

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX052219\
 Data File : VX009822.D
 Acq On : 22 May 2019 23:21
 Operator : JC/SP
 Sample : K2976-01 50X
 Misc : 6.56g/5mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 EP1-SB-E2-3R(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	701	713	727	rBV	96138	282367	3.73%	0.483%
2	5.659	727	739	753	rVV	174045	493414	6.52%	0.844%
3	6.061	793	805	823	rBV	122188	327393	4.33%	0.560%
4	6.860	922	936	950	rBV	283015	671583	8.87%	1.148%
5	8.713	1235	1240	1255	rVB	597732	946423	12.50%	1.618%
6	9.445	1350	1360	1368	rBV	83977	137595	1.82%	0.235%
7	9.561	1373	1379	1385	rBV3	67720	151047	2.00%	0.258%
8	9.622	1385	1389	1393	rVB	121031	164775	2.18%	0.282%
9	9.676	1393	1398	1402	rBV	185584	271438	3.59%	0.464%
10	9.890	1426	1433	1437	rBV3	145115	359297	4.75%	0.614%
11	9.957	1440	1444	1449	rVB	161874	221129	2.92%	0.378%
12	10.067	1455	1462	1465	rBV	317731	452819	5.98%	0.774%
13	10.109	1465	1469	1475	rVB	570402	747280	9.87%	1.278%
14	10.182	1475	1481	1485	rVB5	43416	94297	1.25%	0.161%
15	10.237	1485	1490	1497	rBV2	103108	196926	2.60%	0.337%
16	10.335	1497	1506	1507	rBV2	205437	347022	4.58%	0.593%
17	10.371	1507	1512	1516	rVB2	347145	640778	8.47%	1.096%
18	10.408	1516	1518	1529	rVB	227067	380496	5.03%	0.651%
19	10.512	1529	1535	1541	rBV	143606	214361	2.83%	0.367%
20	10.640	1549	1556	1564	rBV3	571955	1037110	13.70%	1.774%
21	10.725	1564	1570	1573	rBV3	296175	452493	5.98%	0.774%
22	10.768	1574	1577	1580	rVB2	153117	183373	2.42%	0.314%
23	10.835	1580	1588	1592	rBV2	1471915	2132270	28.17%	3.646%
24	10.908	1592	1600	1604	rBV2	1125116	1929194	25.49%	3.299%
25	10.944	1604	1606	1611	rVB	566481	737994	9.75%	1.262%
26	11.012	1611	1617	1621	rBV2	298673	497695	6.58%	0.851%
27	11.054	1621	1624	1626	rBV2	225464	318172	4.20%	0.544%
28	11.085	1626	1629	1632	rVB	381296	460920	6.09%	0.788%
29	11.133	1632	1637	1642	rVB	1324558	1800014	23.78%	3.078%
30	11.182	1642	1645	1648	rVB2	209486	245516	3.24%	0.420%
31	11.219	1648	1651	1652	rBV	119399	123029	1.63%	0.210%
32	11.274	1656	1660	1669	rVB3	884267	1360514	17.97%	2.327%
33	11.359	1670	1674	1677	rBV	597684	736542	9.73%	1.260%
34	11.390	1677	1679	1683	rBV2	155017	185217	2.45%	0.317%

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35	11.450	1683	1689	1692	rBV2	789777	1372891	18.14%	2.348%
36	11.487	1692	1695	1696	rBV	710831	774810	10.24%	1.325%
37	11.505	1696	1698	1701	rVB	665625	617646	8.16%	1.056%
38	11.578	1707	1710	1715	rVB4	140750	180661	2.39%	0.309%
39	11.627	1715	1718	1720	rBV	117423	136984	1.81%	0.234%
40	11.682	1720	1727	1732	rVV6	451851	917097	12.12%	1.568%
41	11.737	1732	1736	1738	rVV2	359660	457228	6.04%	0.782%
42	11.767	1738	1741	1744	rVV	1607257	1877844	24.81%	3.211%
43	11.810	1744	1748	1757	rVB	5450410	7569299	100.00%	12.944%
44	11.889	1757	1761	1763	rBV2	297459	416652	5.50%	0.713%
45	11.920	1763	1766	1767	rVV	356839	427495	5.65%	0.731%
46	11.944	1767	1770	1774	rVV	875747	1184061	15.64%	2.025%
47	11.999	1774	1779	1781	rVV	932318	1237277	16.35%	2.116%
48	12.024	1781	1783	1787	rVV2	538474	661221	8.74%	1.131%
49	12.078	1787	1792	1797	rVV4	696833	1492412	19.72%	2.552%
50	12.139	1797	1802	1809	rVB	1746172	2332946	30.82%	3.990%
51	12.298	1823	1828	1830	rBV	667338	916539	12.11%	1.567%
52	12.322	1830	1832	1835	rVV	605967	656062	8.67%	1.122%
53	12.371	1835	1840	1846	rVV3	1776144	3195258	42.21%	5.464%
54	12.462	1846	1855	1861	rVV3	352632	1026043	13.56%	1.755%
55	12.523	1861	1865	1870	rVB3	309818	575032	7.60%	0.983%
56	12.584	1870	1875	1877	rBV	791129	1101747	14.56%	1.884%
57	12.609	1877	1879	1885	rVV	1052517	1498411	19.80%	2.562%
58	12.664	1885	1888	1892	rVV	949821	1256138	16.60%	2.148%
59	12.694	1892	1893	1897	rVV	168439	177224	2.34%	0.303%
60	12.767	1897	1905	1911	rVB4	527305	1376217	18.18%	2.353%
61	12.847	1914	1918	1920	rVV3	217829	301963	3.99%	0.516%
62	12.901	1924	1927	1930	rVB2	348881	431501	5.70%	0.738%
63	12.944	1930	1934	1936	rBV2	296133	424116	5.60%	0.725%
64	12.975	1936	1939	1942	rVB	444037	515665	6.81%	0.882%
65	13.017	1942	1946	1953	rVB	734313	1011610	13.36%	1.730%
66	13.078	1953	1956	1961	rBV3	192506	348216	4.60%	0.595%
67	13.145	1965	1967	1972	rVB	95157	118844	1.57%	0.203%
68	13.194	1972	1975	1976	rBV2	120758	127495	1.68%	0.218%
69	13.261	1983	1986	1990	rVB3	96417	116090	1.53%	0.199%
70	13.304	1990	1993	1996	rVV2	164608	213241	2.82%	0.365%
71	13.340	1996	1999	2006	rVV2	731374	1054390	13.93%	1.803%

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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	81827	110321	1.46%	0.189%
73	13.474	2017	2021	2031	rVB2	274700	447361	5.91%	0.765%
74	13.560	2031	2035	2038	rBV3	54272	78915	1.04%	0.135%
75	13.639	2044	2048	2052	rBV3	108772	180375	2.38%	0.308%
76	13.718	2059	2061	2067	rVB2	68110	96957	1.28%	0.166%
77	13.834	2076	2080	2087	rVB2	102331	164273	2.17%	0.281%

