

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX070319\
 Data File : VX010674.D
 Acq On : 04 Jul 2019 07:51
 Operator : JC/SP
 Sample : K3664-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FL-0582

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\82X061919W.M

Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.794	99	105	117	rVB	235846	347860	2.48%	0.362%
2	2.001	134	139	149	rVB	248132	349873	2.49%	0.364%
3	2.397	195	204	210	rBV	75344	157696	1.12%	0.164%
4	2.849	260	278	280	rBV	129406	273338	1.94%	0.284%
5	2.891	280	285	289	rVV	463970	1000067	7.12%	1.039%
6	2.934	289	292	304	rVB	248659	476675	3.39%	0.495%
7	3.178	320	332	346	rBV	466565	1088206	7.74%	1.131%
8	3.525	378	389	401	rBV	180880	416974	2.97%	0.433%
9	4.403	523	533	549	rVB	893254	2614272	18.60%	2.717%
10	5.501	699	713	718	rBV2	131840	382572	2.72%	0.398%
11	5.586	718	727	735	rVV	934721	3038978	21.62%	3.158%
12	5.659	735	739	746	rVV	238323	655635	4.67%	0.681%
13	5.726	747	750	766	rVB2	86212	224872	1.60%	0.234%
14	5.933	771	784	797	rVV4	192890	662810	4.72%	0.689%
15	6.061	797	805	811	rVV	127903	335346	2.39%	0.349%
16	6.147	811	819	829	rVV	482565	1276394	9.08%	1.327%
17	6.263	829	838	847	rVV2	154086	488255	3.47%	0.507%
18	6.360	847	854	862	rVV2	156582	429646	3.06%	0.447%
19	6.458	862	870	882	rVB	258762	715411	5.09%	0.744%
20	6.860	925	936	945	rBV	339339	797632	5.68%	0.829%
21	7.464	1022	1035	1047	rBV	2182213	4807208	34.21%	4.996%
22	7.732	1072	1079	1089	rVB	198558	387483	2.76%	0.403%
23	8.665	1225	1232	1235	rBV2	172132	298453	2.12%	0.310%
24	8.713	1235	1240	1248	rVB	742411	1225590	8.72%	1.274%
25	8.841	1254	1261	1265	rBV	100250	169681	1.21%	0.176%
26	9.036	1288	1293	1300	rVB	191921	320229	2.28%	0.333%
27	9.165	1305	1314	1320	rBV	94462	151518	1.08%	0.157%
28	9.622	1385	1389	1394	rVB	169802	229242	1.63%	0.238%
29	10.109	1464	1469	1476	rBV	726057	955888	6.80%	0.993%
30	10.183	1476	1481	1486	rVV	126958	184739	1.31%	0.192%
31	10.250	1486	1492	1498	rVV	4970400	6409470	45.61%	6.661%
32	10.353	1501	1509	1516	rVB	1806261	2483240	17.67%	2.581%
33	11.018	1612	1618	1629	rVB	1728391	2168349	15.43%	2.254%
34	11.134	1631	1637	1642	rBV	638366	831004	5.91%	0.864%

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Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
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 Stop Thrs : 0

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 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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 Title : SW846 8260

35	11.359	1669	1674	1680	rBV	2454240	2986667	21.25%	3.104%
36	11.432	1680	1686	1693	rVV	843011	1334948	9.50%	1.387%
37	11.506	1693	1698	1705	rVB	677895	839211	5.97%	0.872%
38	11.664	1715	1724	1734	rBV	509309	753095	5.36%	0.783%
39	11.810	1742	1748	1753	rVB	11438715	14053966	100.00%	14.606%
40	11.920	1758	1766	1767	rBV	119926	169724	1.21%	0.176%
41	11.944	1767	1770	1778	rVB	389383	475485	3.38%	0.494%
42	12.024	1778	1783	1786	rBV	132004	158651	1.13%	0.165%
43	12.073	1786	1791	1797	rVB2	843280	1211823	8.62%	1.259%
44	12.140	1797	1802	1808	rBV	6296994	7800029	55.50%	8.106%
45	12.231	1812	1817	1821	rBV	187808	226502	1.61%	0.235%
46	12.298	1821	1828	1835	rBV	1513718	2047151	14.57%	2.128%
47	12.365	1835	1839	1849	rVB3	707160	1193227	8.49%	1.240%
48	12.457	1849	1854	1859	rBV	426100	525843	3.74%	0.546%
49	12.518	1859	1864	1870	rBV	1108560	1324459	9.42%	1.376%
50	12.585	1870	1875	1877	rBV	946947	1197849	8.52%	1.245%
51	12.609	1877	1879	1883	rVB	842025	855306	6.09%	0.889%
52	12.664	1883	1888	1892	rBV	1327376	1622109	11.54%	1.686%
53	12.700	1892	1894	1898	rVB	361703	366277	2.61%	0.381%
54	12.755	1898	1903	1910	rBV2	927000	1569710	11.17%	1.631%
55	12.902	1922	1927	1932	rVB2	959491	1193906	8.50%	1.241%
56	12.975	1932	1939	1942	rBV	887008	1099258	7.82%	1.142%
57	13.017	1942	1946	1952	rVB	1572730	1840510	13.10%	1.913%
58	13.078	1952	1956	1961	rBV2	125312	217164	1.55%	0.226%
59	13.194	1971	1975	1977	rBV2	167104	210541	1.50%	0.219%
60	13.225	1977	1980	1983	rVB	304641	337790	2.40%	0.351%
61	13.304	1989	1993	1995	rBV2	143929	185492	1.32%	0.193%
62	13.341	1995	1999	2005	rVB2	2532544	3471406	24.70%	3.608%
63	13.475	2017	2021	2027	rVB	1119502	1432417	10.19%	1.489%
64	13.560	2027	2035	2038	rBV3	99204	153496	1.09%	0.160%
65	13.597	2038	2041	2044	rBV	118416	142447	1.01%	0.148%
66	13.639	2044	2048	2053	rVV3	237297	418217	2.98%	0.435%
67	13.719	2058	2061	2068	rVB2	271806	435896	3.10%	0.453%
68	13.834	2071	2080	2088	rVB	1615385	2276720	16.20%	2.366%
69	13.908	2088	2092	2097	rBV	146489	191313	1.36%	0.199%
70	13.981	2097	2104	2108	rBV2	232409	340686	2.42%	0.354%
71	14.286	2146	2154	2159	rVB2	331030	579014	4.12%	0.602%

