

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX052219\
 Data File : VX009838.D
 Acq On : 23 May 2019 05:34
 Operator : JC/SP
 Sample : K2976-19 50X
 Misc : 6.08g/5mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 EP1-SB-SBR-3BR(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	5.500	702	713	726	rBV	102546	296434	6.97%	0.740%
2	5.659	729	739	760	rVB	184735	528437	12.42%	1.319%
3	6.061	792	805	823	rBV	132524	350465	8.24%	0.875%
4	6.860	925	936	951	rBV	302562	722480	16.98%	1.803%
5	8.713	1235	1240	1254	rVB	650940	1026005	24.11%	2.561%
6	9.042	1288	1294	1303	rVB	27080	47015	1.11%	0.117%
7	9.445	1350	1360	1369	rBV	67720	112572	2.65%	0.281%
8	9.560	1373	1379	1385	rBV4	55078	123429	2.90%	0.308%
9	9.621	1385	1389	1393	rVB	77658	106060	2.49%	0.265%
10	9.676	1393	1398	1402	rBV	143891	215031	5.05%	0.537%
11	9.890	1425	1433	1437	rVV4	118315	295962	6.96%	0.739%
12	9.957	1440	1444	1449	rVB	125900	170694	4.01%	0.426%
13	10.060	1455	1461	1465	rBV	240253	348293	8.19%	0.869%
14	10.109	1465	1469	1474	rVV	630300	860591	20.23%	2.148%
15	10.182	1474	1481	1485	rVB5	32073	78519	1.85%	0.196%
16	10.237	1485	1490	1497	rBV2	76240	141145	3.32%	0.352%
17	10.335	1497	1506	1508	rBV2	157455	302478	7.11%	0.755%
18	10.371	1508	1512	1516	rVV2	273012	450867	10.60%	1.125%
19	10.408	1516	1518	1530	rVB	183780	306658	7.21%	0.765%
20	10.512	1530	1535	1541	rBV	119100	179025	4.21%	0.447%
21	10.585	1541	1547	1549	rBV2	50833	67887	1.60%	0.169%
22	10.640	1549	1556	1564	rBV3	500010	874265	20.55%	2.182%
23	10.725	1564	1570	1573	rBV3	269776	417513	9.81%	1.042%
24	10.768	1574	1577	1580	rVB	140894	176128	4.14%	0.440%
25	10.835	1580	1588	1592	rBV	1266104	1832660	43.07%	4.574%
26	10.908	1592	1600	1604	rBV	914762	1564631	36.77%	3.905%
27	10.944	1604	1606	1611	rVB	522659	675282	15.87%	1.685%
28	11.011	1611	1617	1621	rBV3	197385	340383	8.00%	0.850%
29	11.054	1621	1624	1626	rBV2	197376	278200	6.54%	0.694%
30	11.085	1626	1629	1632	rVV	431168	575839	13.53%	1.437%
31	11.133	1632	1637	1641	rVB	1289813	1701614	39.99%	4.247%
32	11.182	1641	1645	1648	rVB2	177892	221990	5.22%	0.554%
33	11.219	1648	1651	1653	rBV	115539	147725	3.47%	0.369%
34	11.274	1653	1660	1669	rVB2	707105	1205027	28.32%	3.007%

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Integration Parameters: RTEINT.P

Integrator: RTE
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35	11.359	1670	1674	1676	rBV	261234	312906	7.35%	0.781%
36	11.395	1676	1680	1684	rVV3	183677	306895	7.21%	0.766%
37	11.450	1684	1689	1692	rVV2	499311	794252	18.67%	1.982%
38	11.481	1692	1694	1699	rVV	674203	917883	21.57%	2.291%
39	11.530	1699	1702	1707	rVB3	372542	593764	13.96%	1.482%
40	11.578	1707	1710	1715	rVB4	101632	142796	3.36%	0.356%
41	11.627	1715	1718	1720	rBV	96461	110270	2.59%	0.275%
42	11.682	1720	1727	1732	rVB4	374620	676788	15.91%	1.689%
43	11.737	1732	1736	1738	rBV2	322488	422404	9.93%	1.054%
44	11.767	1738	1741	1744	rVV	1375318	1588272	37.33%	3.964%
45	11.804	1744	1747	1757	rVB	3059621	4254684	100.00%	10.619%
46	11.889	1757	1761	1763	rBV	286527	382264	8.98%	0.954%
47	11.920	1763	1766	1767	rVV	215496	254978	5.99%	0.636%
48	11.944	1767	1770	1774	rVV2	566992	765091	17.98%	1.909%
49	11.999	1774	1779	1782	rVV2	726827	985343	23.16%	2.459%
50	12.030	1782	1784	1786	rVV2	153543	158583	3.73%	0.396%
51	12.078	1786	1792	1797	rVB3	774463	1425793	33.51%	3.558%
52	12.139	1797	1802	1806	rBV2	569595	824953	19.39%	2.059%
53	12.298	1823	1828	1833	rBV2	417686	603965	14.20%	1.507%
54	12.365	1834	1839	1846	rVV2	749862	1370089	32.20%	3.419%
55	12.468	1846	1856	1862	rVV3	247356	694465	16.32%	1.733%
56	12.523	1862	1865	1870	rVB3	258441	443948	10.43%	1.108%
57	12.584	1870	1875	1877	rBV	403509	589086	13.85%	1.470%
58	12.609	1877	1879	1885	rVV4	493332	708413	16.65%	1.768%
59	12.664	1885	1888	1891	rVB	389438	424368	9.97%	1.059%
60	12.767	1897	1905	1911	rVB4	219466	676990	15.91%	1.690%
61	12.846	1915	1918	1920	rBV3	67165	89228	2.10%	0.223%
62	12.871	1920	1922	1925	rVV2	88825	90844	2.14%	0.227%
63	12.944	1930	1934	1936	rBV2	250590	358485	8.43%	0.895%
64	12.974	1936	1939	1942	rVB	239067	290252	6.82%	0.724%
65	13.017	1942	1946	1949	rBV	324737	401992	9.45%	1.003%
66	13.078	1953	1956	1964	rVB4	140030	256363	6.03%	0.640%
67	13.145	1964	1967	1972	rBV2	78403	118185	2.78%	0.295%
68	13.194	1972	1975	1977	rBV	65981	84421	1.98%	0.211%
69	13.261	1983	1986	1989	rVB3	54292	61341	1.44%	0.153%
70	13.304	1989	1993	1996	rVV2	88697	117022	2.75%	0.292%
71	13.340	1996	1999	2006	rVB2	373400	576740	13.56%	1.439%

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ClientSampleId :
EP1-SB-SBR-3BR(14-17)

Integration Parameters: RTEINT.P

Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Title : SW846 8260

72	13.426	2010	2013	2017	rVV2	49993	74035	1.74%	0.185%
73	13.480	2017	2022	2030	rVB6	80945	180725	4.25%	0.451%
74	13.560	2031	2035	2038	rBV3	32150	47643	1.12%	0.119%
75	13.633	2044	2047	2053	rBV3	54224	93156	2.19%	0.232%
76	13.718	2059	2061	2070	rVB2	36078	56801	1.34%	0.142%

