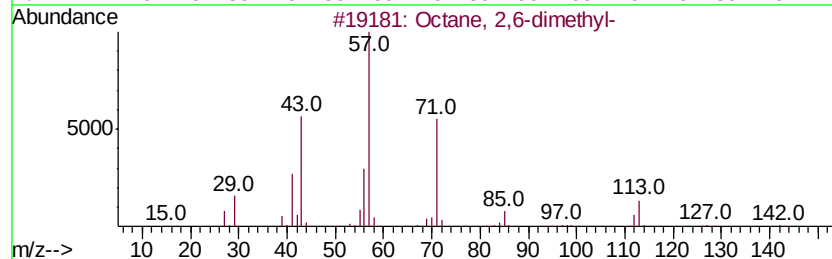
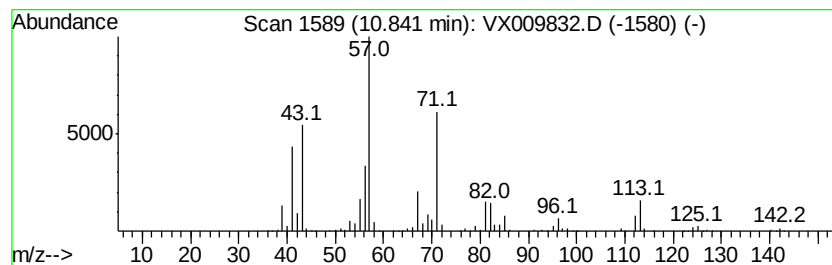
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 Octane, 2,6-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.84	33.02 ug/l	547632	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 76
3		Octane, 3,6-dimethyl-	142 C10H22	015869-94-0 76
4		Undecane, 6,6-dimethyl-	184 C13H28	017312-76-4 53
5		Sulfurous acid, di(2-ethylhexyl)...	306 C16H34O3S	1000309-19-1 50



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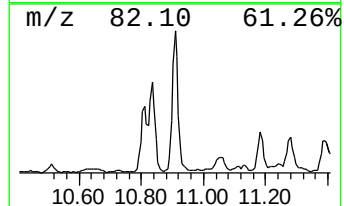
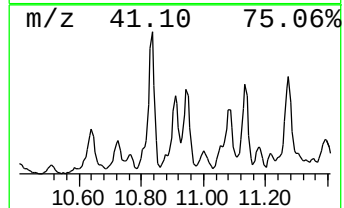
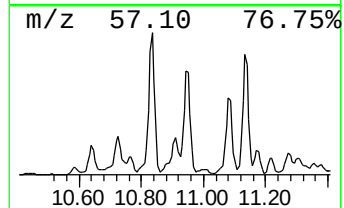
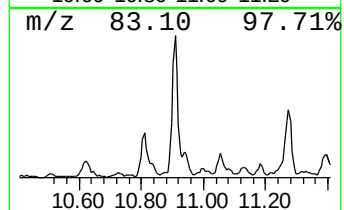
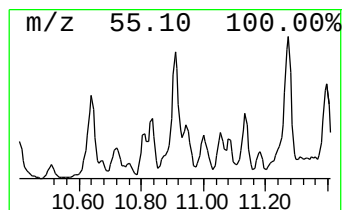
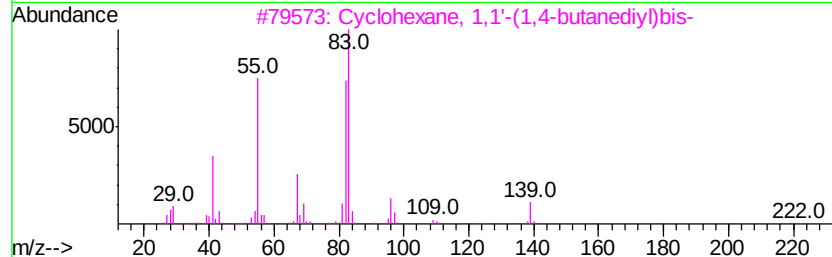
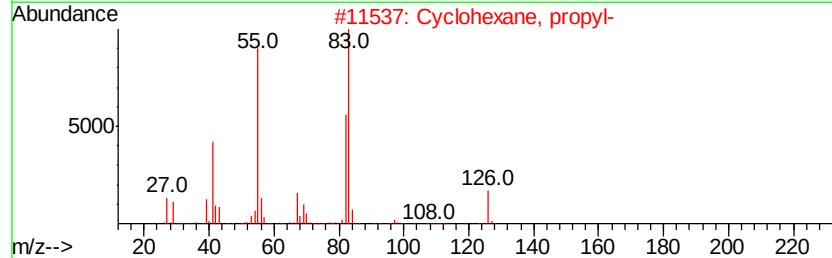
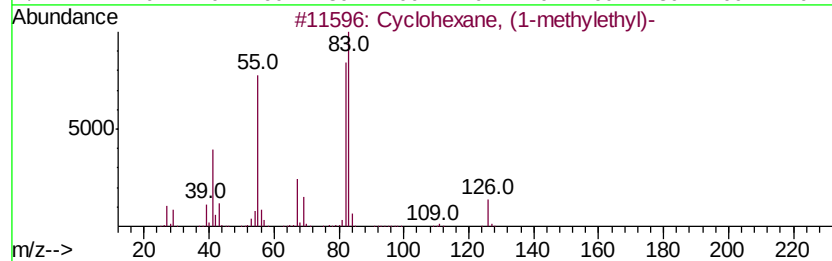
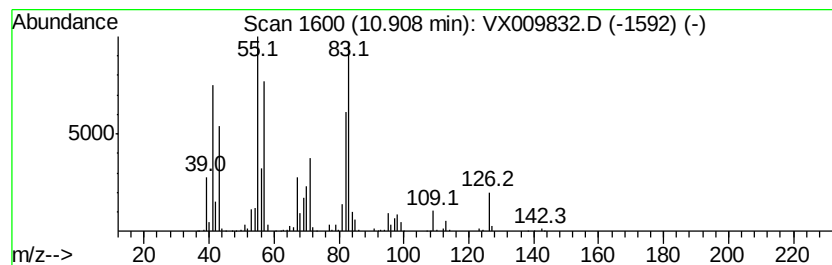
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.91	22.00 ug/l	364876	Chlorobenzene-d5	10.11		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cvclohexane, (1-methylethvl)-	126	C9H18	000696-29-7	64
2		Cvclohexane, propyl-	126	C9H18	001678-92-8	52
3		Cvclohexane, 1,1'-(1,4-butanediv...	222	C16H30	006165-44-2	47
4		trans-7-Oxabicyclo[4.3.0]nonane	126	C8H14O	027345-70-6	43
5		Ethanone, 1-cyclohexyl-	126	C8H14O	000823-76-7	38