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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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72	14.535	2190	2195	2200	rBV2	352880	447810	3.19%	0.465%
73	14.694	2208	2221	2226	rBV	1710667	2144953	15.26%	2.229%
74	14.834	2239	2244	2253	rBV	1142376	1480758	10.54%	1.539%
75	15.267	2305	2315	2320	rBV2	71351	145346	1.03%	0.151%
76	15.541	2354	2360	2365	rBV	99649	169095	1.20%	0.176%
77	15.682	2374	2383	2387	rBV	129429	240133	1.71%	0.250%

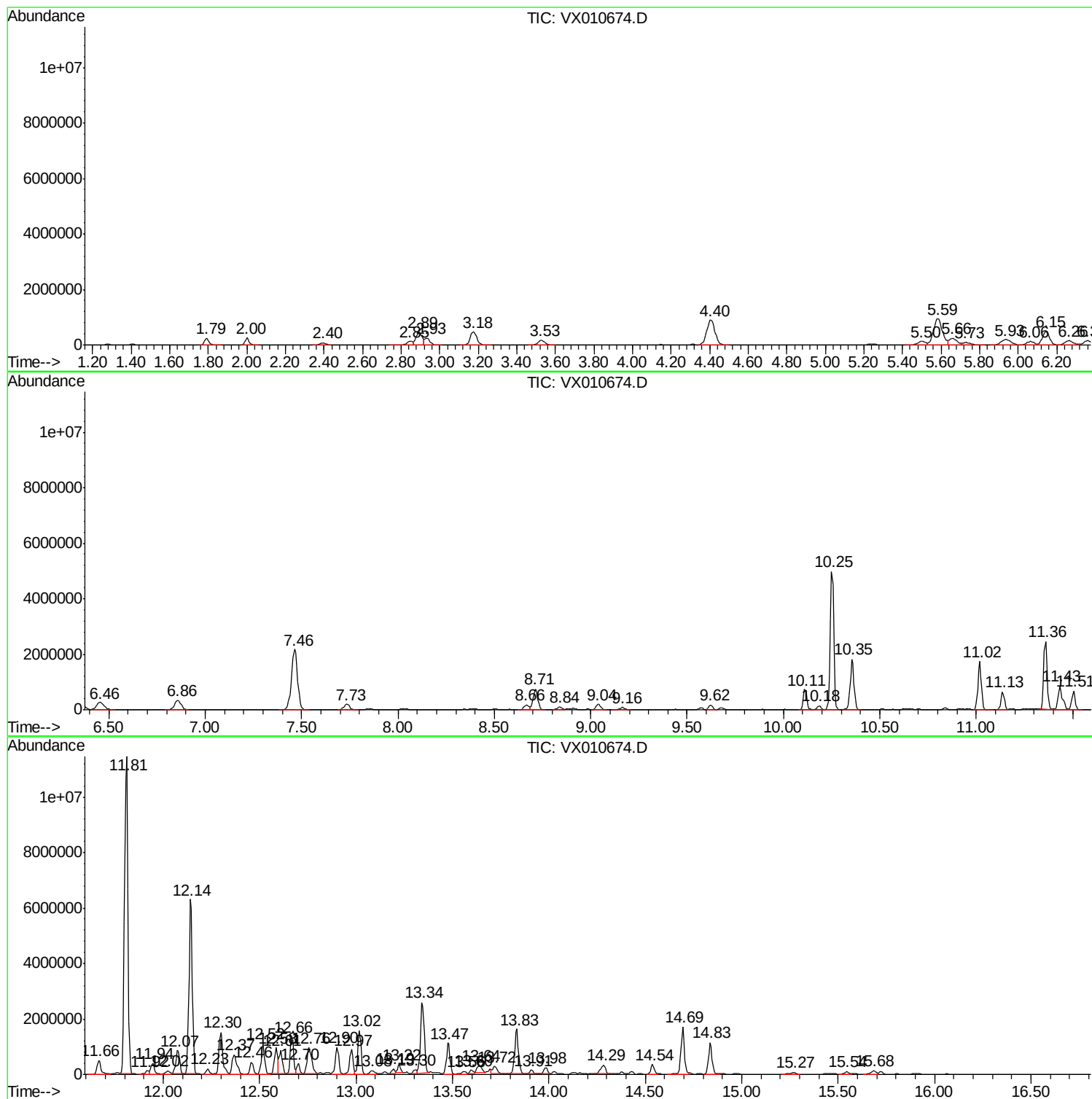
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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