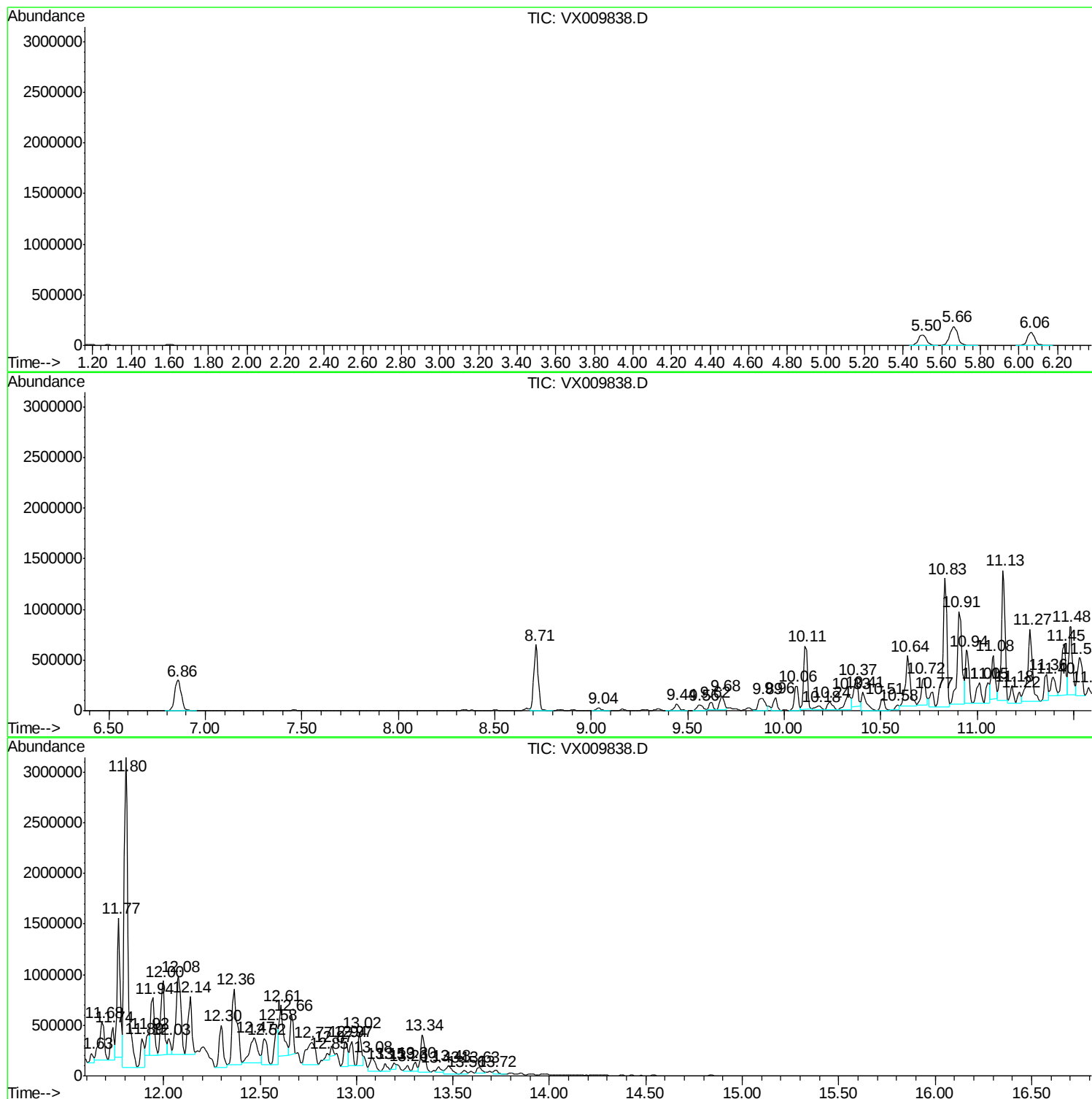
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Instrument :
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ClientSampleId :
EP1-SB-SBR-3BR(14-17)

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



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ALS Vial : 50 Sample Multiplier: 1

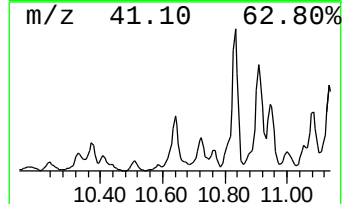
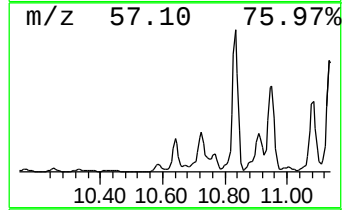
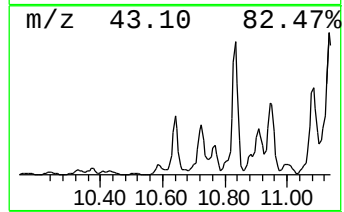
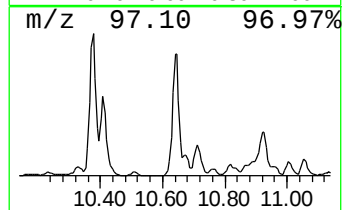
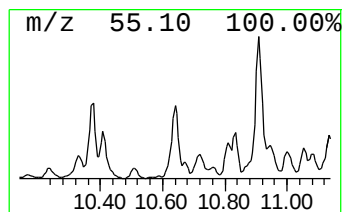
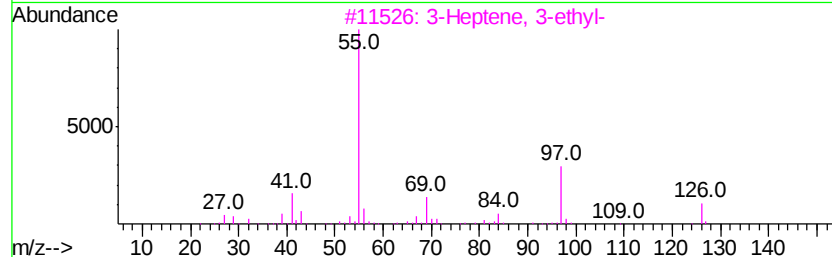
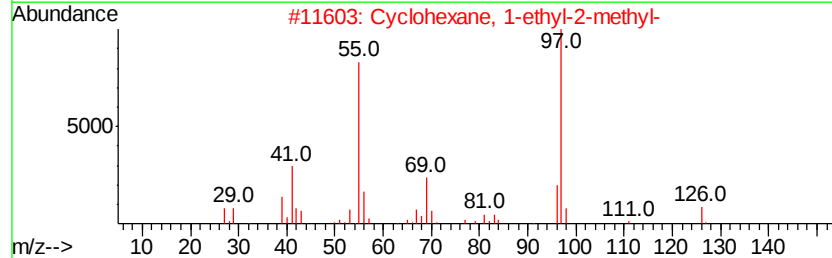
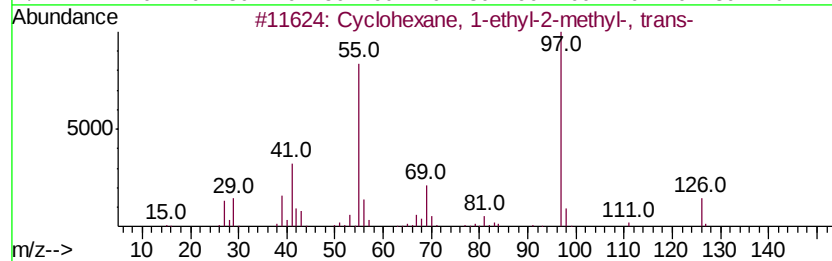
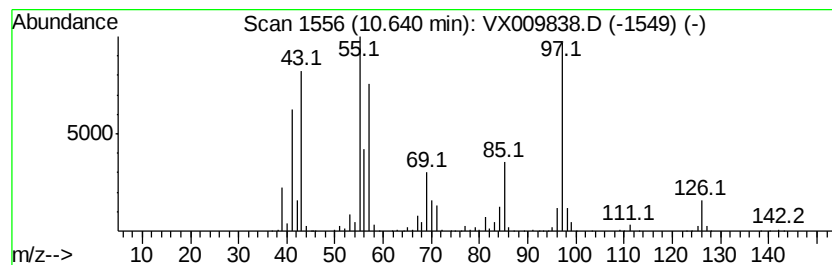
Instrument :
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EP1-SB-SBR-3BR(14-17)