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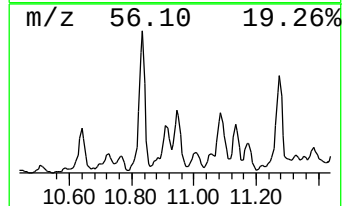
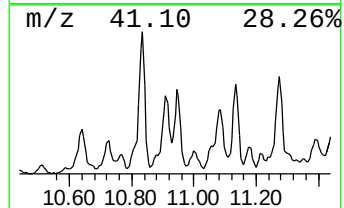
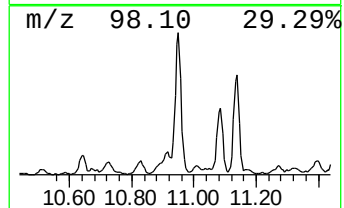
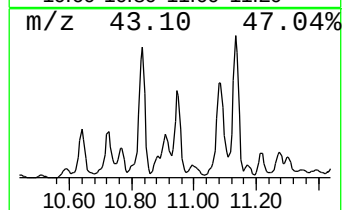
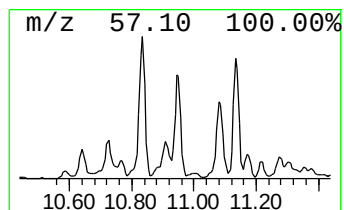
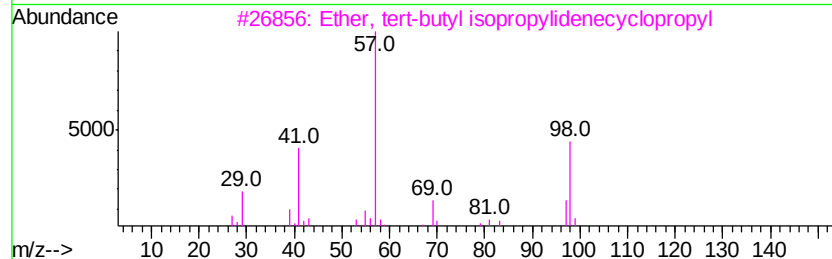
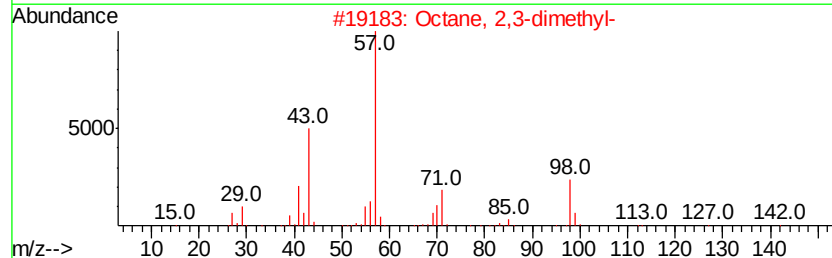
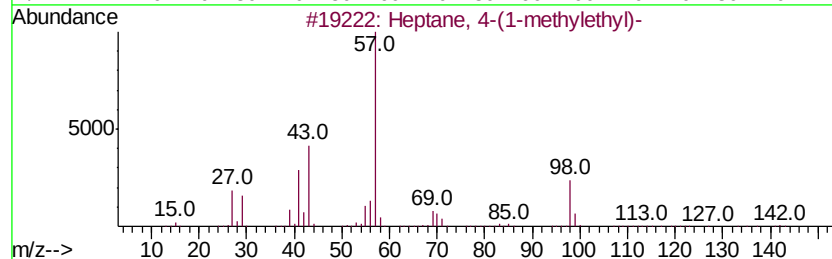
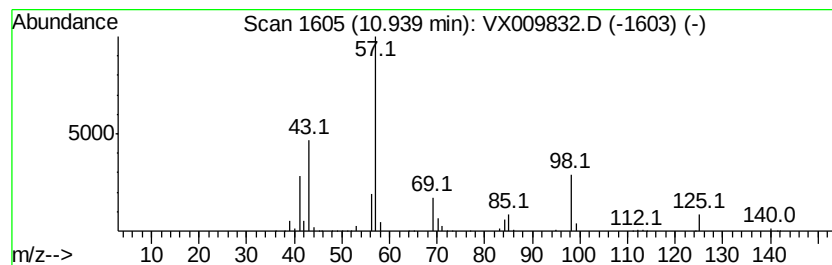
Instrument :
MSVOA_X
ClientSampleId :
EP1-SB-MW-HR(14-17)