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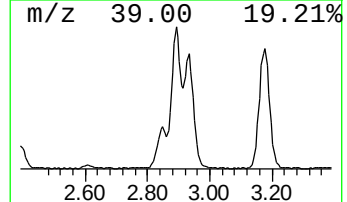
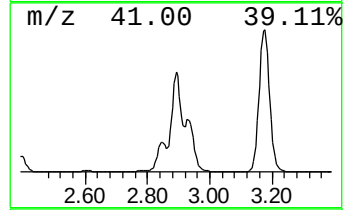
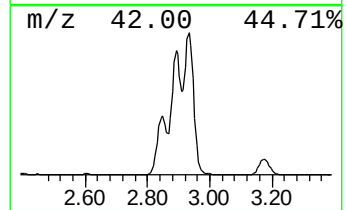
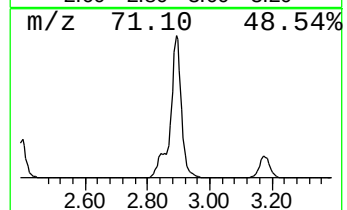
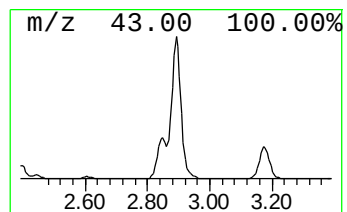
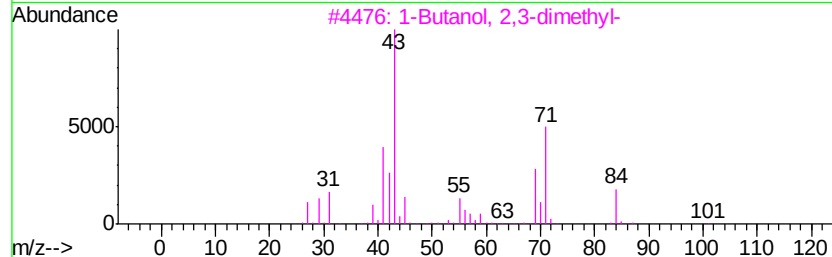
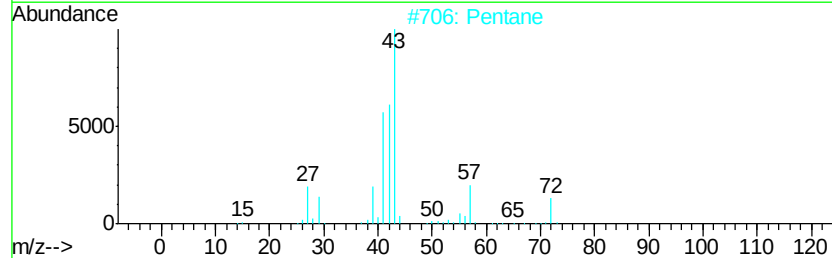
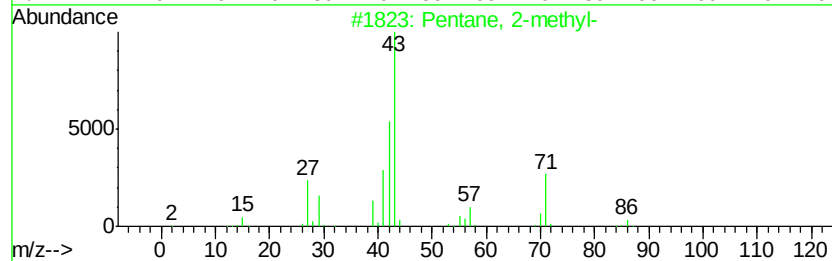
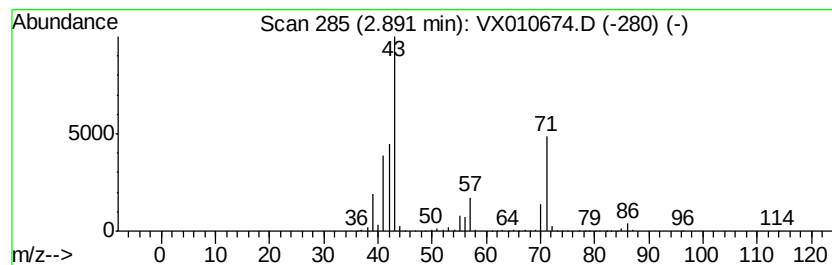
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

Peak Number 1 Pentane, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.89	76.27 ug/l	1000070	Pentafluorobenzene	5.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2-methyl-	86	C6H14	000107-83-5	72
2		Pentane	72	C5H12	000109-66-0	46
3		1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	38
4		Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	38
5		Pentane, 3,3-dimethyl-	100	C7H16	000562-49-2	38



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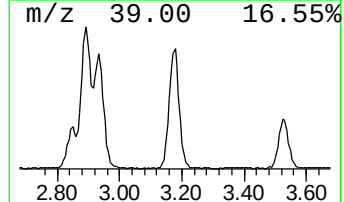
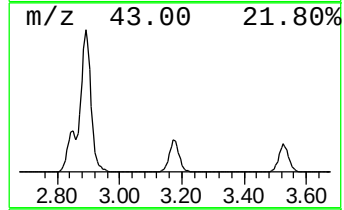
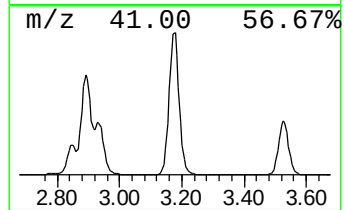
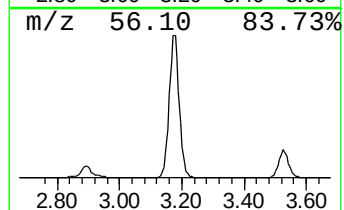
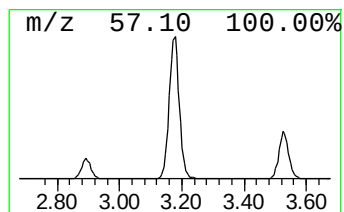
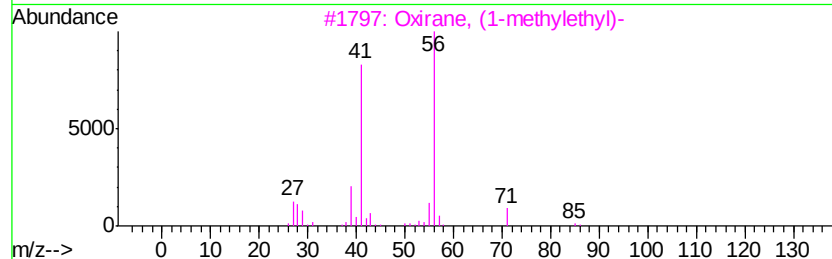
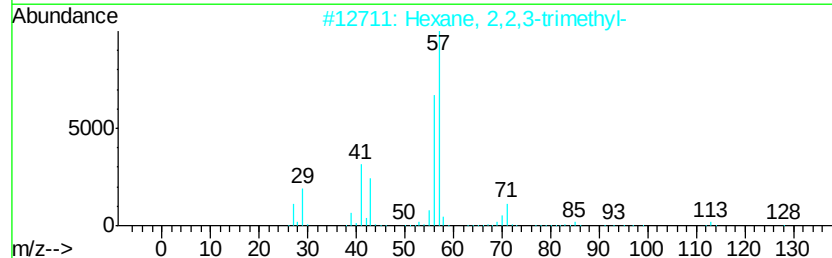
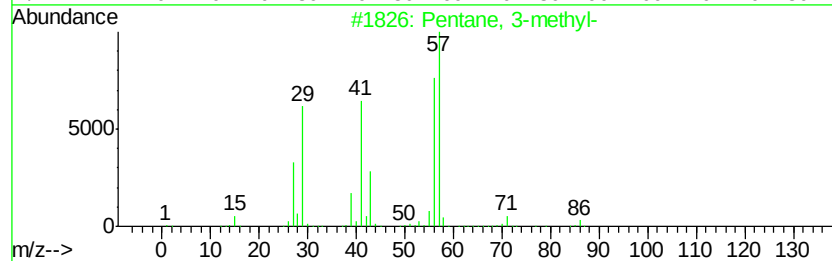
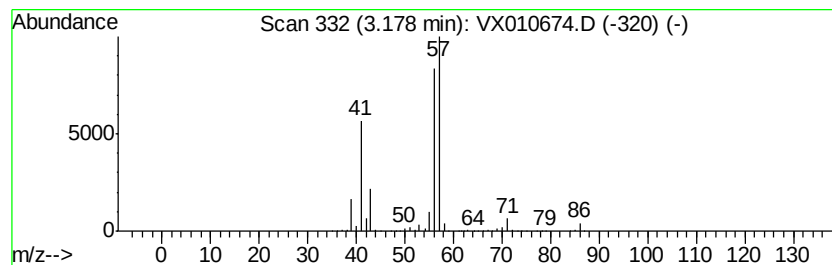
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