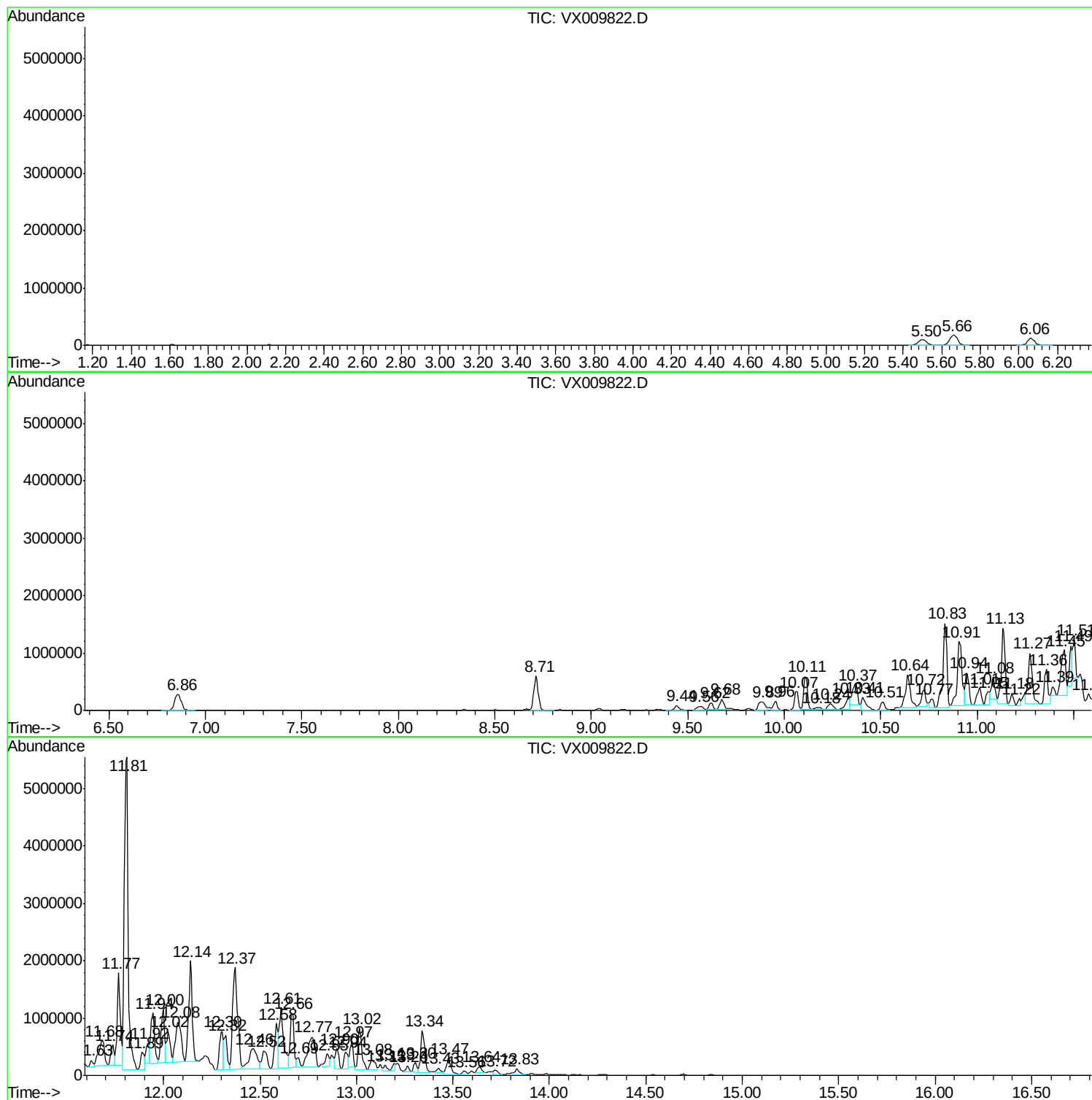
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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



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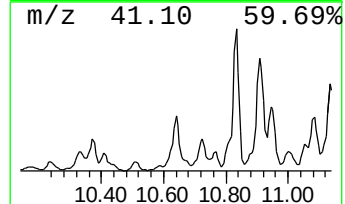
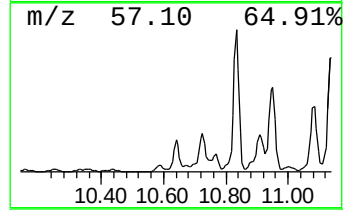
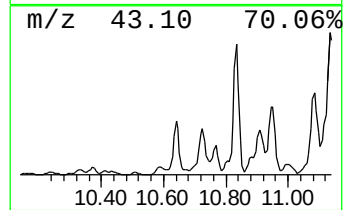
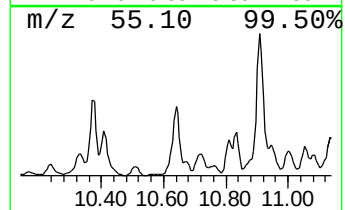
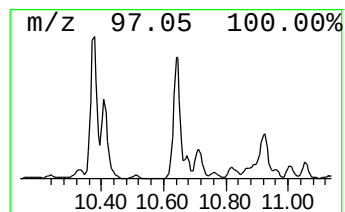
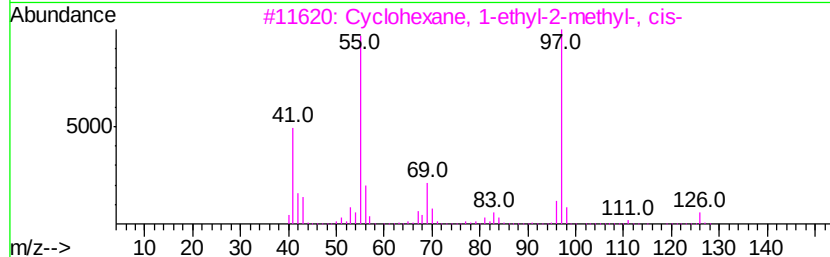
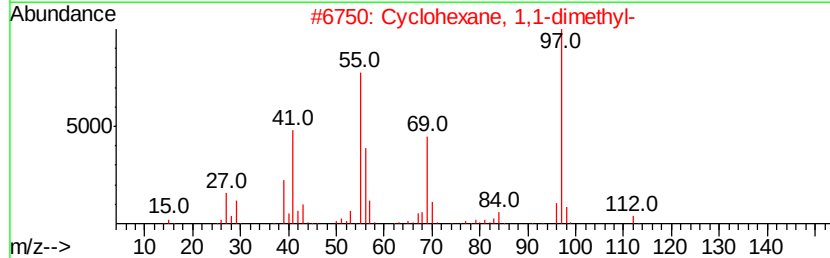
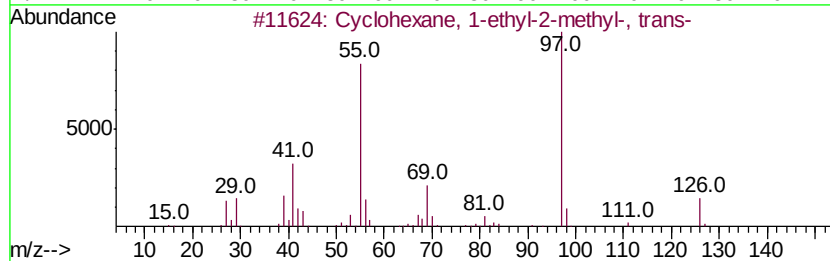
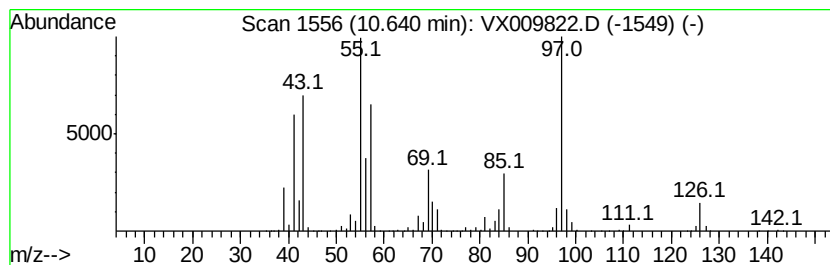
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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	69.39 ug/l	1037110	Chlorobenzene-d5	10.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	64
2		Cyclohexane, 1,1-dimethyl-	112	C8H16	000590-66-9	53
3		Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-77-7	53
4		1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	52
5		Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	52