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 4 Heptane, 4-(1-methylethyl)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.94	16.74 ug/l	277631	Chlorobenzene-d5	10.11		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptane, 4-(1-methylethyl)-	142	C10H22	052896-87-4	64
2		Octane, 2,3-dimethyl-	142	C10H22	007146-60-3	53
3		Ether, tert-butyl isopropylidene...	154	C10H18O	024524-56-9	50
4		Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	47
5		Undecane, 5,6-dimethyl-	184	C13H28	017615-91-7	37



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009832.D
Acq On : 23 May 2019 03:14
Operator : JC/SP
Sample : K2976-13 50X
Misc : 6.02g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 44 Sample Multiplier: 1

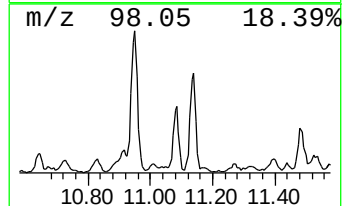
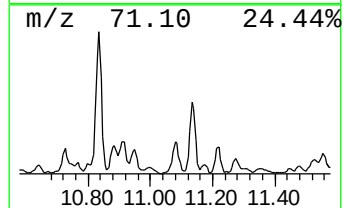
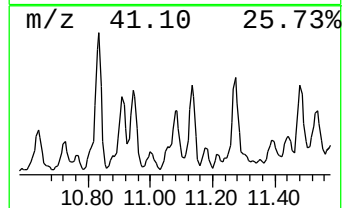
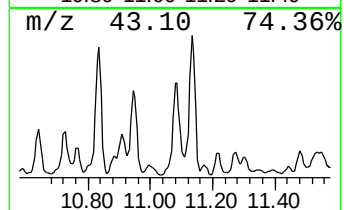
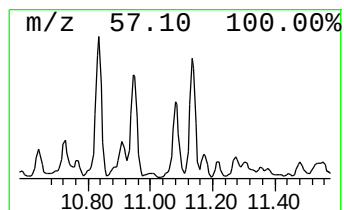
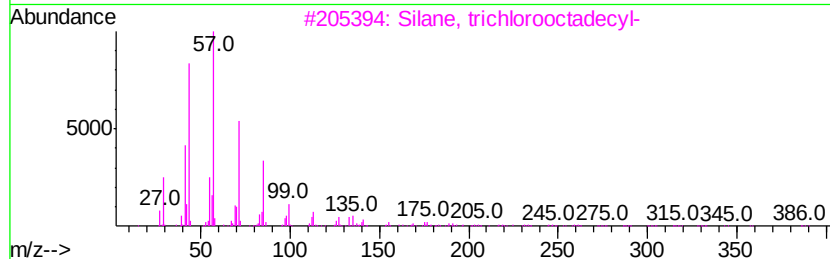
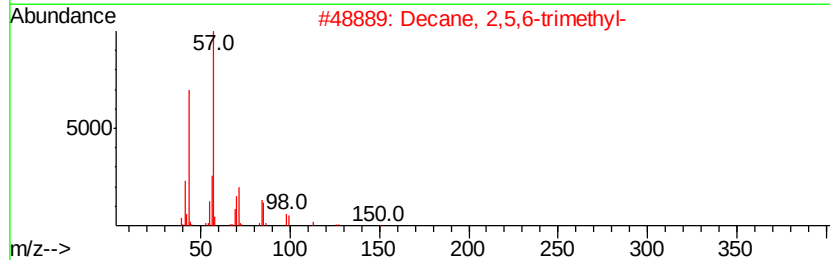
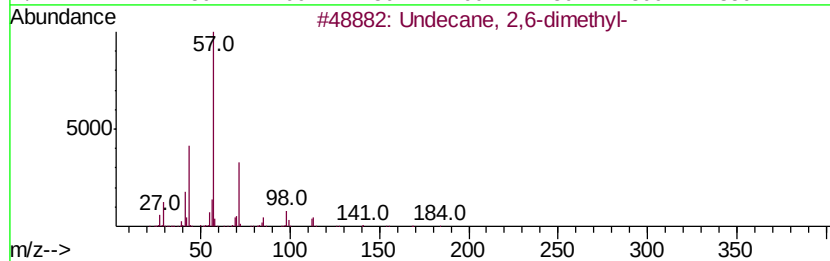
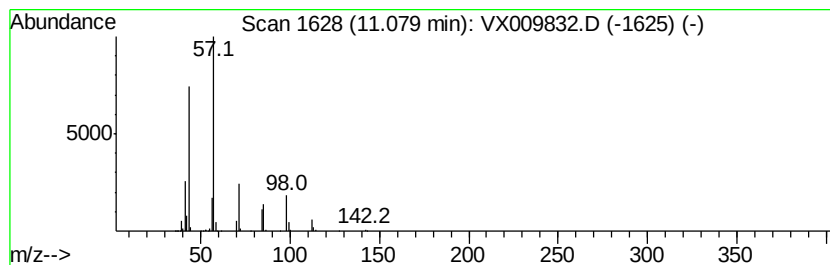
Instrument :
MSVOA_X
ClientSampleID :
EP1-SB-MW-HR(14-17)