Peak Number 2 Pentane, 3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.18	82.99 ug/l	1088210	Pentafluorobenzene	5.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2		Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	78
3		Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	45
4		Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	43
5		Sulphuric acid dibutyl ester	210	C8H18O4S	000625-22-9	40



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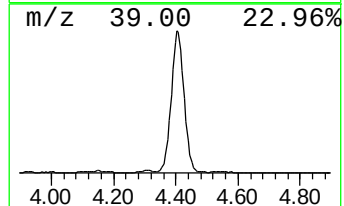
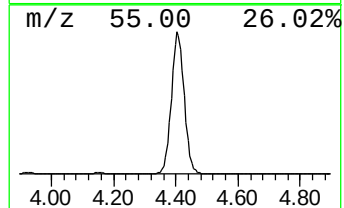
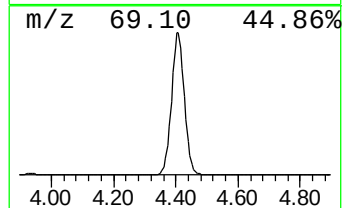
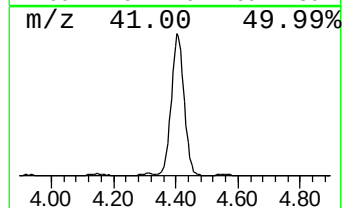
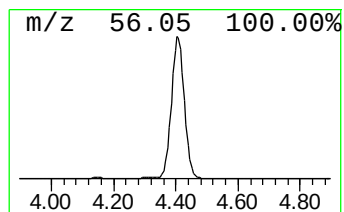
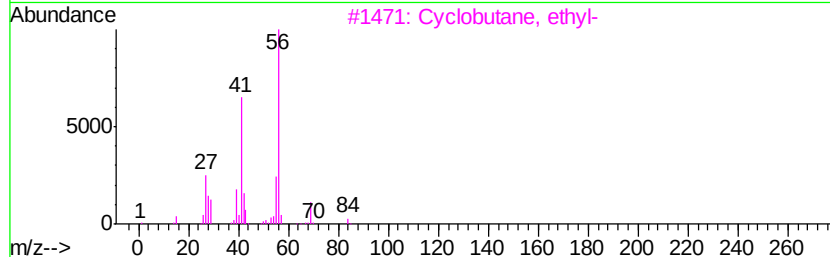
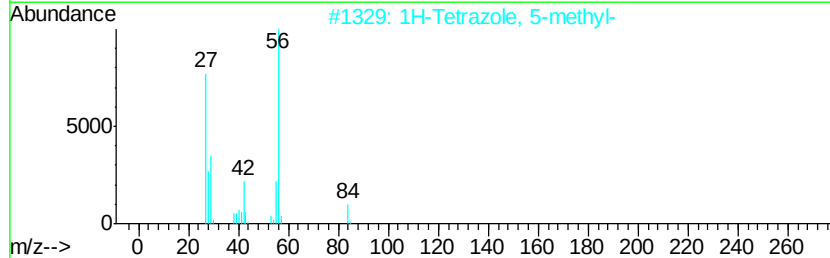
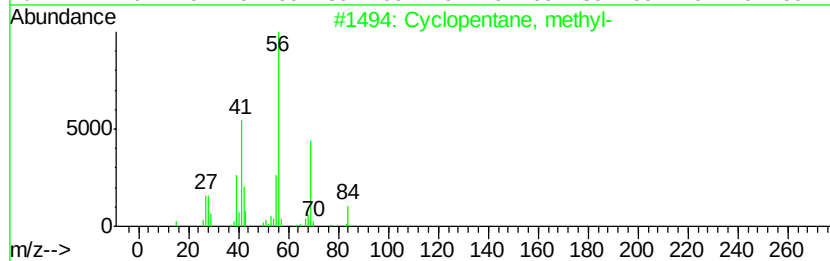
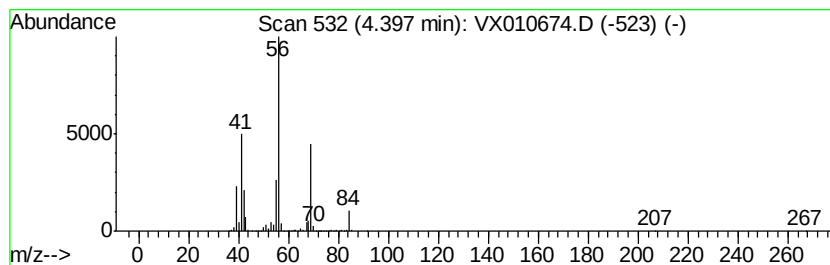
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Quant Title : SW846 8260