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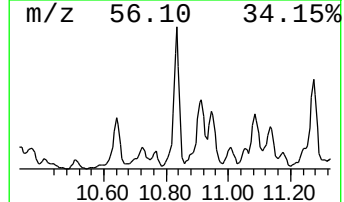
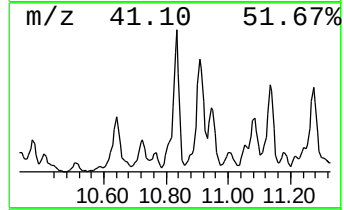
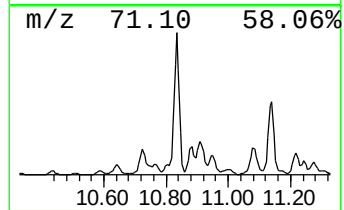
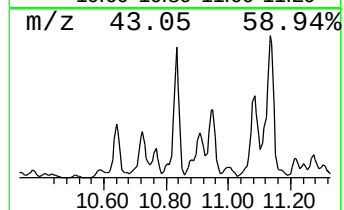
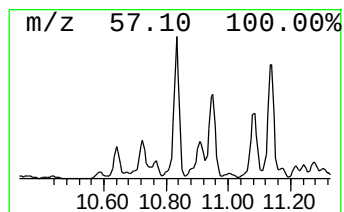
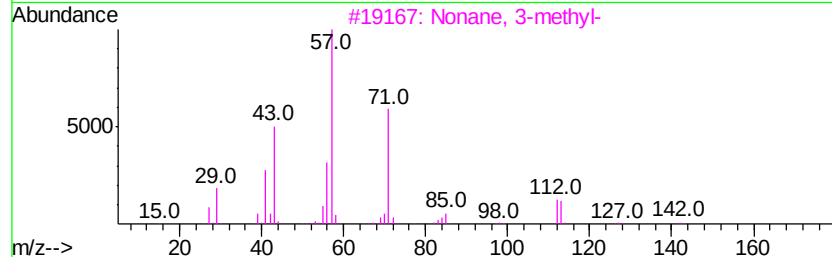
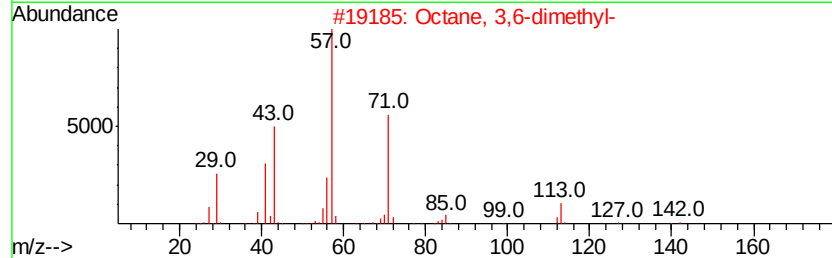
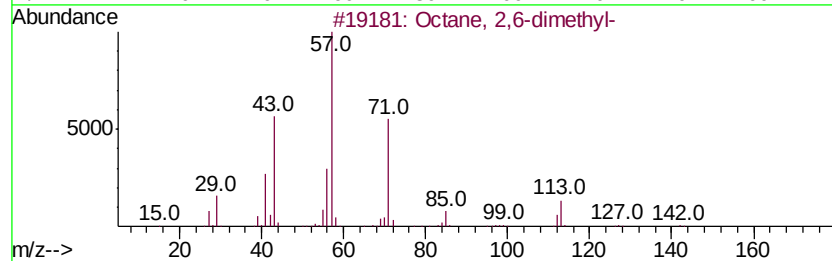
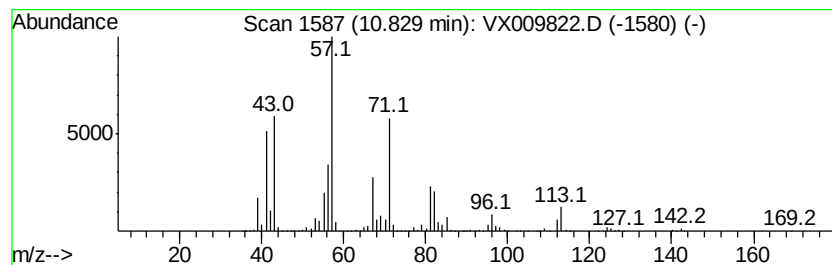
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Peak Number 2 Octane, 2,6-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.83	142.67 ug/l	2132270	Chlorobenzene-d5	10.11		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	92
2		Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	64
3		Nonane, 3-methyl-	142	C10H22	005911-04-6	60
4		Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	46
5		Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1000309-19-2	46



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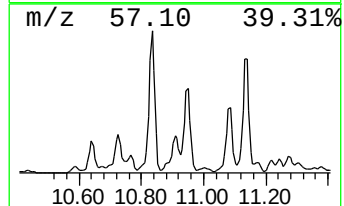
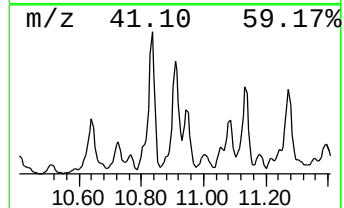
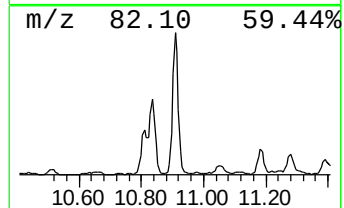
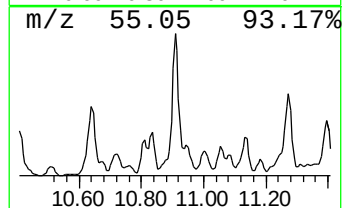
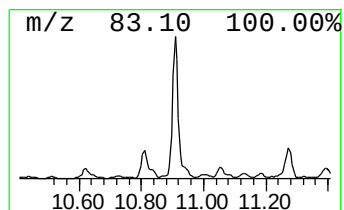
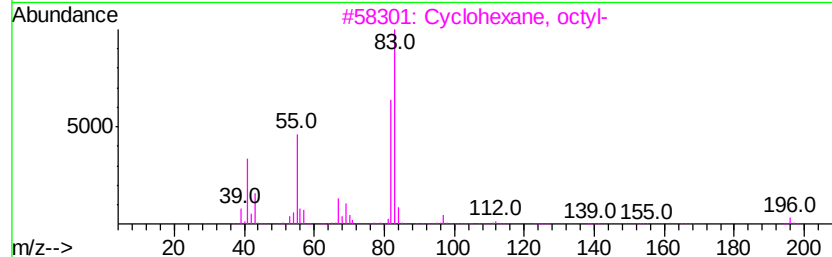
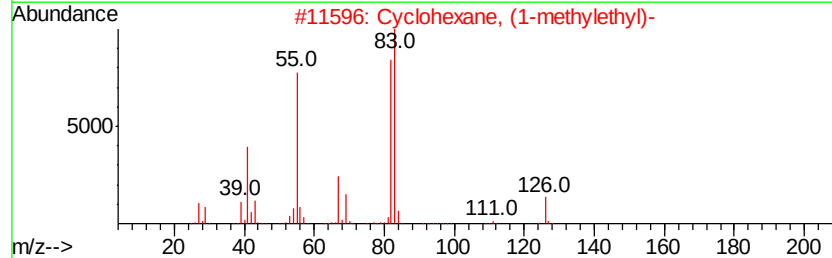
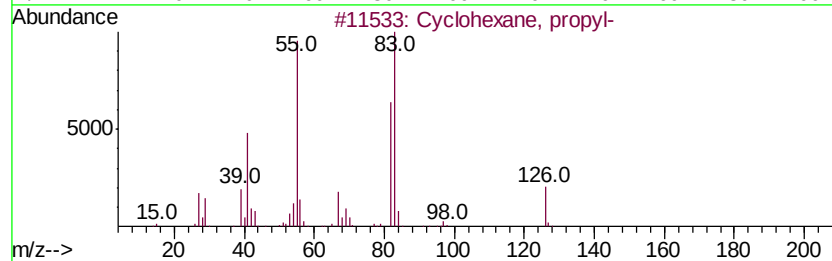
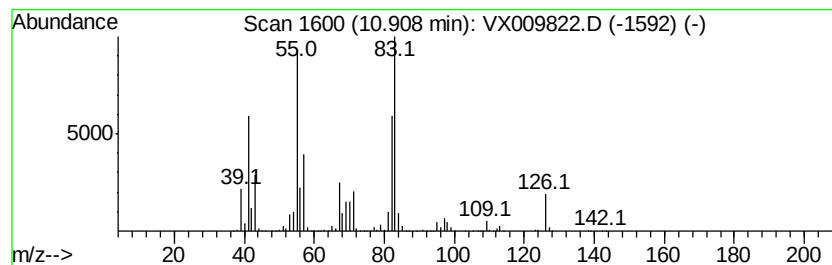
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Quant Title : SW846 8260