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 Undecane, 2,6-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.08	12.24 ug/l	203027	Chlorobenzene-d5	10.11
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Undecane, 2,6-dimethyl-	184 C13H28	017301-23-4	64
2	Decane, 2,5,6-trimethyl-	184 C13H28	062108-23-0	53
3	Silane, trichlorooctadecyl-	386 C18H37Cl3Si	000112-04-9	53
4	Heptane, 3-ethyl-2-methyl-	142 C10H22	014676-29-0	50
5	Octane, 2,3-dimethyl-	142 C10H22	007146-60-3	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009832.D
Acq On : 23 May 2019 03:14
Operator : JC/SP
Sample : K2976-13 50X
Misc : 6.02µ/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 44 Sample Multiplier: 1

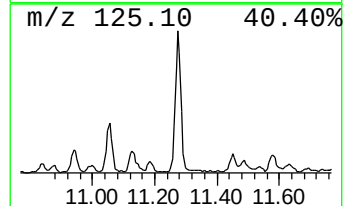
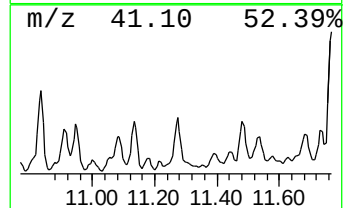
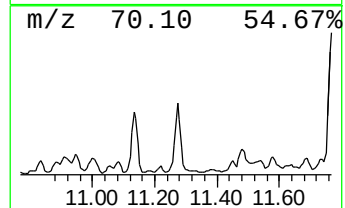
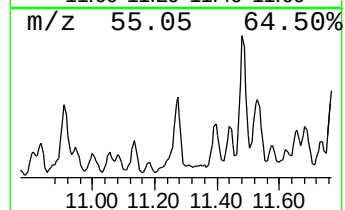
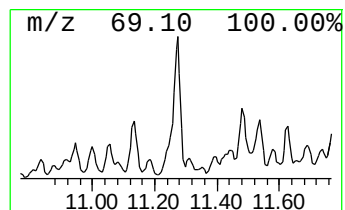
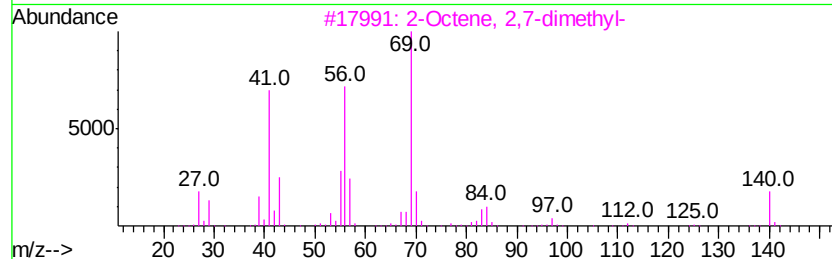
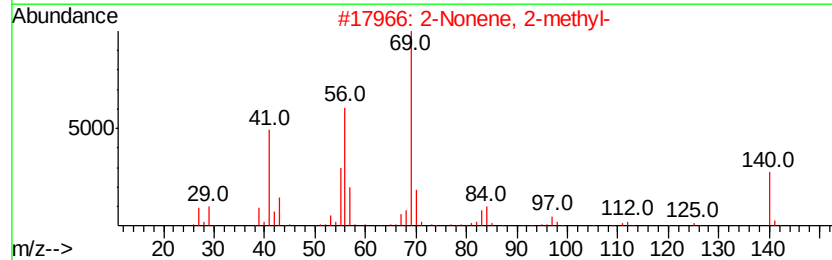
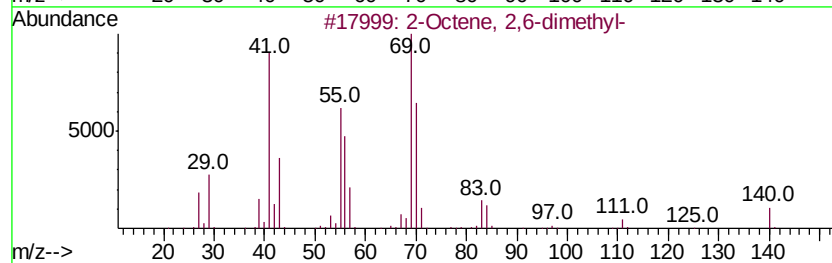
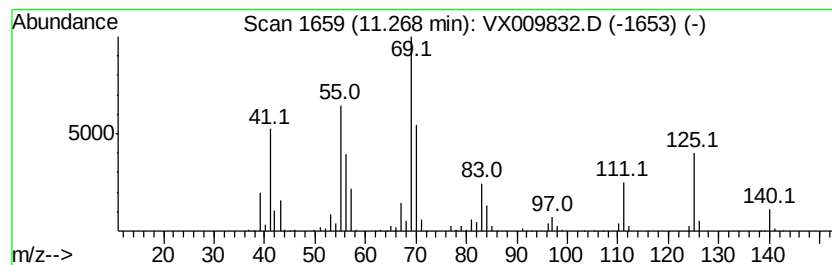
Instrument :
MSVOA_X
ClientSampleId :
EP1-SB-MW-HR(14-17)