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

Peak Number 3 Cyclopentane, methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.40	199.37 ug/l	2614270	Pentafluorobenzene	5.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
3		Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
4		Cyclohexane	84	C6H12	000110-82-7	64
5		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	49



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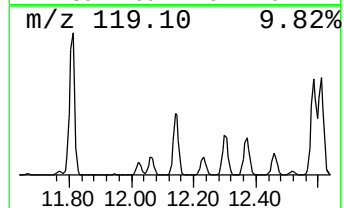
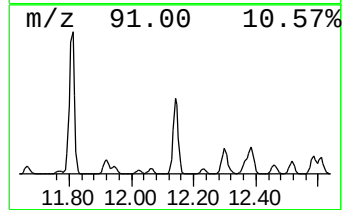
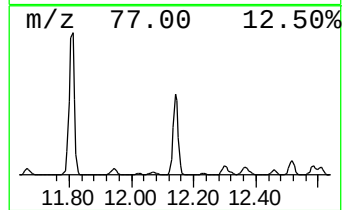
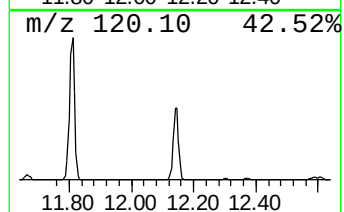
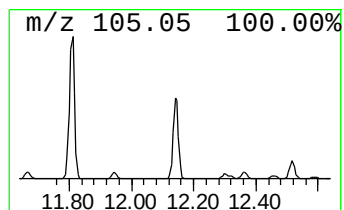
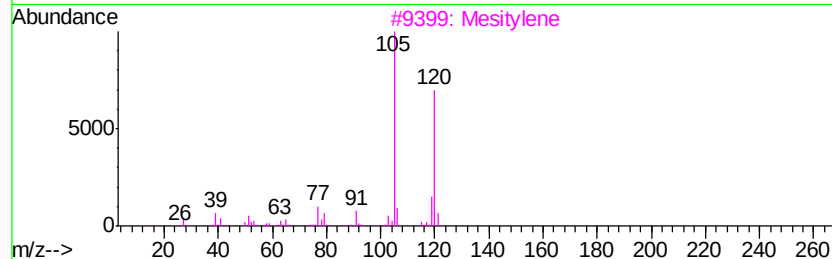
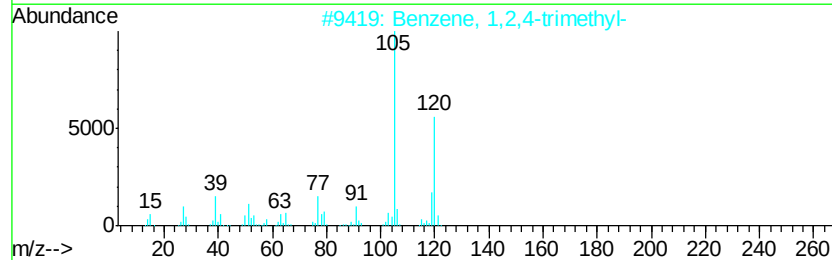
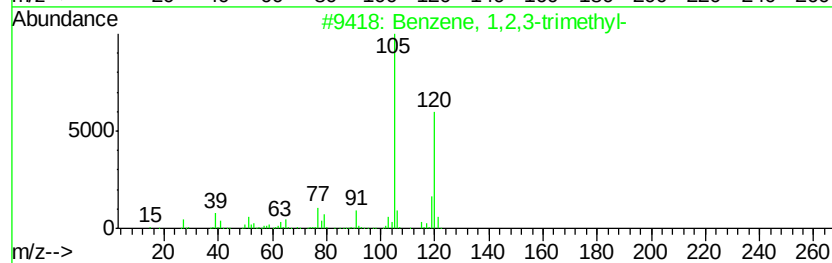
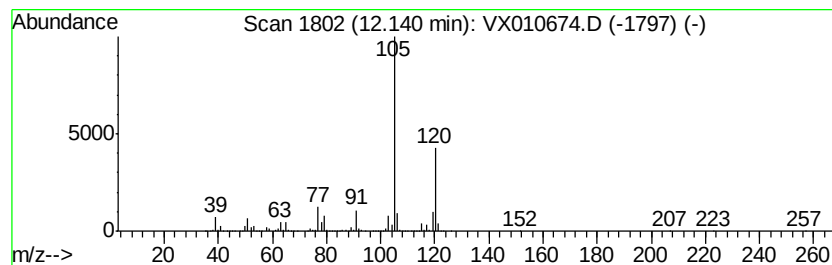
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1,2,3-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	321.83 ug/l	7800030	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
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\VX070319\
Data File : VX010674.D
Acq On : 04 Jul 2019 07:51
Operator : JC/SP
Sample : K3664-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 45 Sample Multiplier: 1