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 1 Cyclohexane, 1-ethyl-2-meth... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.64	50.79 ug/l	874265	Chlorobenzene-d5	10.11		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cvclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	55
2		Cvclohexane, 1-ethyl-2-methyl-	126	C9H18	003728-54-9	55
3		3-Heptene, 3-ethyl-	126	C9H18	074764-46-8	49
4		2-Hexene, 3,4,4-trimethyl-	126	C9H18	053941-19-8	49
5		1-Ethyl-3-methylcyclohexane (c,t)	126	C9H18	003728-55-0	46



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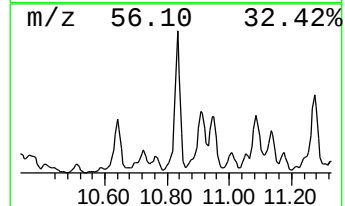
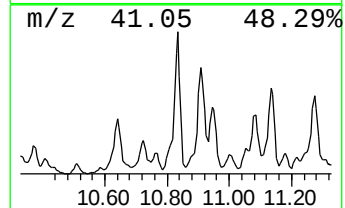
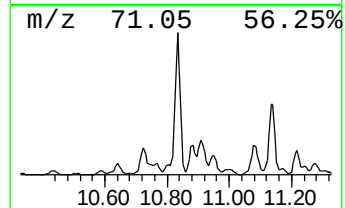
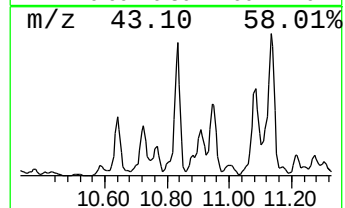
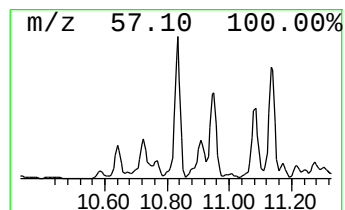
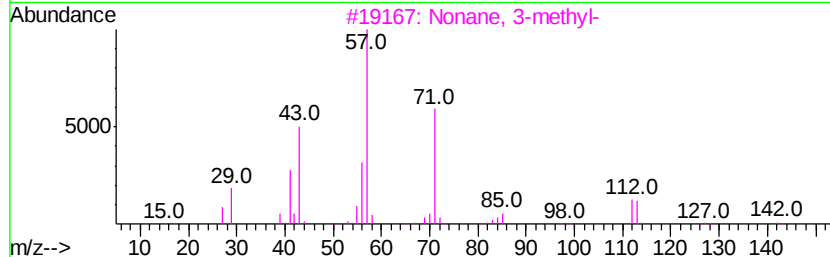
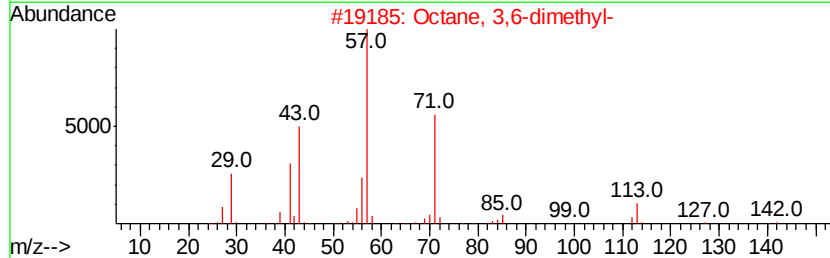
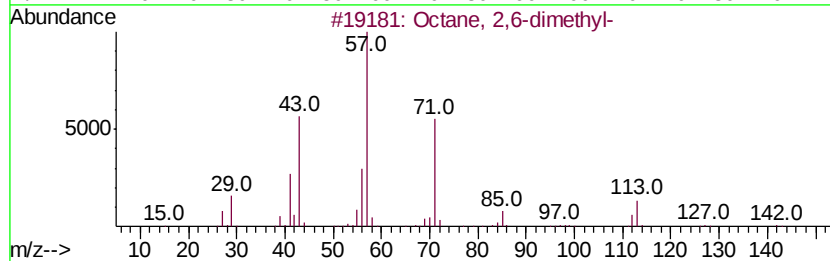
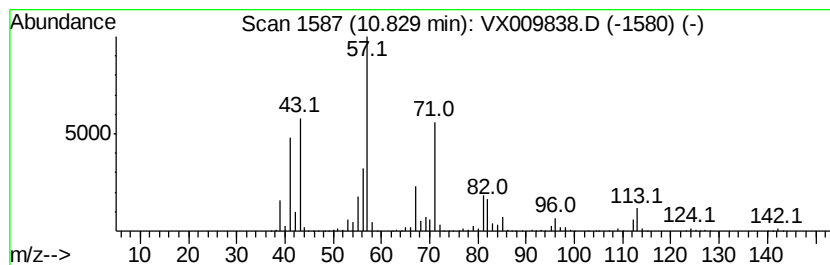
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.83	106.48 ug/l	1832660	Chlorobenzene-d5	10.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	94
2		Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	81
3		Nonane, 3-methyl-	142	C10H22	005911-04-6	76
4		Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	52
5		Sulfurous acid, 2-ethylhexyl hex...	418	C24H50O3S	1000309-19-9	50



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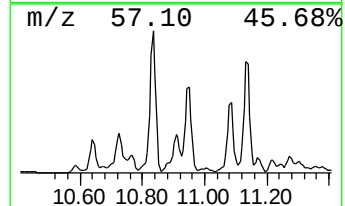
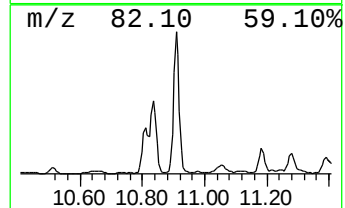
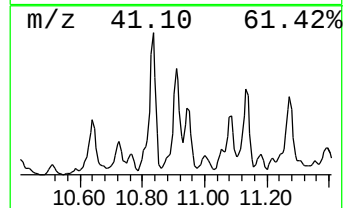
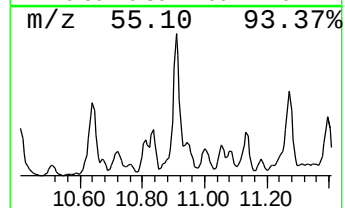
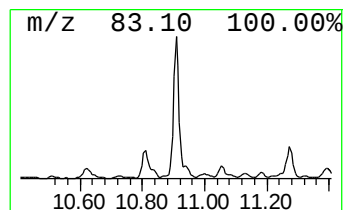
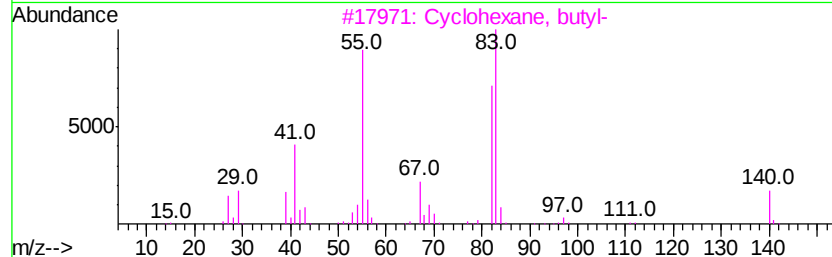
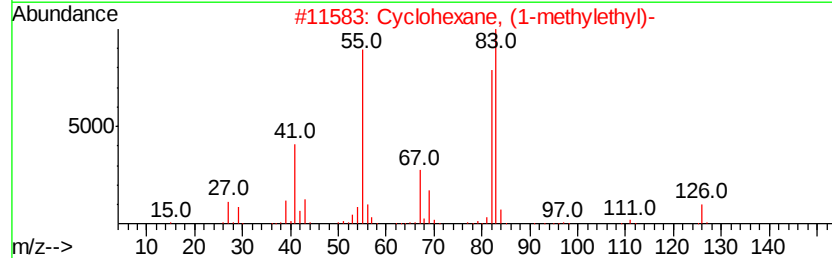
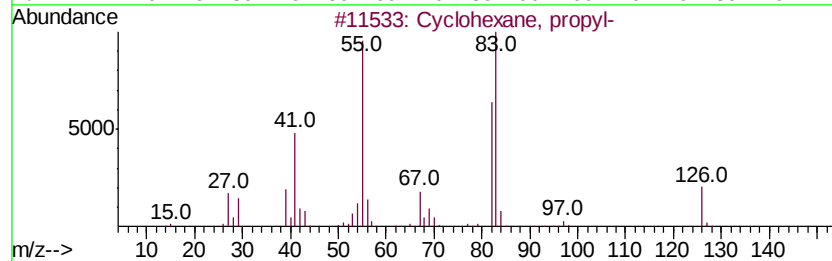
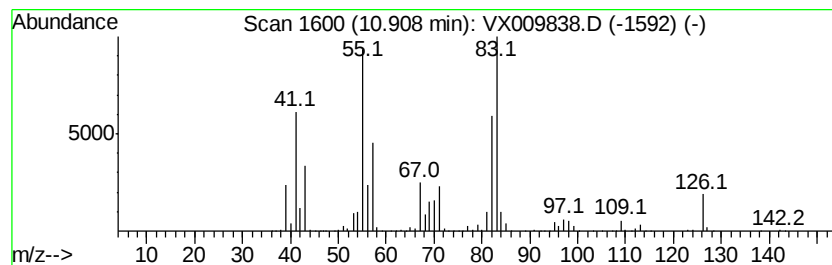
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Quant Title : SW846 8260