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TIC Integration Parameters: LSCINT.P

Peak Number 3 Cyclohexane, propyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.91	129.08 ug/l	1929190	Chlorobenzene-d5	10.11
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclohexane, propyl-	126 C9H18	001678-92-8 76
2		Cyclohexane, (1-methylethyl)-	126 C9H18	000696-29-7 60
3		Cyclohexane, octyl-	196 C14H28	001795-15-9 59
4		Cyclohexane, 2-propenyl-	124 C9H16	002114-42-3 59
5		Cyclohexane, butyl-	140 C10H20	001678-93-9 59



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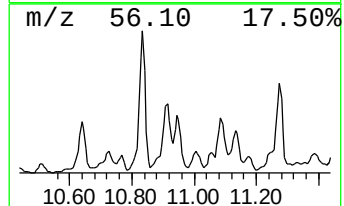
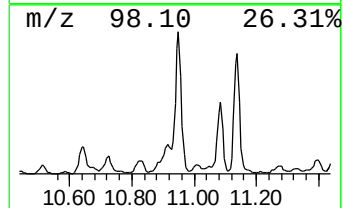
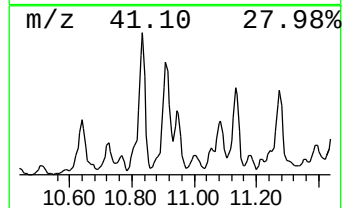
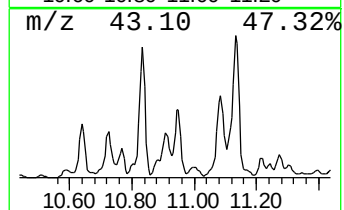
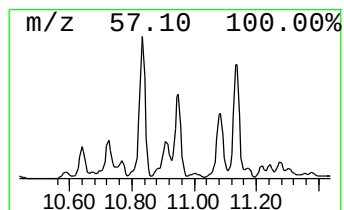
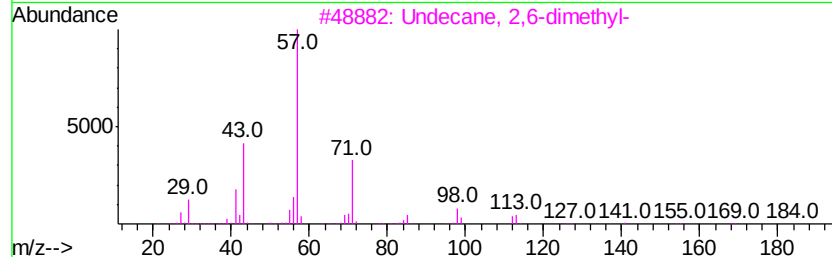
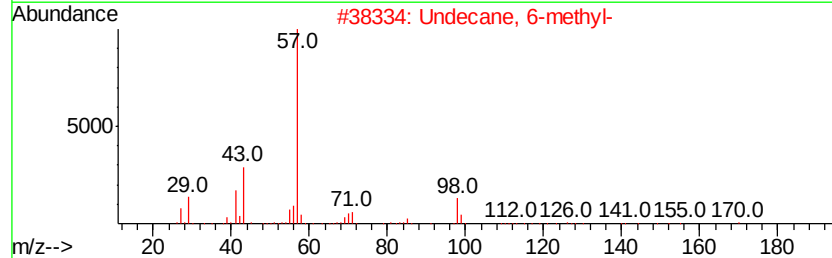
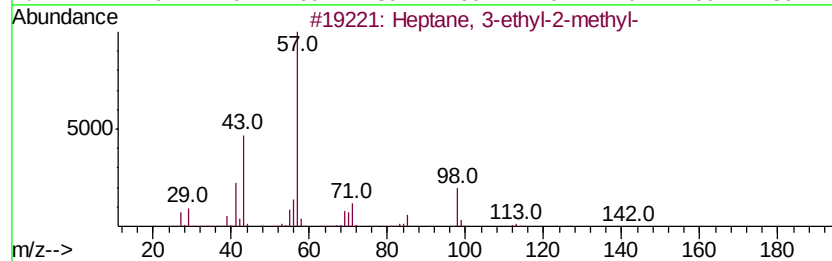
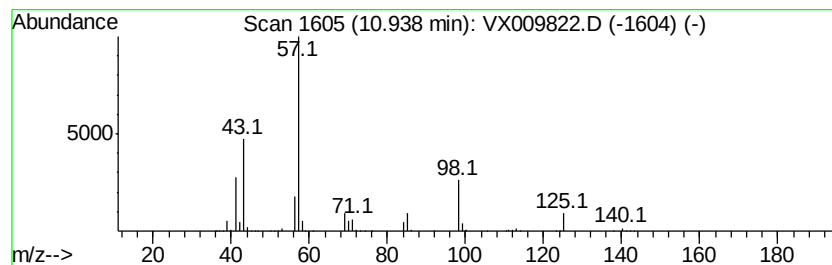
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Quant Title : SW846 8260

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

Peak Number 4 Heptane, 3-ethyl-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.94	49.38 ug/l	737994	Chlorobenzene-d5	10.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	64
2		Undecane, 6-methyl-	170	C12H26	017302-33-9	64
3		Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	50
4		Heptane, 4-propyl-	142	C10H22	003178-29-8	45
5		Hexane, 3-ethyl-2,5-dimethyl-	142	C10H22	052897-04-8	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009822.D
Acq On : 22 May 2019 23:21
Operator : JC/SP
Sample : K2976-01 50X
Misc : 6.56g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 34 Sample Multiplier: 1