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 6 2-Octene, 2,6-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.27	21.01 ug/l	444646	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	62
2	2-Nonene, 2-methyl-	140	C10H20	002129-95-5	49
3	2-Octene, 2,7-dimethyl-	140	C10H20	002050-81-9	46
4	3-Ethyl-4-octene	140	C10H20	019781-32-9	45
5	3-Nonene, 2-methyl-	140	C10H20	053966-53-3	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009832.D
Acq On : 23 May 2019 03:14
Operator : JC/SP
Sample : K2976-13 50X
Misc : 6.02g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 44 Sample Multiplier: 1

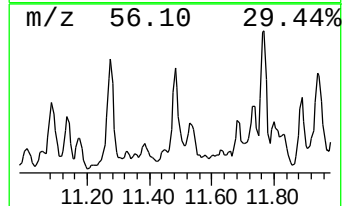
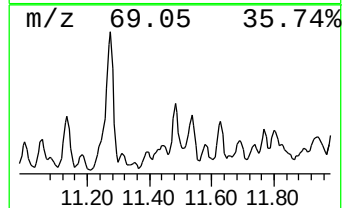
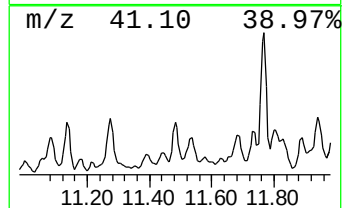
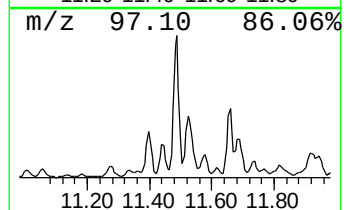
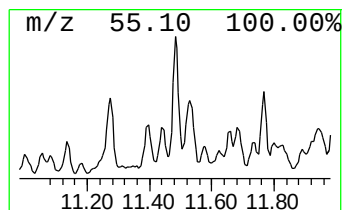
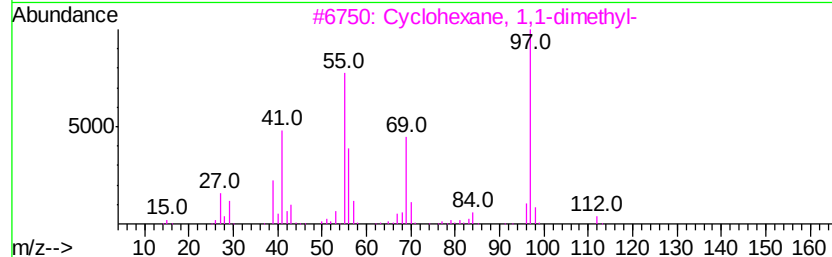
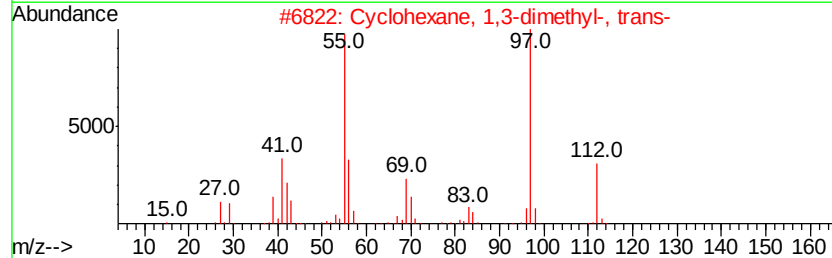
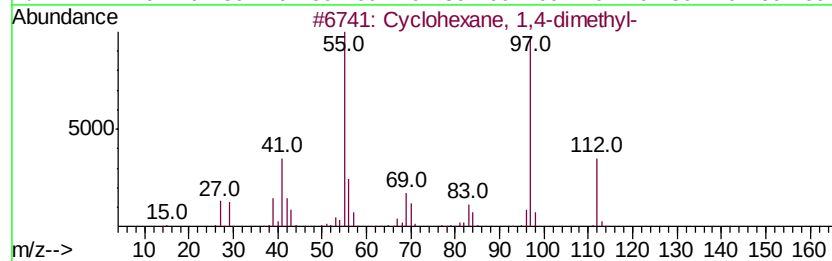
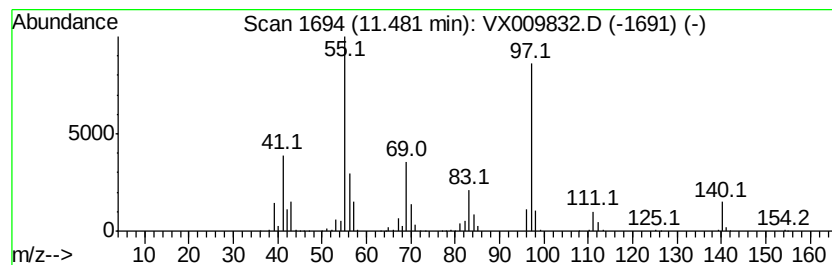
Instrument :
MSVOA_X
ClientSampleId :
EP1-SB-MW-HR(14-17)

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, 1,4-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.48	14.52 ug/l	307282	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclohexane, 1,4-dimethyl-	112 C8H16	000589-90-2 87
2		Cyclohexane, 1,3-dimethyl-, trans-	112 C8H16	002207-03-6 74
3		Cyclohexane, 1,1-dimethyl-	112 C8H16	000590-66-9 70
4		Cyclohexane, 1,3-dimethyl-, cis-	112 C8H16	000638-04-0 64
5		3-Hexene, 3-ethyl-2,5-dimethyl-	140 C10H20	062338-08-3 64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009832.D
Acq On : 23 May 2019 03:14
Operator : JC/SP
Sample : K2976-13 50X
Misc : 6.02g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 44 Sample Multiplier: 1