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

Peak Number 3 Cyclohexane, propyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.91	90.90 ug/l	1564630	Chlorobenzene-d5	10.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, propyl-	126	C9H18	001678-92-8	76
2		Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	74
3		Cyclohexane, butyl-	140	C10H20	001678-93-9	59
4		Cyclohexane, 2-propenyl-	124	C9H16	002114-42-3	59
5		Cyclohexane, ethyl-	112	C8H16	001678-91-7	53



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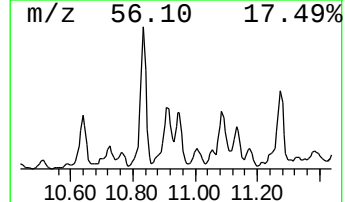
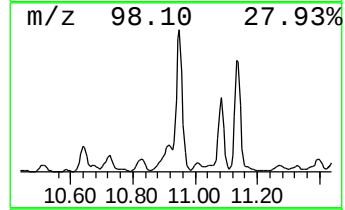
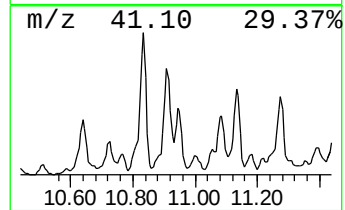
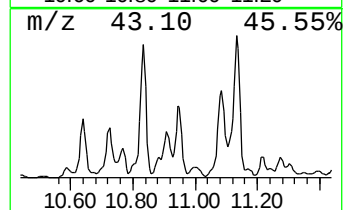
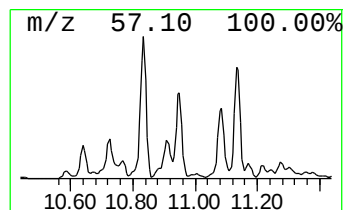
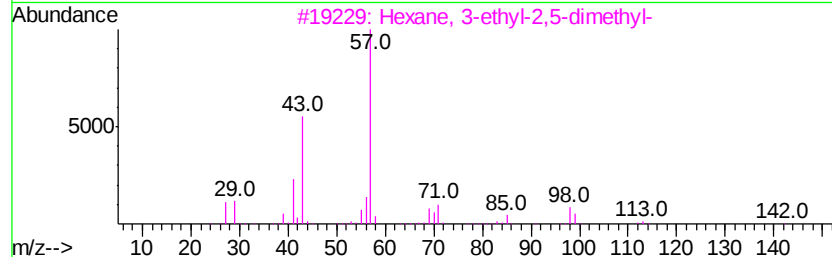
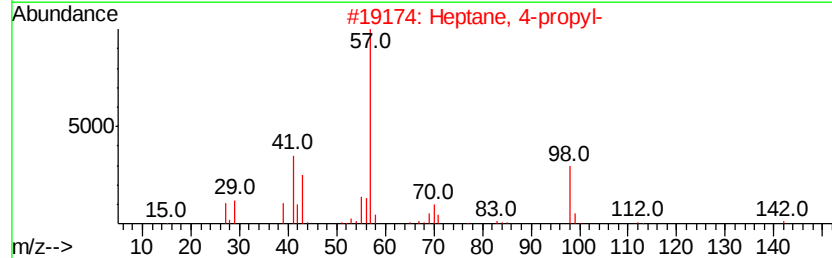
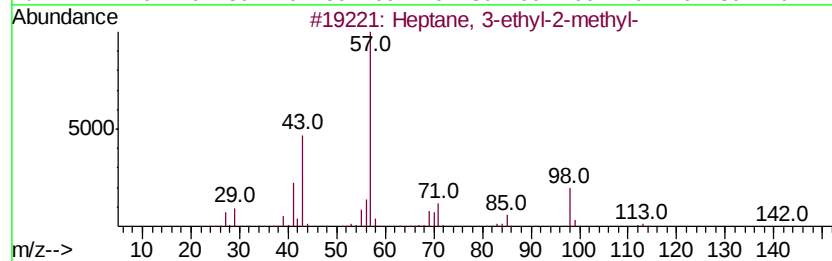
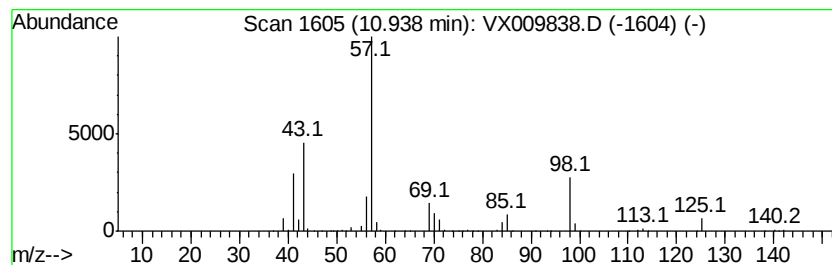
Instrument :
MSVOA_X
ClientSampleId :
EP1-SB-SBR-3BR(14-17)

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, 3-ethyl-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.94	39.23 ug/l	675282	Chlorobenzene-d5	10.11	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	90
2	Heptane, 4-propyl-	142	C10H22	003178-29-8	64
3	Hexane, 3-ethyl-2,5-dimethyl-	142	C10H22	052897-04-8	53
4	Heptane, 4-(1-methyl-ethyl)-	142	C10H22	052896-87-4	43
5	Octane, 2,3-dimethyl-	142	C10H22	007146-60-3	43



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009838.D
Acq On : 23 May 2019 05:34
Operator : JC/SP
Sample : K2976-19 50X
Misc : 6.08g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EP1-SB-SBR-3BR(14-17)