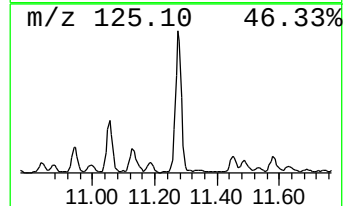
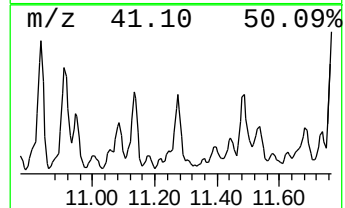
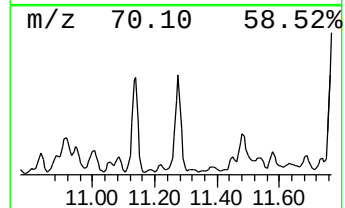
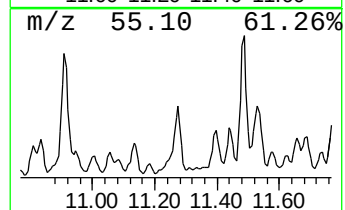
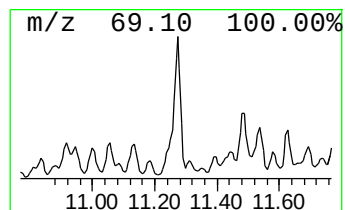
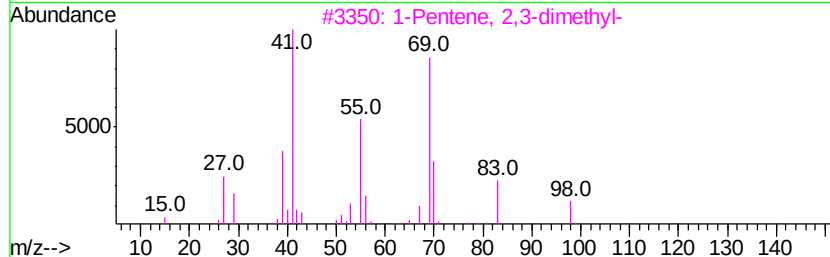
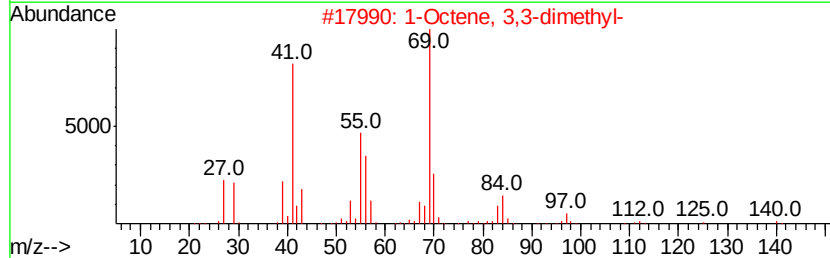
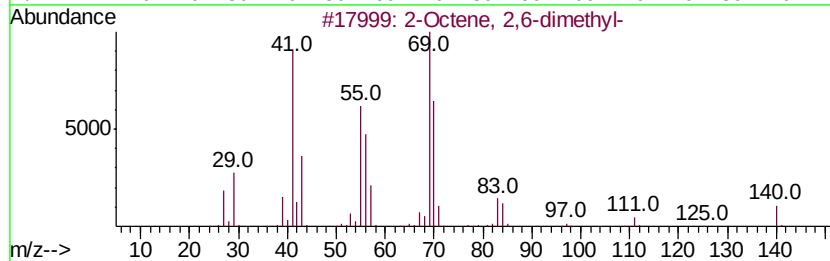
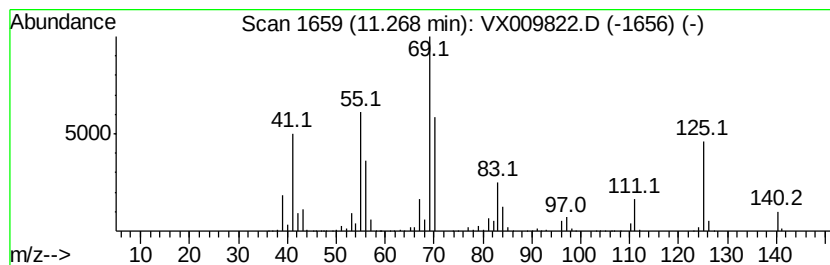
Instrument :
MSVOA_X
ClientSampleId :
EP1-SB-E2-3R(14-17)

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 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.27	45.58 ug/l	1360510	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	1-Octene, 3,3-dimethyl-	140 C10H20	074511-51-6	38
3	1-Pentene, 2,3-dimethyl-	98 C7H14	003404-72-6	38
4	3-Nonene, 2-methyl-	140 C10H20	053966-53-3	38
5	1,1,4-Trimethylcyclohexane	126 C9H18	007094-27-1	35



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009822.D
Acq On : 22 May 2019 23:21
Operator : JC/SP
Sample : K2976-01 50X
Misc : 6.56g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 34 Sample Multiplier: 1

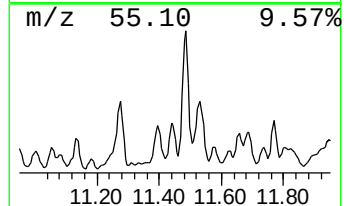
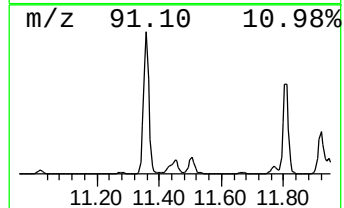
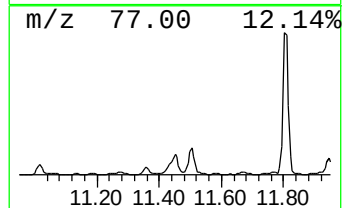
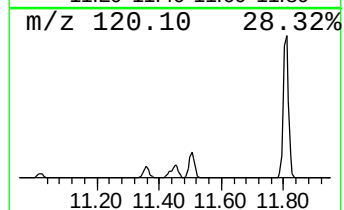
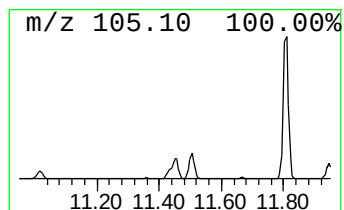
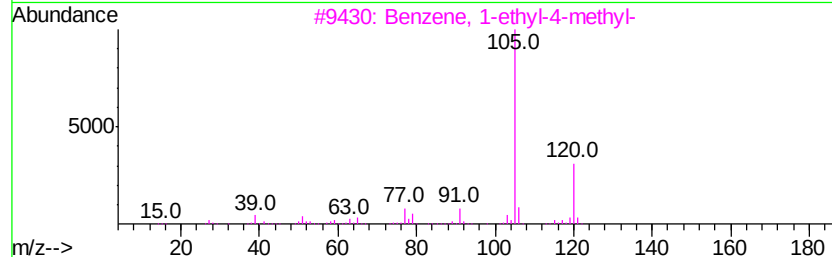
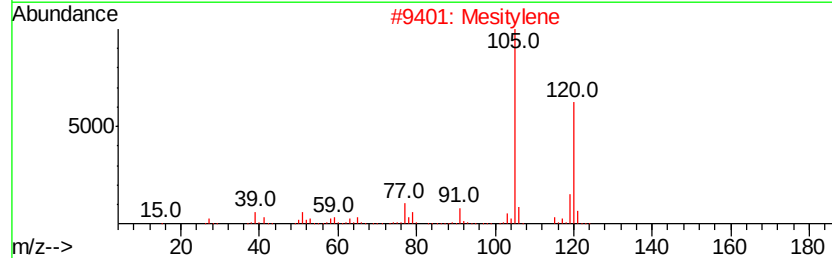
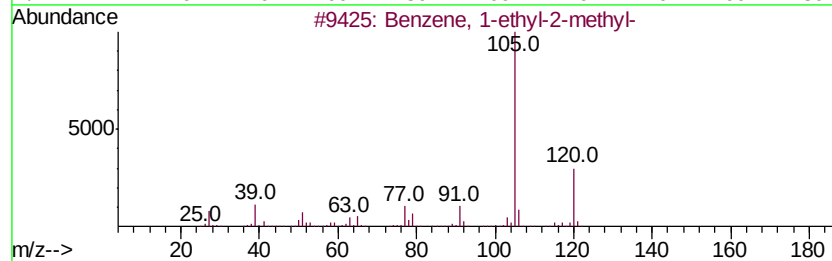
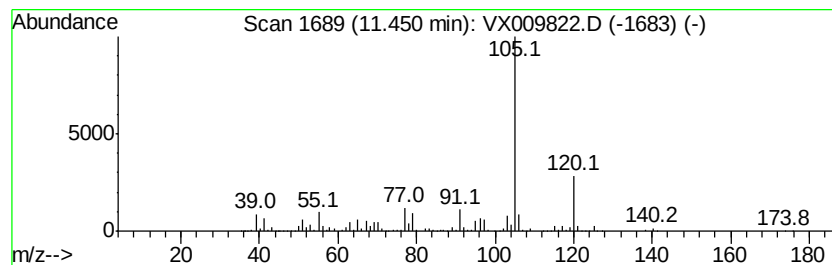
Instrument :
MSVOA_X
ClientSampleId :
EP1-SB-E2-3R(14-17)