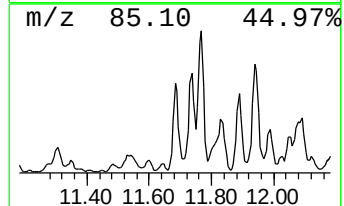
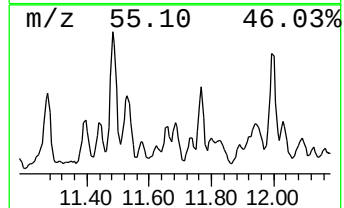
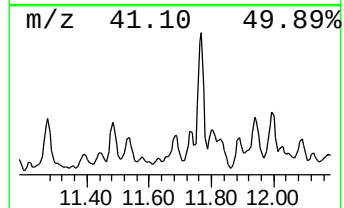
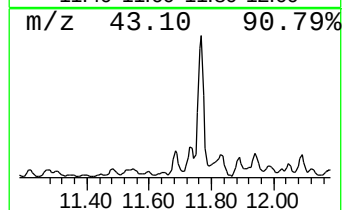
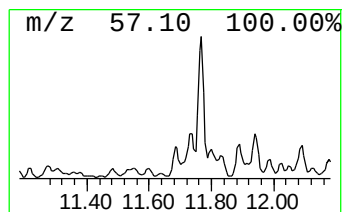
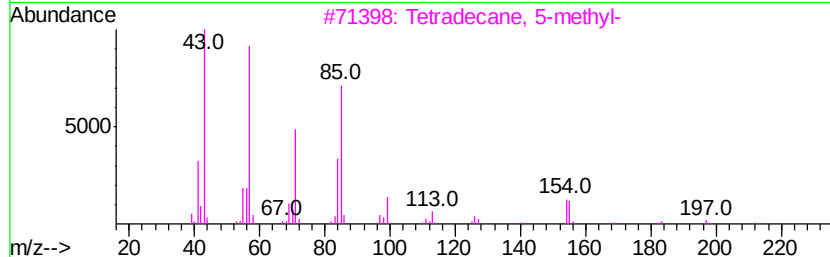
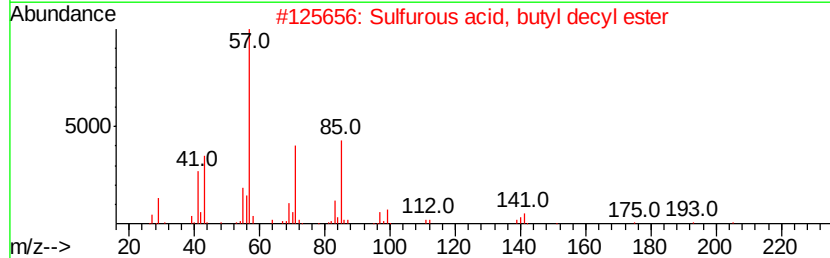
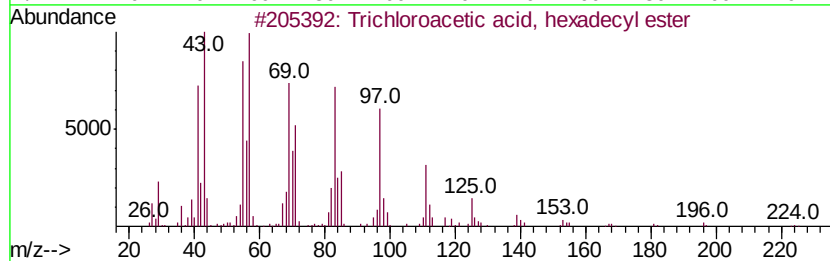
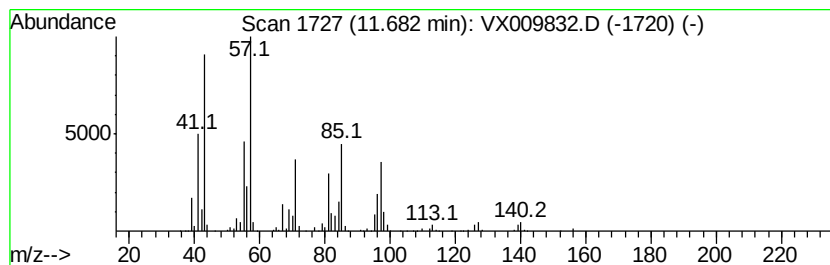
Instrument :
MSVOA_X
ClientSampleId :
EP1-SB-MW-HR(14-17)

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 unknown11.68 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.68	12.51 ug/l	264789	1,4-Dichlorobenzene-d4	12.08		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	47
2		Sulfurous acid, butyl decyl ester	278	C14H30O3S	1000309-17-7	47
3		Tetradecane, 5-methyl-	212	C15H32	025117-32-2	43
4		2,4-Dimethyldodecane	198	C14H30	006117-99-3	43
5		Tridecane, 5-methyl-	198	C14H30	025117-31-1	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009832.D
Acq On : 23 May 2019 03:14
Operator : JC/SP
Sample : K2976-13 50X
Misc : 6.02g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 44 Sample Multiplier: 1