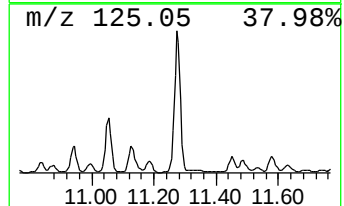
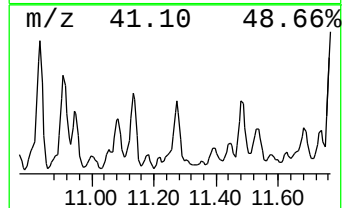
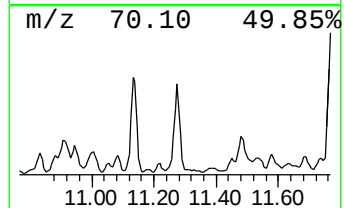
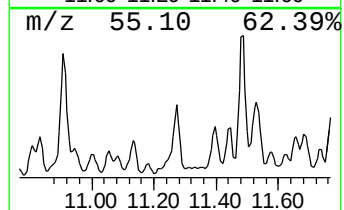
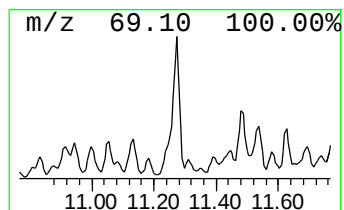
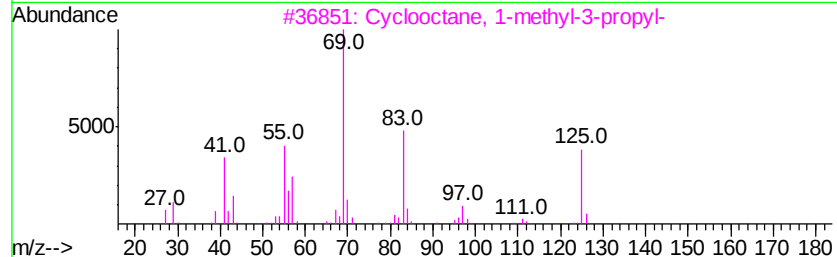
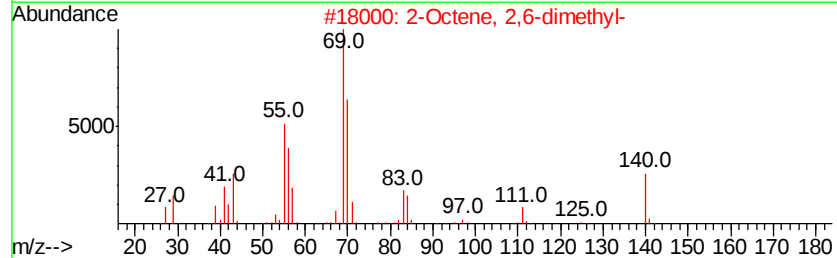
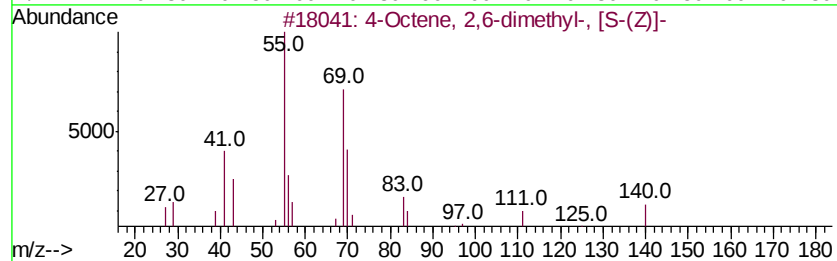
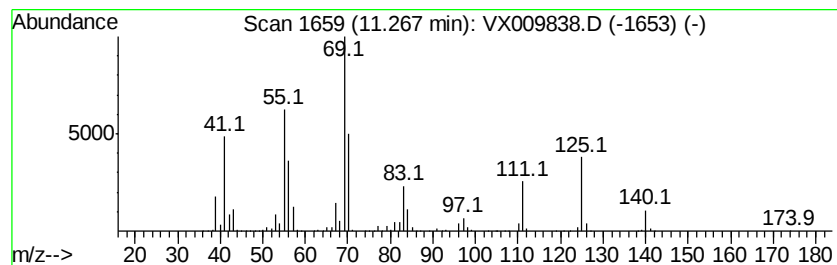
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Quant Title : SW846 8260

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

Peak Number 5 4-Octene, 2,6-dimethyl-, [S... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.27	42.26 ug/l	1205030	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Octene, 2,6-dimethyl-, [S-(Z)]-	140	C10H20	062960-77-4	68
2		2-Octene, 2,6-dimethyl-	140	C10H20	004057-42-5	64
3		Cyclooctane, 1-methyl-3-propyl-	168	C12H24	255885-37-1	50
4		Cyclohexane, 1,1,3,5-tetramethyl...	140	C10H20	050876-31-8	50
5		3-Ethyl-4-octene	140	C10H20	019781-32-9	46



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009838.D
Acq On : 23 May 2019 05:34
Operator : JC/SP
Sample : K2976-19 50X
Misc : 6.08g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 50 Sample Multiplier: 1

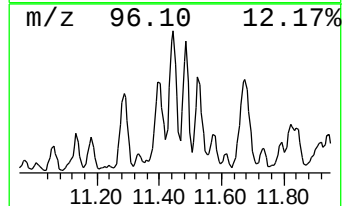
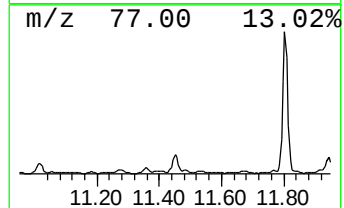
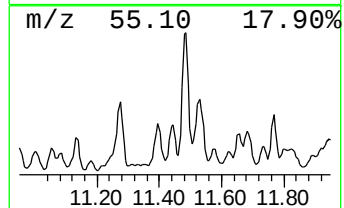
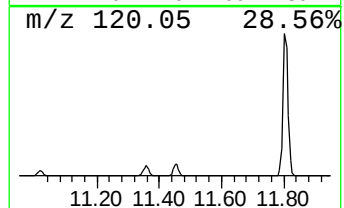
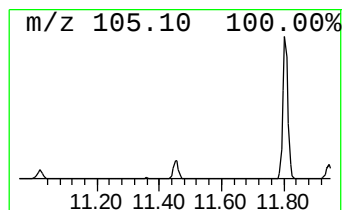
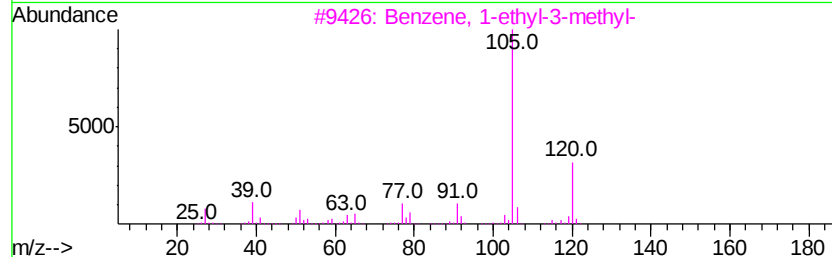
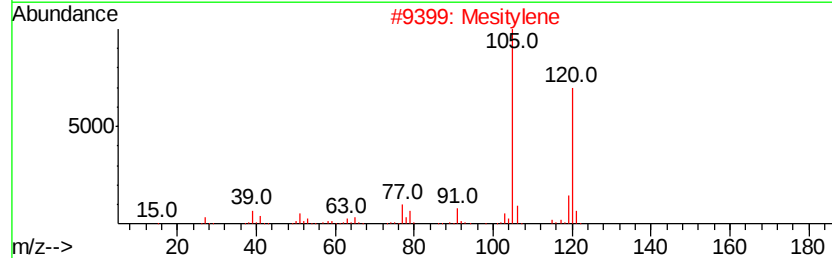
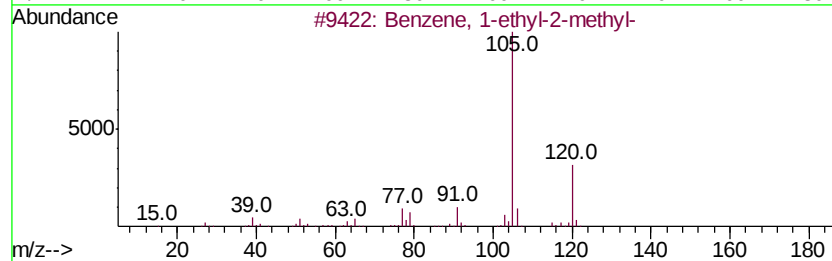
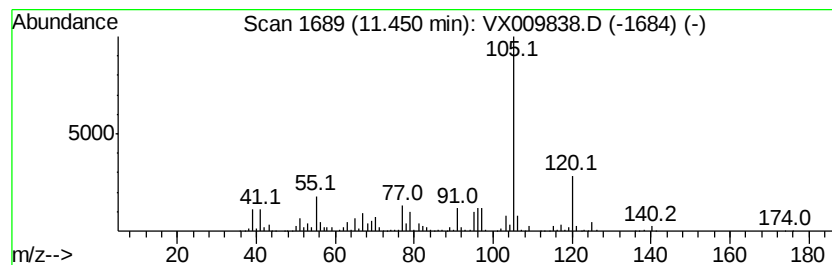
Instrument :
MSVOA_X
ClientSampleId :
EP1-SB-SBR-3BR(14-17)