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 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.45	46.00 ug/l	1372890	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 94
2		Mesitylene	120 C9H12	000108-67-8 81
3		Benzene, 1-ethyl-4-methyl-	120 C9H12	000622-96-8 81
4		Benzene, 1-ethyl-3-methyl-	120 C9H12	000620-14-4 81
5		Benzene, 1,2,4-trimethyl-	120 C9H12	000095-63-6 81



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009822.D
Acq On : 22 May 2019 23:21
Operator : JC/SP
Sample : K2976-01 50X
Misc : 6.56g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 34 Sample Multiplier: 1

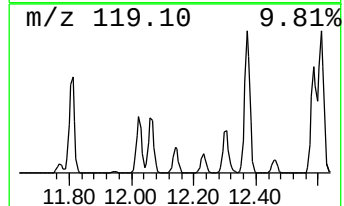
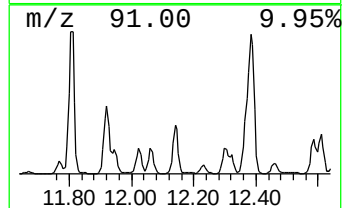
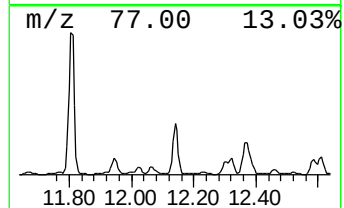
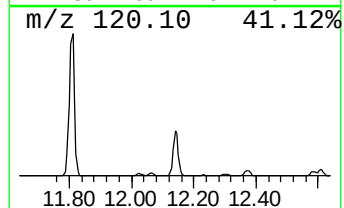
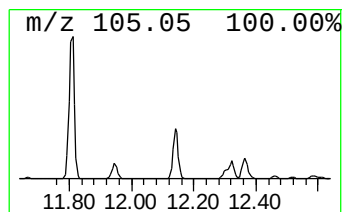
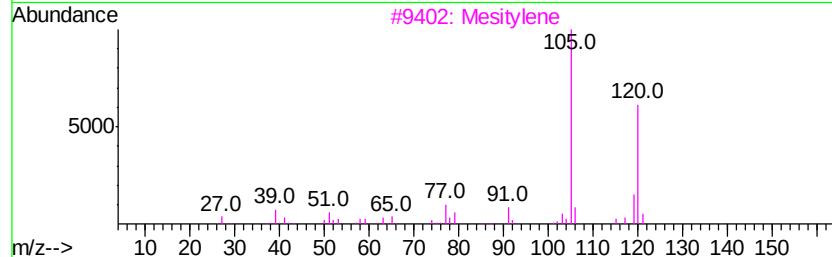
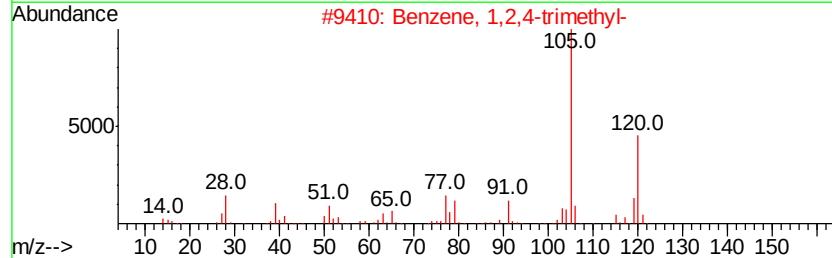
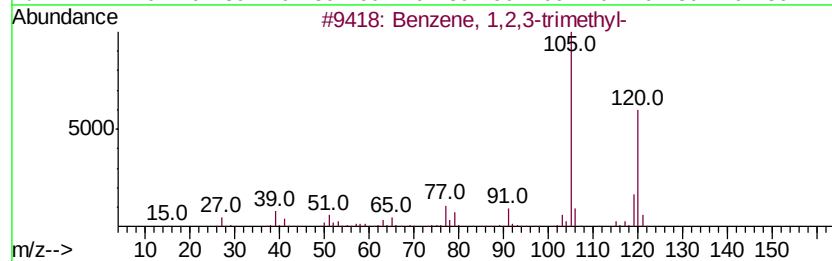
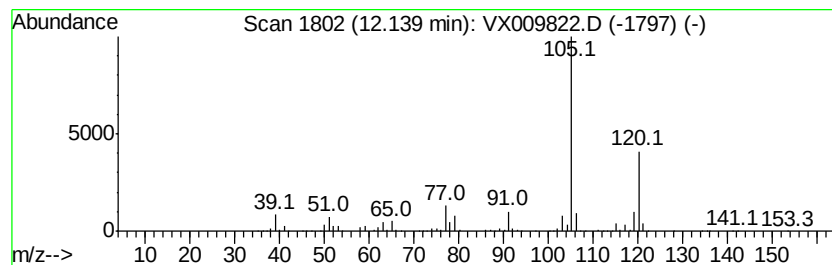
Instrument :
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ClientSampleId :
EP1-SB-E2-3R(14-17)

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	78.16 ug/l	2332950	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-4-methyl-	120 C9H12	000622-96-8 90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009822.D
Acq On : 22 May 2019 23:21
Operator : JC/SP
Sample : K2976-01 50X
Misc : 6.56g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 34 Sample Multiplier: 1