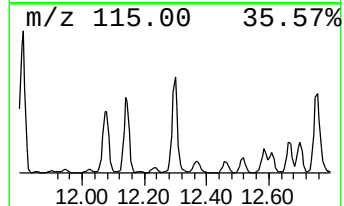
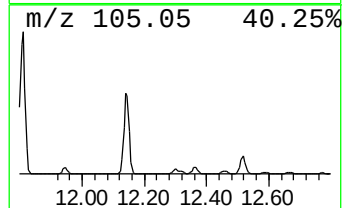
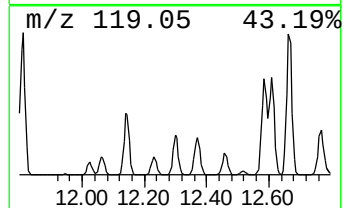
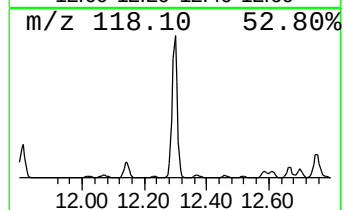
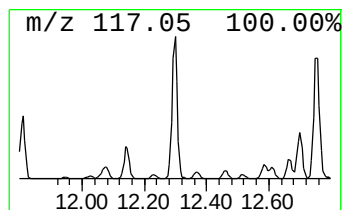
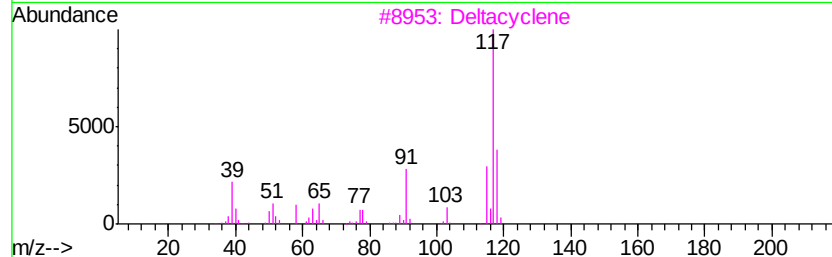
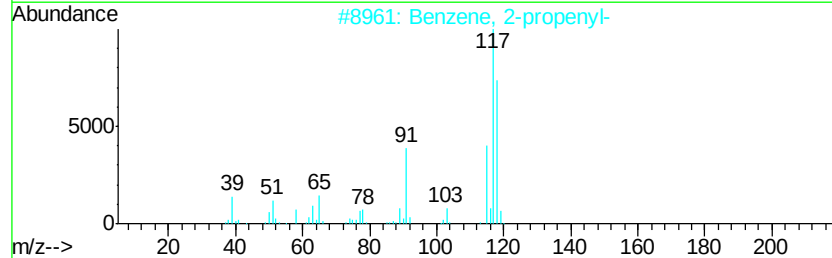
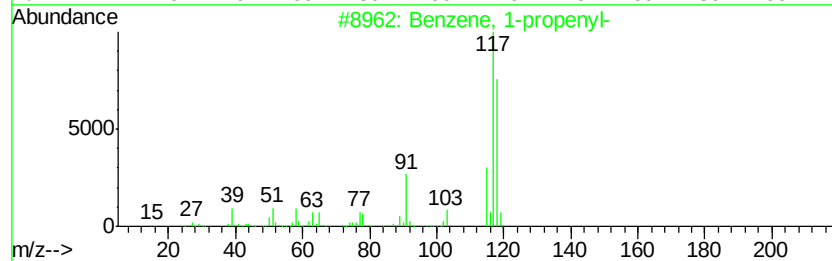
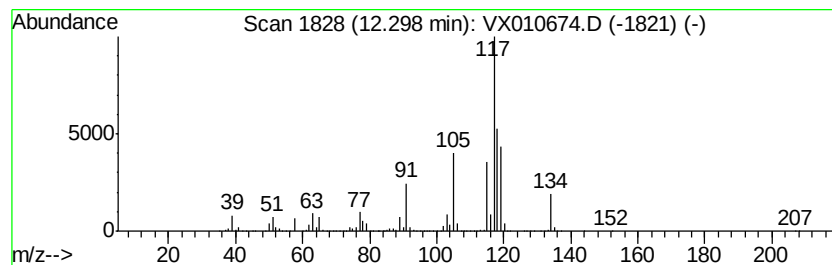
Instrument :
MSVOA_X
ClientSampleId :
FL-0582

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-propenyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	84.47 ug/l	2047150	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-propenyl-	118 C9H10	000637-50-3 90
2		Benzene, 2-propenyl-	118 C9H10	000300-57-2 55
3		Deltacyclene	118 C9H10	007785-10-6 55
4		Benzene, 1-ethenyl-2-methyl-	118 C9H10	000611-15-4 55
5		Indane	118 C9H10	000496-11-7 55



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX070319\
Data File : VX010674.D
Acq On : 04 Jul 2019 07:51
Operator : JC/SP
Sample : K3664-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 45 Sample Multiplier: 1

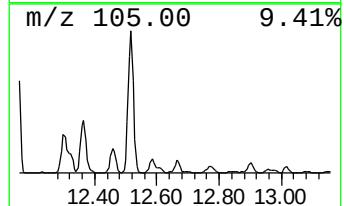
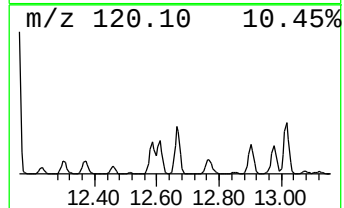
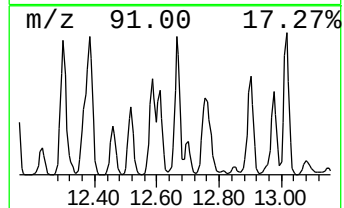
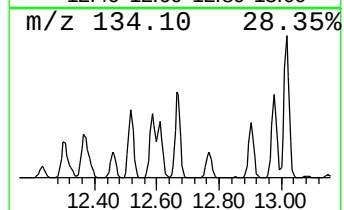
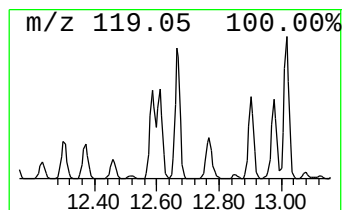
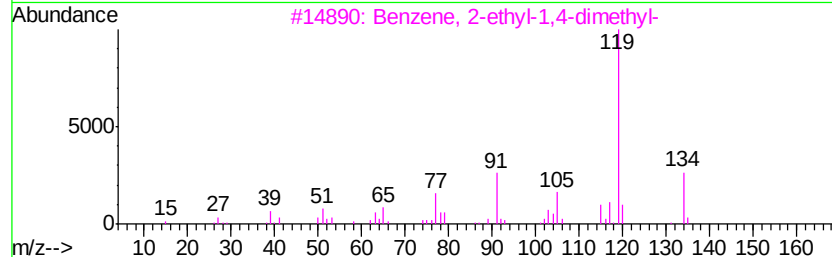
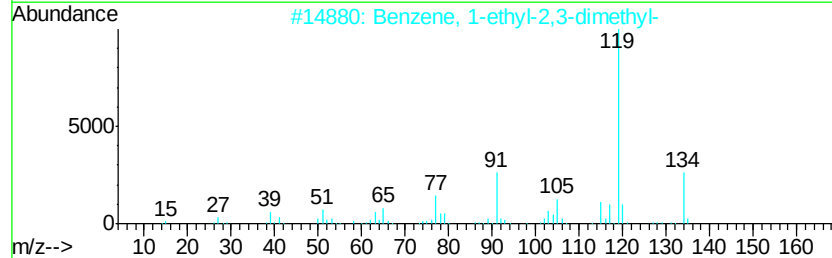
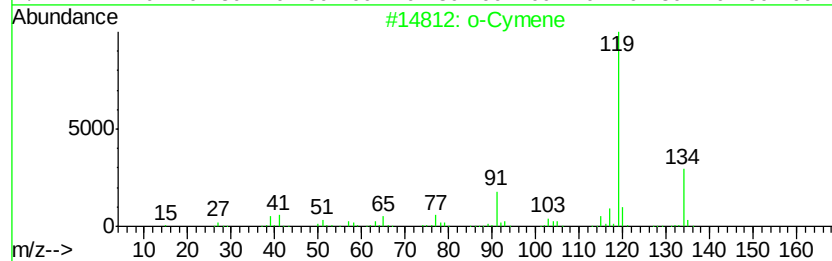
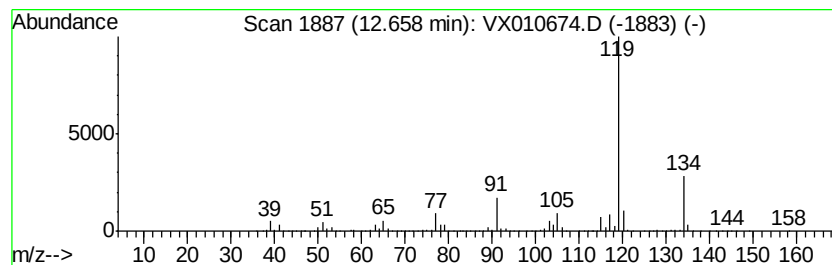
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ClientSampleId :
FL-0582