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 Benzene, 1-ethyl-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.45	27.85 ug/l	794252	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethyl-2-methyl-	120 C9H12	000611-14-3 92
2		Mesitylene	120 C9H12	000108-67-8 64
3		Benzene, 1-ethyl-3-methyl-	120 C9H12	000620-14-4 64
4		Benzene, 1,2,3-trimethyl-	120 C9H12	000526-73-8 64
5		Benzene, 1-ethyl-4-methyl-	120 C9H12	000622-96-8 64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009838.D
Acq On : 23 May 2019 05:34
Operator : JC/SP
Sample : K2976-19 50X
Misc : 6.08g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 50 Sample Multiplier: 1

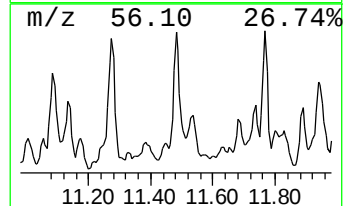
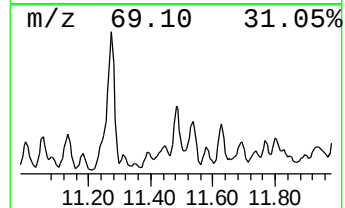
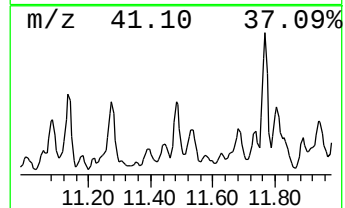
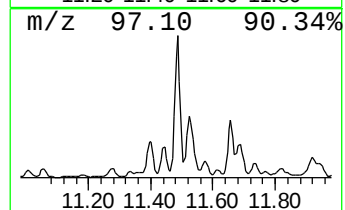
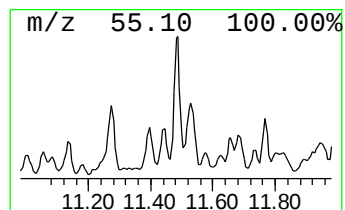
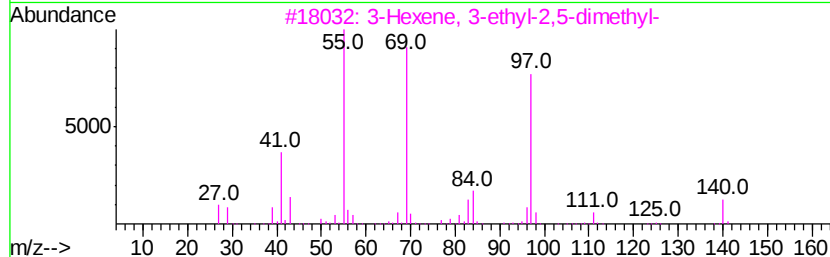
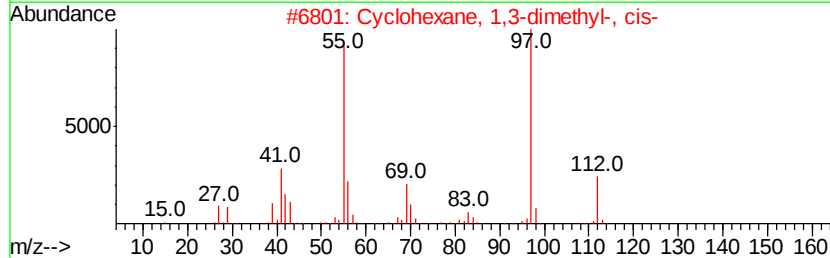
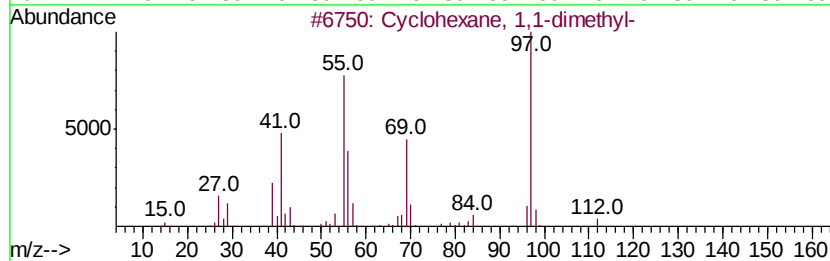
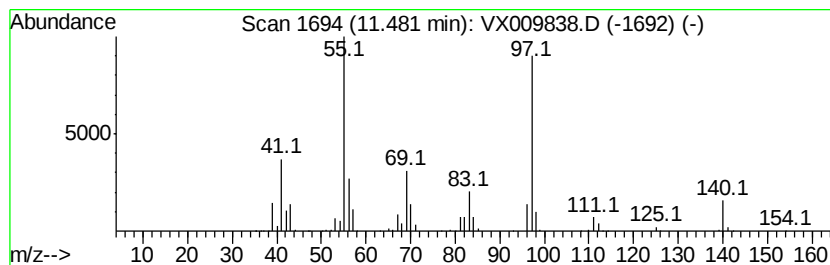
Instrument :
MSVOA_X
ClientSampleId :
EP1-SB-SBR-3BR(14-17)

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, 1,1-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.48	32.19 ug/l	917883	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclohexane, 1,1-dimethyl-	112 C8H16	000590-66-9 76
2		Cyclohexane, 1,3-dimethyl-, cis-	112 C8H16	000638-04-0 72
3		3-Hexene, 3-ethyl-2,5-dimethyl-	140 C10H20	062338-08-3 62
4		Sulfurous acid, cyclohexylmethyl...	276 C14H28O3S	1000309-21-7 59
5		Cyclohexane, 1-methyl-4-(1-methyl...	140 C10H20	006069-98-3 58



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009838.D
Acq On : 23 May 2019 05:34
Operator : JC/SP
Sample : K2976-19 50X
Misc : 6.08g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 50 Sample Multiplier: 1