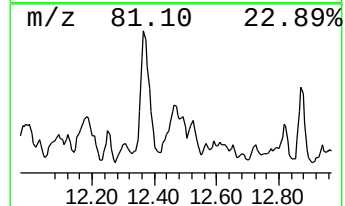
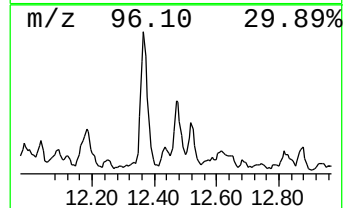
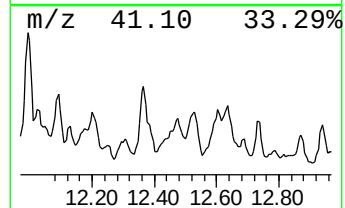
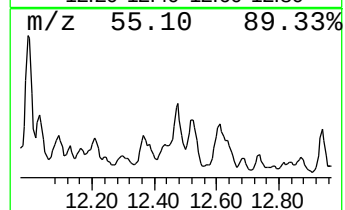
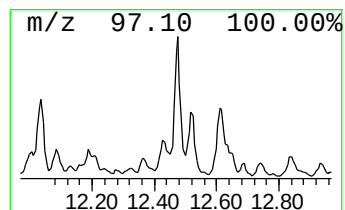
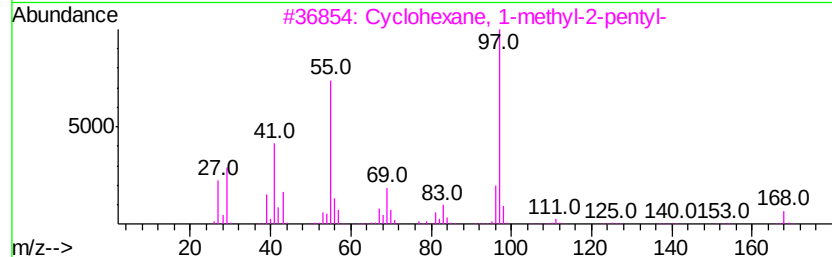
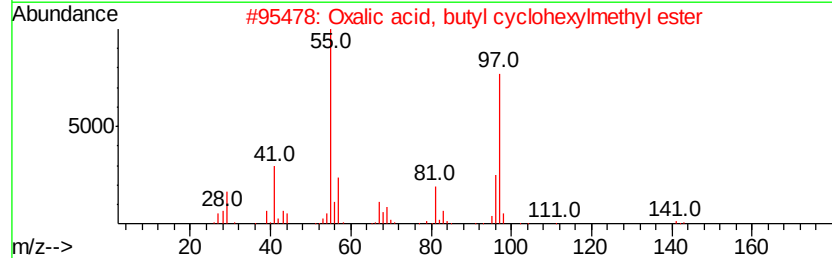
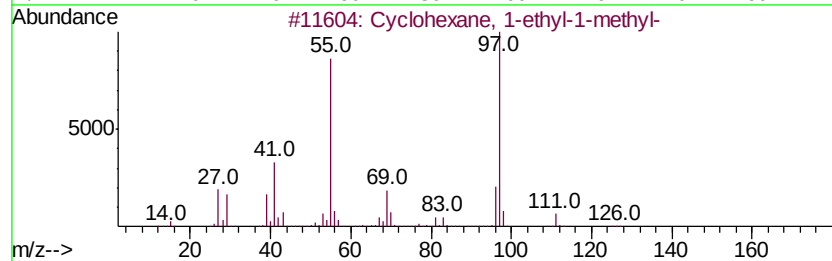
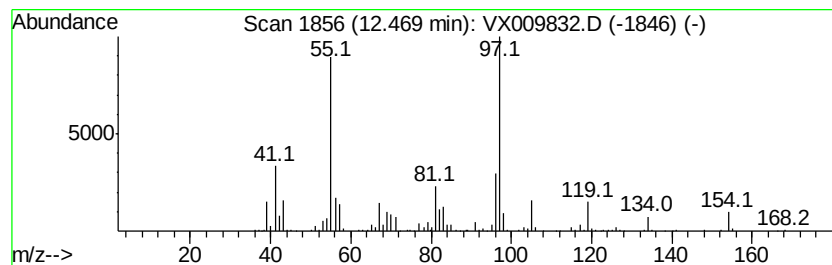
Instrument :
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ClientSampleId :
EP1-SB-MW-HR(14-17)

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 Cyclohexane, 1-ethyl-1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.47	14.17 ug/l	299930	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclohexane, 1-ethyl-1-methyl-	126 C9H18	004926-90-3 59
2		Oxalic acid, butyl cyclohexylmet...	242 C13H22O4	1000309-68-2 59
3		Cyclohexane, 1-methyl-2-pentyl-	168 C12H24	054411-01-7 59
4		Oxalic acid, cyclohexylmethyl pr...	228 C12H20O4	1000309-68-1 59
5		Cyclohexane, 1-isopropyl-1-methyl-	140 C10H20	016580-26-0 59



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009832.D
Acq On : 23 May 2019 03:14
Operator : JC/SP
Sample : K2976-13 50X
Misc : 6.02g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 44 Sample Multiplier: 1