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

Peak Number 6 o-Cymene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.66	66.93 ug/l	1622110	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		o-Cymene	134 C10H14	000527-84-4 97
2		Benzene, 1-ethyl-2,3-dimethyl-	134 C10H14	000933-98-2 96
3		Benzene, 2-ethyl-1,4-dimethyl-	134 C10H14	001758-88-9 96
4		Benzene, 2-ethyl-1,3-dimethyl-	134 C10H14	002870-04-4 95
5		Benzene, 1-ethyl-2,4-dimethyl-	134 C10H14	000874-41-9 95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX070319\
Data File : VX010674.D
Acq On : 04 Jul 2019 07:51
Operator : JC/SP
Sample : K3664-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 45 Sample Multiplier: 1

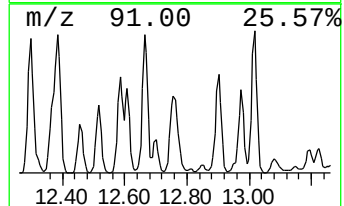
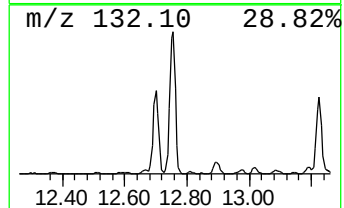
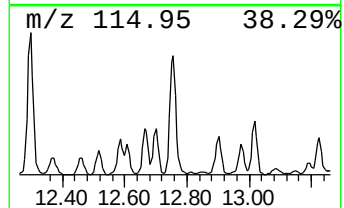
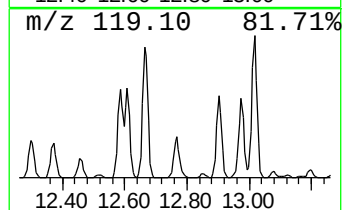
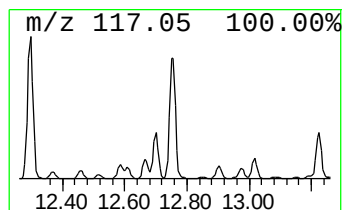
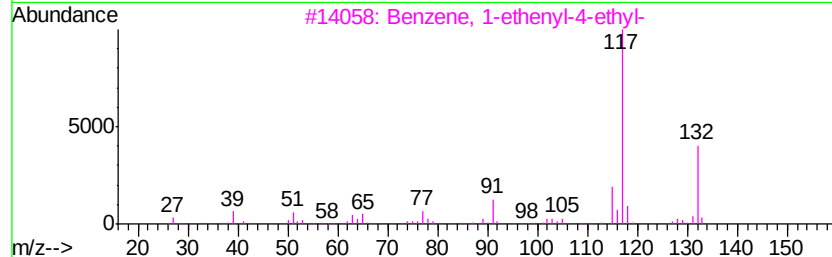
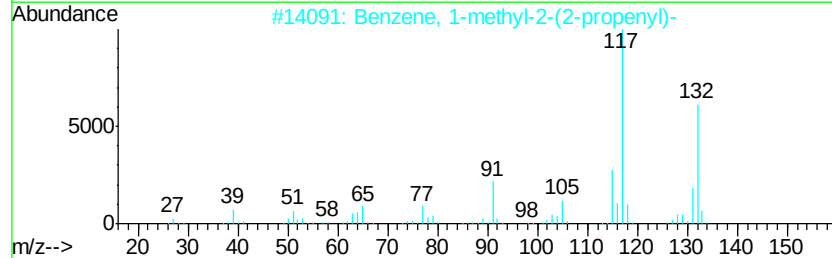
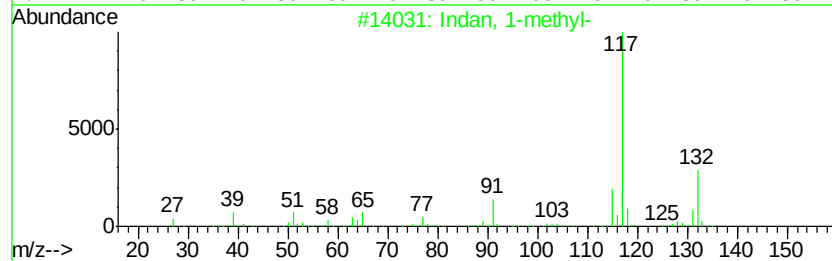
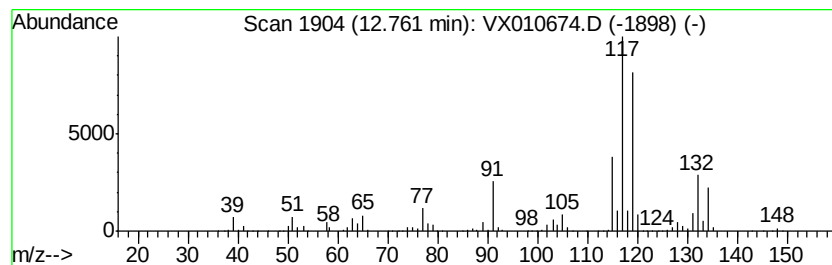
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ClientSampleId :
FL-0582