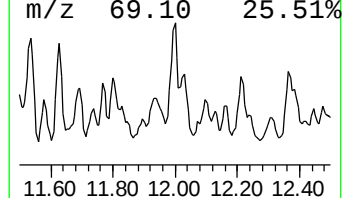
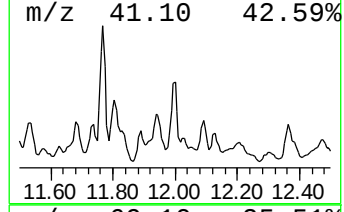
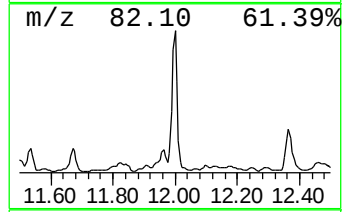
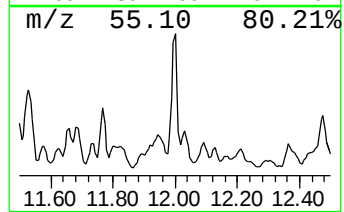
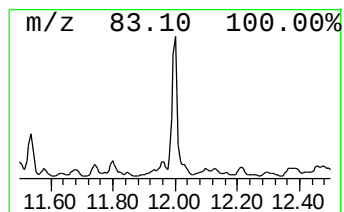
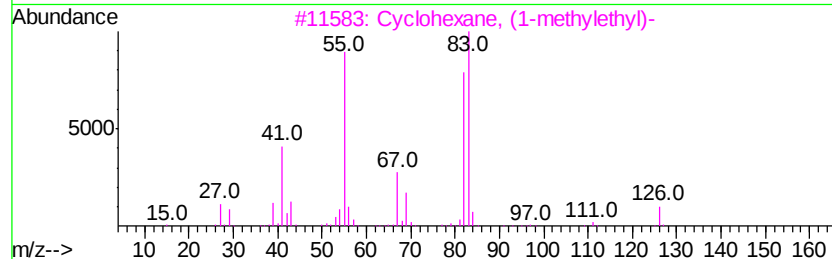
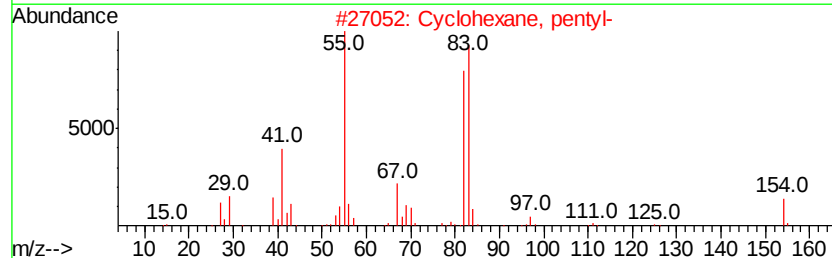
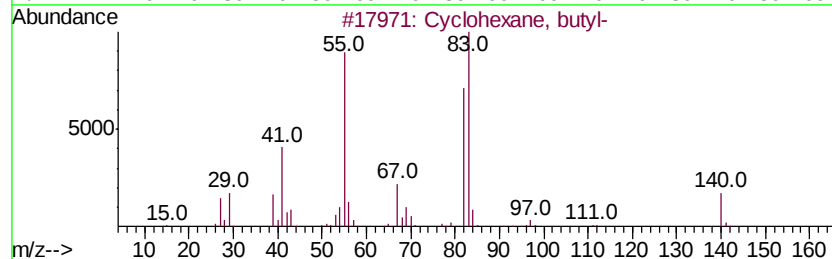
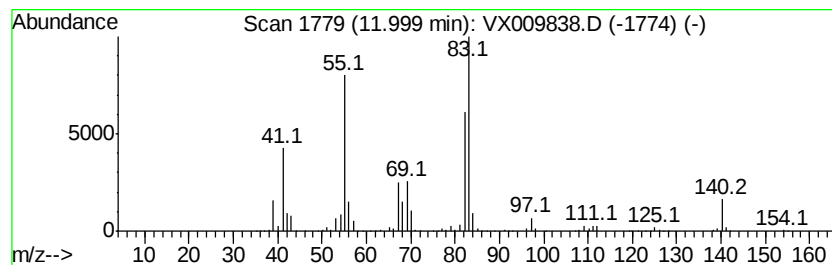
Instrument :
MSVOA_X
ClientSampleId :
EP1-SB-SBR-3BR(14-17)

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 Cyclohexane, butyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.00	34.55 ug/l	985343	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclohexane, butyl-	140 C10H20	001678-93-9 91
2		Cyclohexane, pentyl-	154 C11H22	004292-92-6 72
3		Cyclohexane, (1-methylethyl)-	126 C9H18	000696-29-7 72
4		Cyclohexane, ethyl-	112 C8H16	001678-91-7 64
5		Cyclohexane, octyl-	196 C14H28	001795-15-9 64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009838.D
Acq On : 23 May 2019 05:34
Operator : JC/SP
Sample : K2976-19 50X
Misc : 6.08g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 50 Sample Multiplier: 1

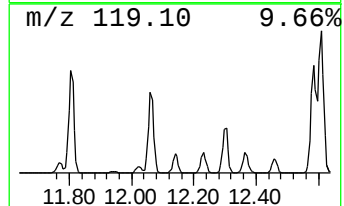
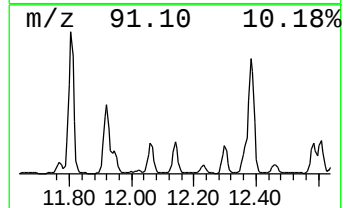
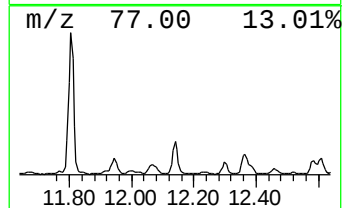
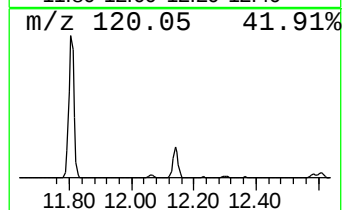
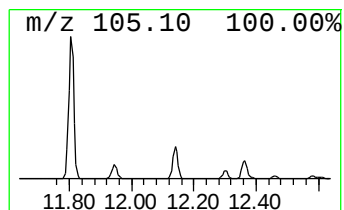
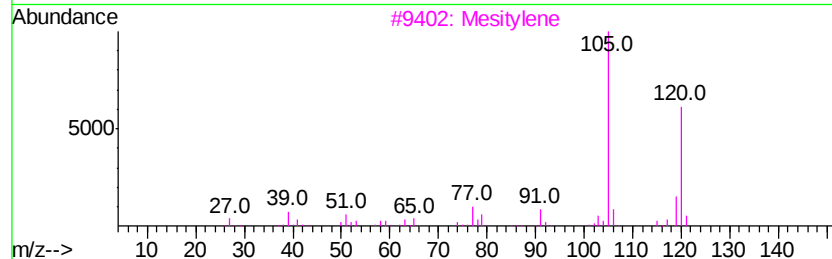
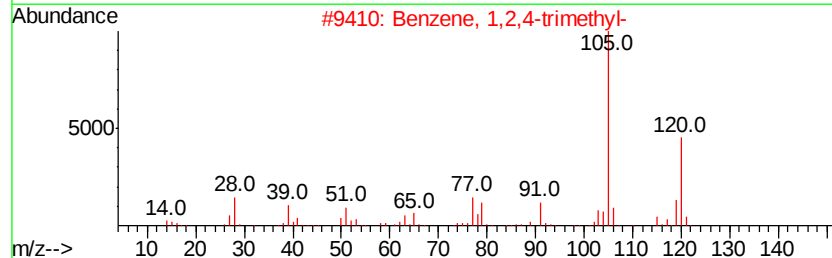
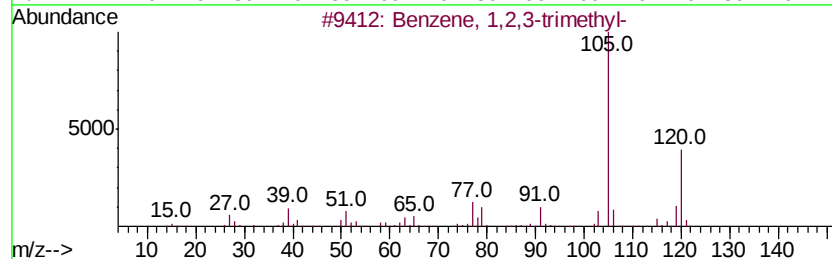
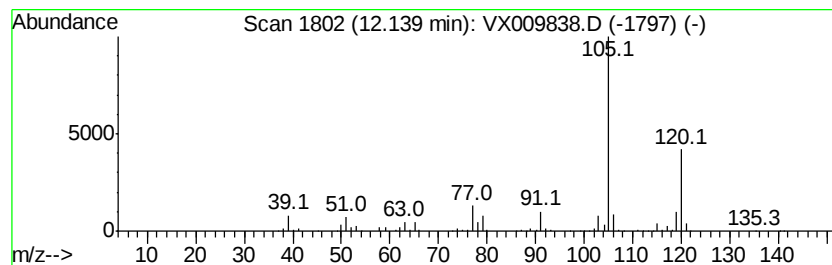
Instrument :
MSVOA_X
ClientSampleId :
EP1-SB-SBR-3BR(14-17)