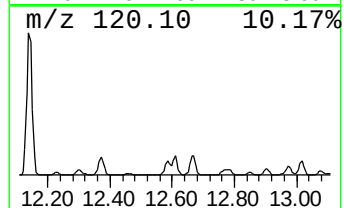
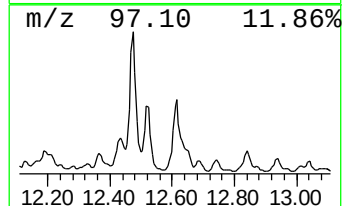
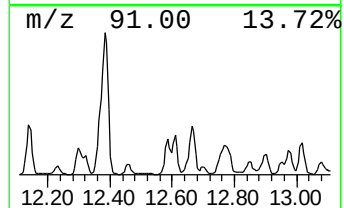
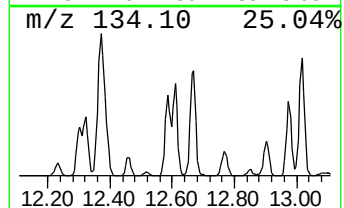
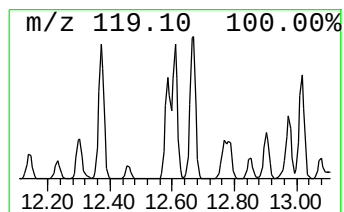
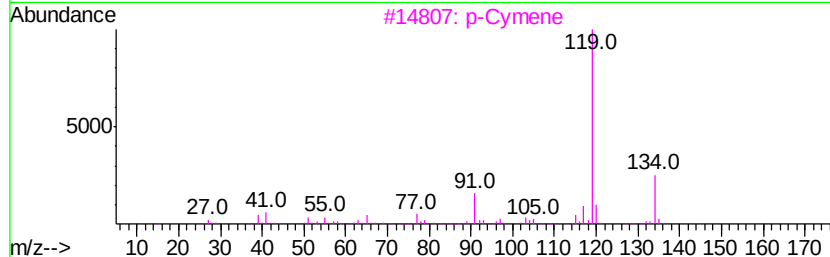
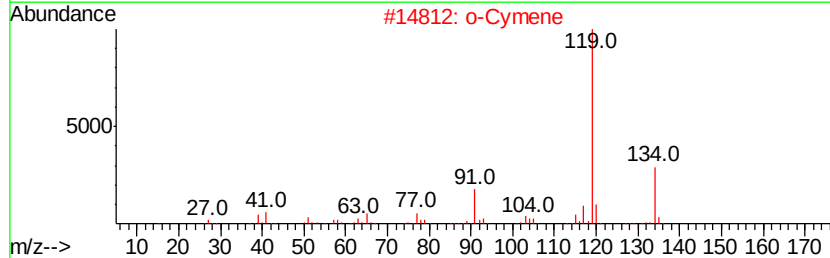
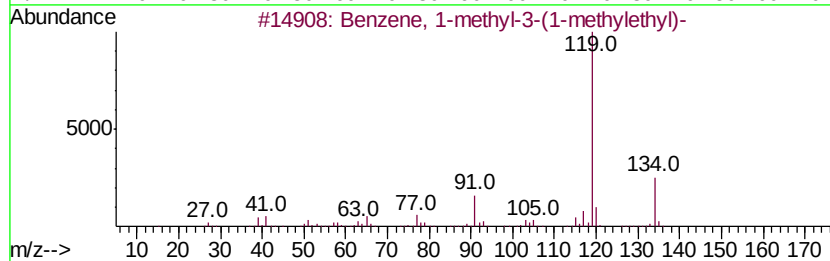
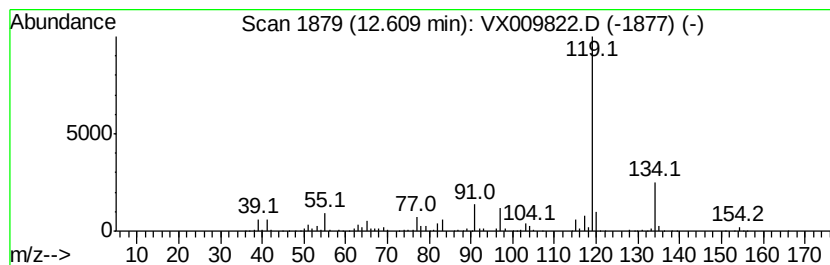
Instrument :
MSVOA_X
ClientSampleId :
EP1-SB-E2-3R(14-17)

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-methyl-3-(1-meth... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.61	50.20 ug/l	1498410	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	95
3	p-Cymene	134	C10H14	000099-87-6	93
4	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009822.D
Acq On : 22 May 2019 23:21
Operator : JC/SP
Sample : K2976-01 50X
Misc : 6.56g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 34 Sample Multiplier: 1

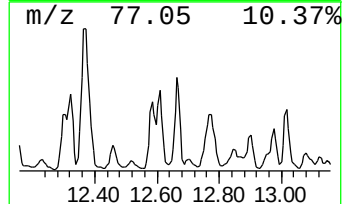
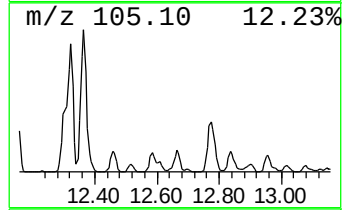
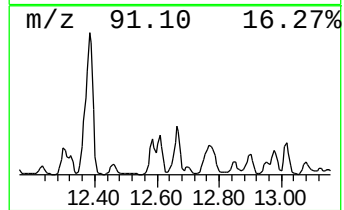
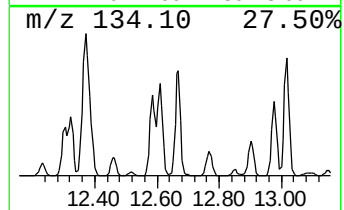
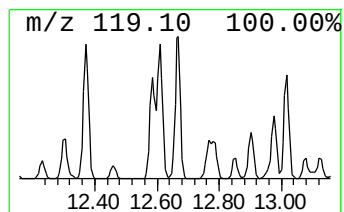
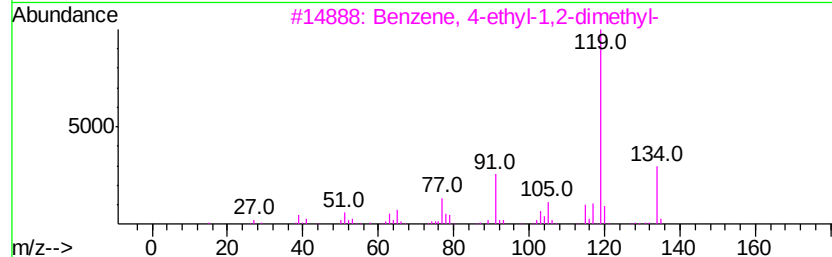
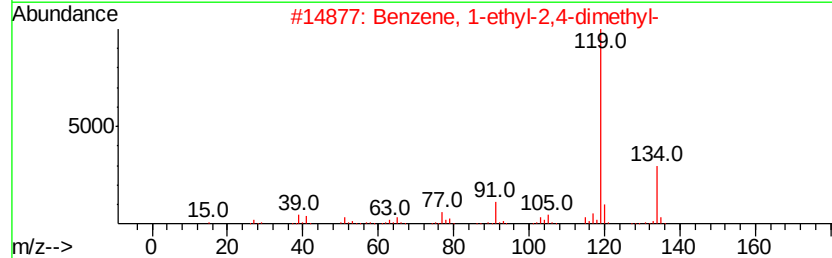
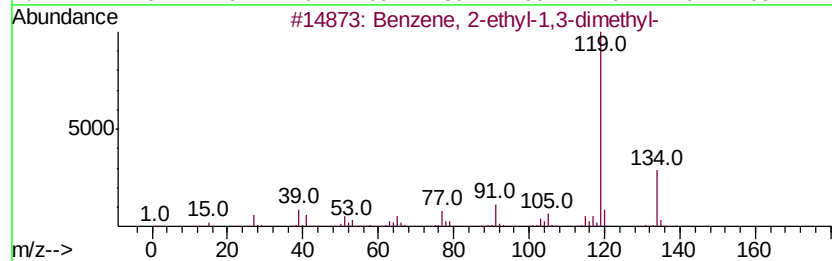
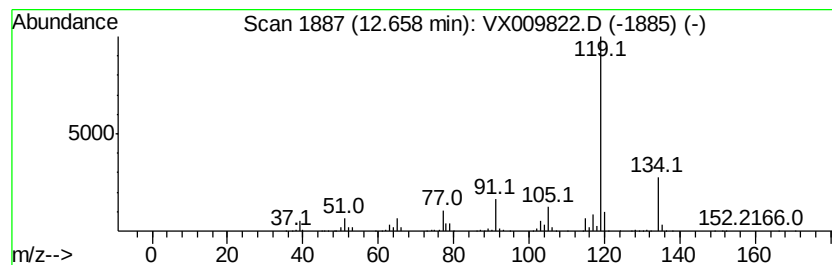
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ClientSampleId :
EP1-SB-E2-3R(14-17)