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

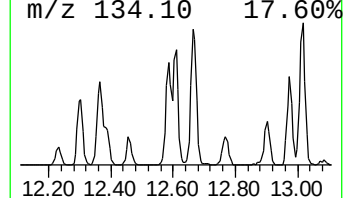
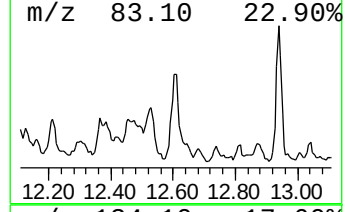
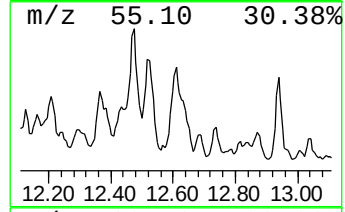
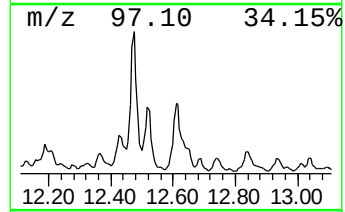
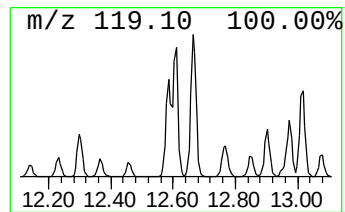
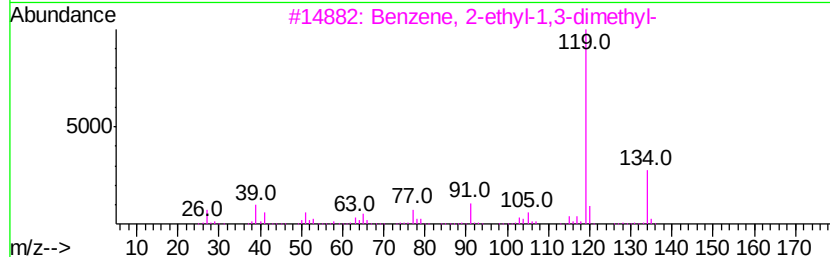
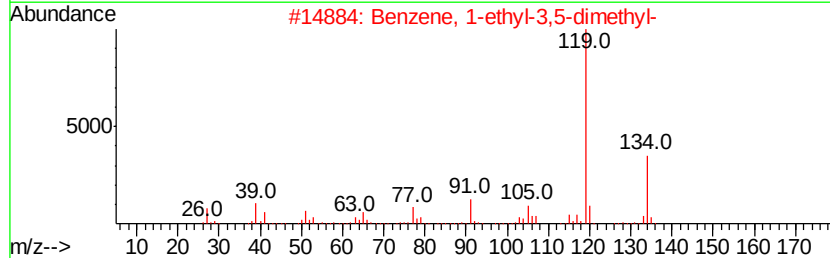
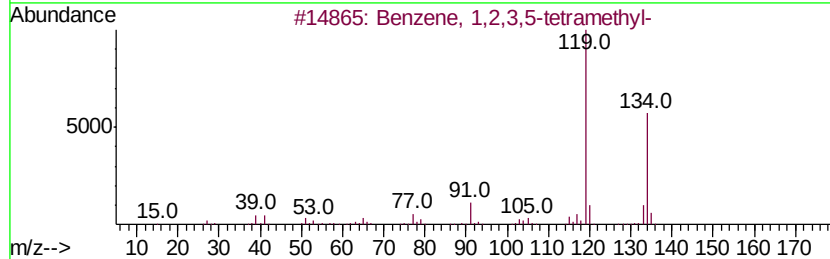
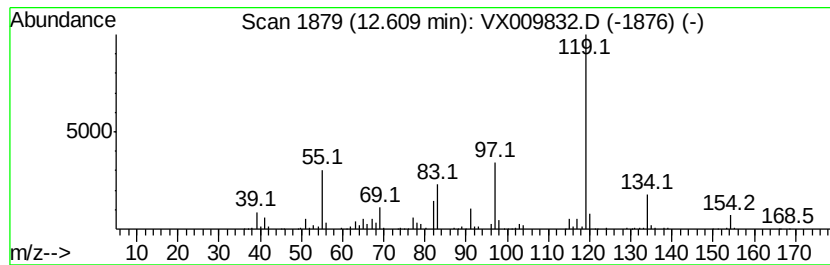
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.61	13.33 ug/l	282201	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	64
2		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	64
3		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	58
4		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	53
5		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	52



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX052219\
Data File : VX009832.D
Acq On : 23 May 2019 03:14
Operator : JC/SP
Sample : K2976-13 50X
Misc : 6.02g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EP1-SB-MW-HR(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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown10.64	10.64	13.9	ug/l	231085	3	10.11	829233	50.0
Octane, 2,6-dimet...	10.84	33.0	ug/l	547632	3	10.11	829233	50.0
Cyclohexane, (1-m...	10.91	22.0	ug/l	364876	3	10.11	829233	50.0
Heptane, 4-(1-met...	10.94	16.7	ug/l	277631	3	10.11	829233	50.0
Undecane, 2,6-dim...	11.08	12.2	ug/l	203027	3	10.11	829233	50.0
2-Octene, 2,6-dim...	11.27	21.0	ug/l	444646	4	12.08	1058260	50.0
Cyclohexane, 1,4-...	11.48	14.5	ug/l	307282	4	12.08	1058260	50.0
unknown11.68	11.68	12.5	ug/l	264789	4	12.08	1058260	50.0
Cyclohexane, 1-et...	12.47	14.2	ug/l	299930	4	12.08	1058260	50.0
Benzene, 1,2,3,5-...	12.61	13.3	ug/l	282201	4	12.08	1058260	50.0