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, 1,2,3-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	28.93 ug/l	824953	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 91
4		Benzene, 1-ethyl-3-methyl-	120 C9H12	000620-14-4 91
5		Benzene, 1-ethyl-4-methyl-	120 C9H12	000622-96-8 90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009838.D
Acq On : 23 May 2019 05:34
Operator : JC/SP
Sample : K2976-19 50X
Misc : 6.08g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 50 Sample Multiplier: 1

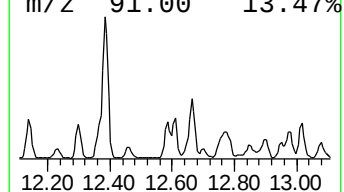
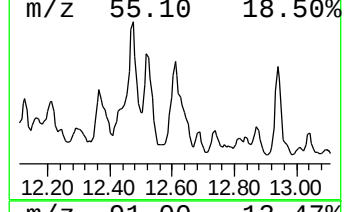
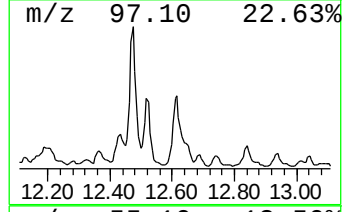
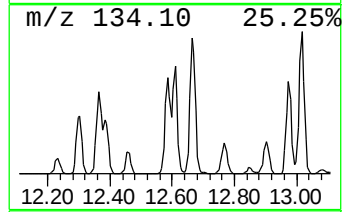
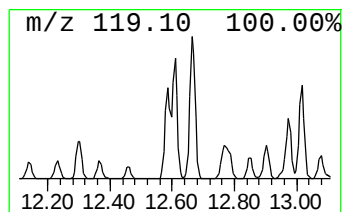
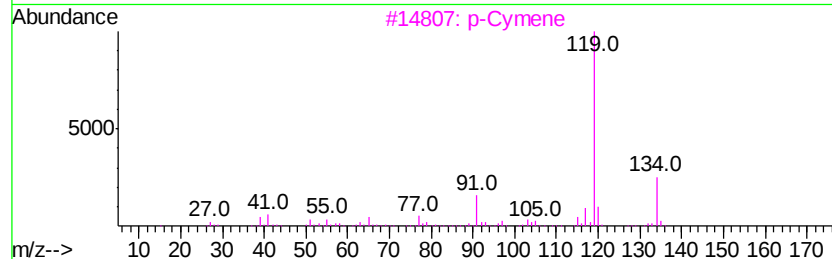
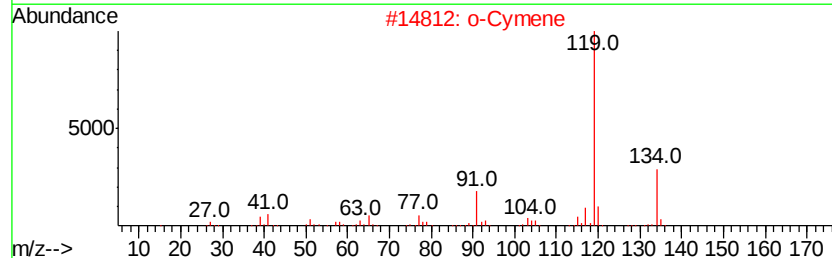
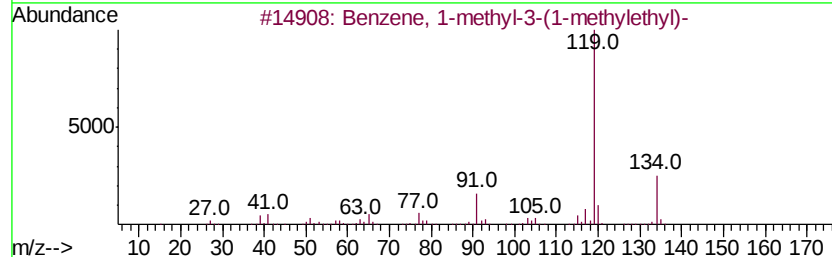
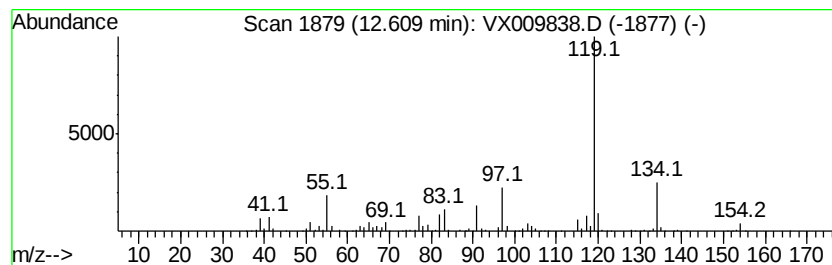
Instrument :
MSVOA_X
ClientSampleId :
EP1-SB-SBR-3BR(14-17)

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-methyl-3-(1-meth... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.61	24.84 ug/l	708413	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzene, 1-methyl-3-(1-methyleth...	134 C10H14	000535-77-3	95
2	o-Cymene	134 C10H14	000527-84-4	94
3	p-Cymene	134 C10H14	000099-87-6	93
4	Benzene, 1-ethyl-2,4-dimethyl-	134 C10H14	000874-41-9	90
5	Benzene, 1,2,3,5-tetramethyl-	134 C10H14	000527-53-7	81



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX052219\
Data File : VX009838.D
Acq On : 23 May 2019 05:34
Operator : JC/SP
Sample : K2976-19 50X
Misc : 6.08g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 50 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EP1-SB-SBR-3BR(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
Cyclohexane, 1-et...	10.64	50.8	ug/l	874265	3	10.11	860591	50.0
Octane, 2,6-dimet...	10.83	106.5	ug/l	1832660	3	10.11	860591	50.0
Cyclohexane, propyl-	10.91	90.9	ug/l	1564630	3	10.11	860591	50.0
Heptane, 3-ethyl-...	10.94	39.2	ug/l	675282	3	10.11	860591	50.0
4-Octene, 2,6-dim...	11.27	42.3	ug/l	1205030	4	12.08	1425790	50.0
Benzene, 1-ethyl-...	11.45	27.9	ug/l	794252	4	12.08	1425790	50.0
Cyclohexane, 1,1-...	11.48	32.2	ug/l	917883	4	12.08	1425790	50.0
Cyclohexane, butyl-	12.00	34.5	ug/l	985343	4	12.08	1425790	50.0
Benzene, 1,2,3-tr...	12.14	28.9	ug/l	824953	4	12.08	1425790	50.0
Benzene, 1-methyl...	12.61	24.8	ug/l	708413	4	12.08	1425790	50.0