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, 2-ethyl-1,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.66	42.08 ug/l	1256140	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzene, 2-ethyl-1,3-dimethyl-	134 C10H14	002870-04-4	95
2	Benzene, 1-ethyl-2,4-dimethyl-	134 C10H14	000874-41-9	95
3	Benzene, 4-ethyl-1,2-dimethyl-	134 C10H14	000934-80-5	95
4	Benzene, 1-ethyl-2,3-dimethyl-	134 C10H14	000933-98-2	95
5	Benzene, 2-ethyl-1,4-dimethyl-	134 C10H14	001758-88-9	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009822.D
Acq On : 22 May 2019 23:21
Operator : JC/SP
Sample : K2976-01 50X
Misc : 6.56g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 34 Sample Multiplier: 1

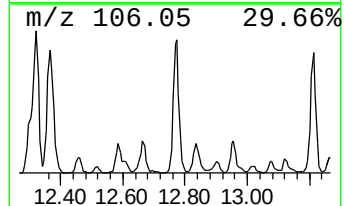
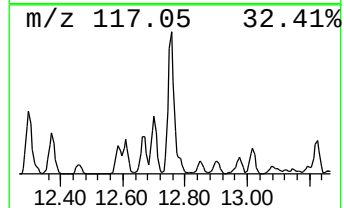
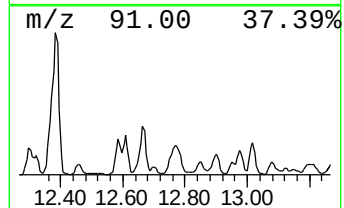
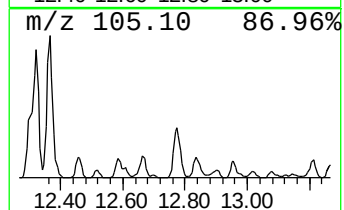
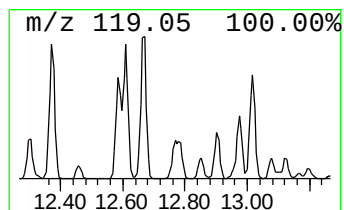
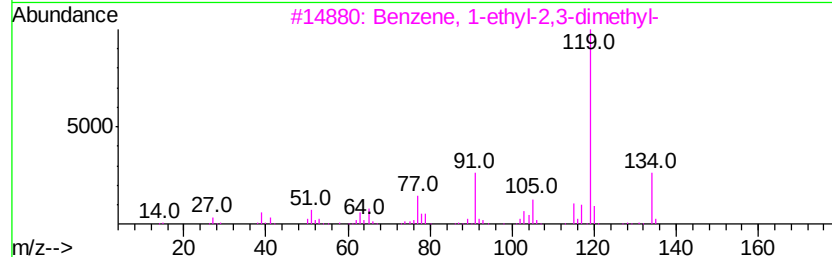
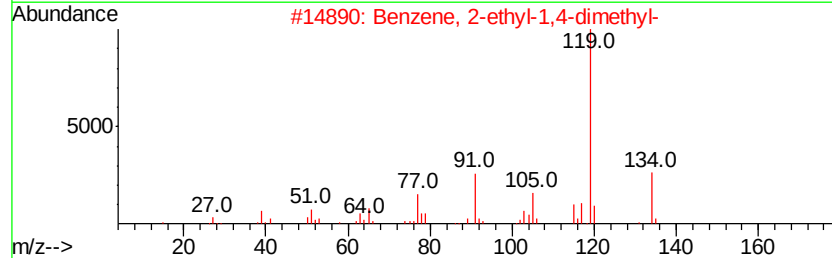
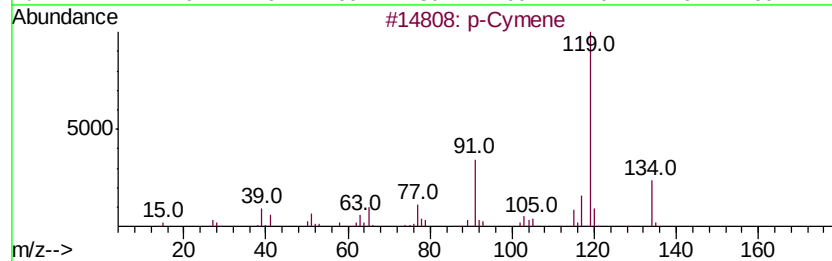
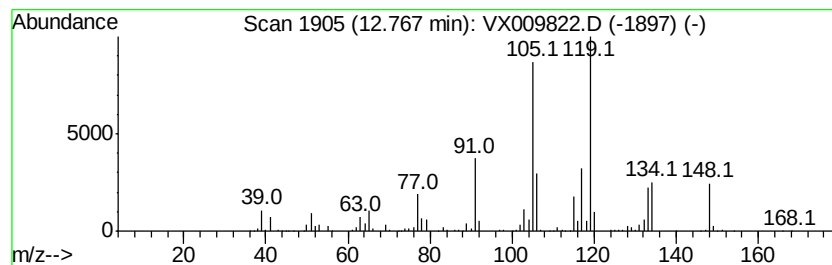
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MSVOA_X
ClientSampleId :
EP1-SB-E2-3R(14-17)

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 p-Cymene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.77	46.11 ug/l	1376220	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	p-Cymene	134 C10H14	000099-87-6	60
2	Benzene, 2-ethyl-1,4-dimethyl-	134 C10H14	001758-88-9	60
3	Benzene, 1-ethyl-2,3-dimethyl-	134 C10H14	000933-98-2	60
4	Benzene, 1,4-diethyl-	134 C10H14	000105-05-5	58
5	o-Cymene	134 C10H14	000527-84-4	55



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX052219\
Data File : VX009822.D
Acq On : 22 May 2019 23:21
Operator : JC/SP
Sample : K2976-01 50X
Misc : 6.56g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 34 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EP1-SB-E2-3R(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	69.4	ug/l	1037110	3	10.11	747280	50.0
Octane, 2,6-dimet...	10.83	142.7	ug/l	2132270	3	10.11	747280	50.0
Cyclohexane, propyl-	10.91	129.1	ug/l	1929190	3	10.11	747280	50.0
Heptane, 3-ethyl-...	10.94	49.4	ug/l	737994	3	10.11	747280	50.0
2-Octene, 2,6-dim...	11.27	45.6	ug/l	1360510	4	12.08	1492410	50.0
Benzene, 1-ethyl-...	11.45	46.0	ug/l	1372890	4	12.08	1492410	50.0
Benzene, 1,2,3-tr...	12.14	78.2	ug/l	2332950	4	12.08	1492410	50.0
Benzene, 1-methyl...	12.61	50.2	ug/l	1498410	4	12.08	1492410	50.0
Benzene, 2-ethyl-...	12.66	42.1	ug/l	1256140	4	12.08	1492410	50.0
p-Cymene	12.77	46.1	ug/l	1376220	4	12.08	1492410	50.0