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

Peak Number 7 Indan, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	64.77 ug/l	1569710	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Indan, 1-methyl-	132 C10H12	000767-58-8	93
2	Benzene, 1-methyl-2-(2-propenyl)-	132 C10H12	001587-04-8	81
3	Benzene, 1-ethenyl-4-ethyl-	132 C10H12	003454-07-7	76
4	Benzene, 2-butenyl-	132 C10H12	001560-06-1	74
5	Benzene, (2-methyl-1-propenyl)-	132 C10H12	000768-49-0	74



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX070319\
Data File : VX010674.D
Acq On : 04 Jul 2019 07:51
Operator : JC/SP
Sample : K3664-01
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ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0582

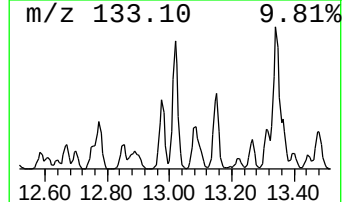
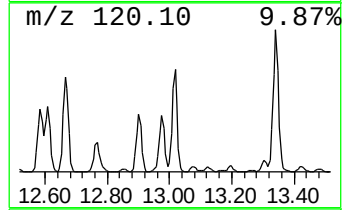
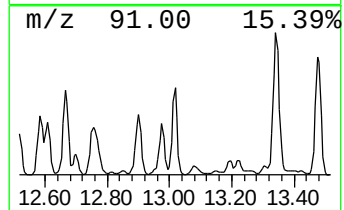
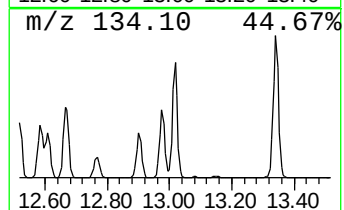
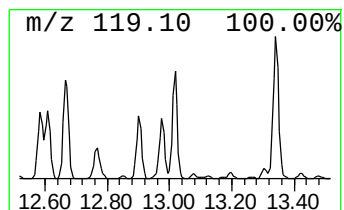
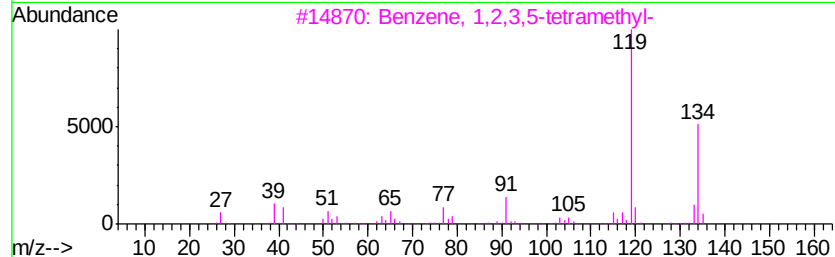
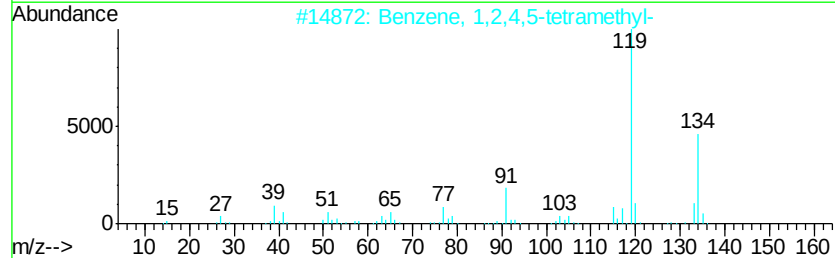
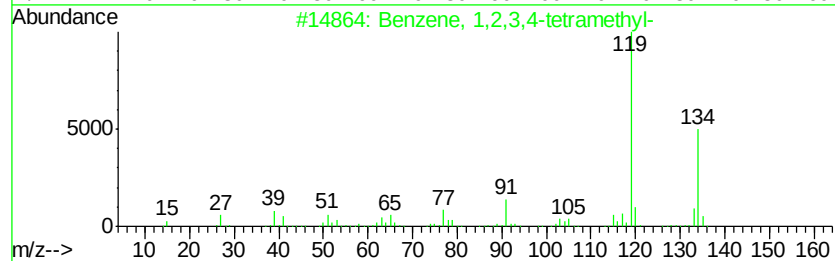
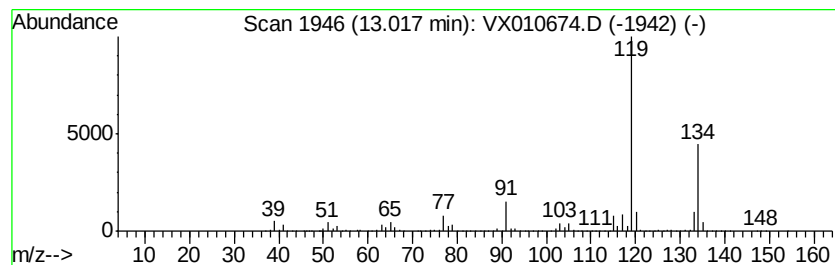
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	75.94 ug/l	1840510	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4		o-Cymene	134	C10H14	000527-84-4	95
5		1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX070319\
Data File : VX010674.D
Acq On : 04 Jul 2019 07:51
Operator : JC/SP
Sample : K3664-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
FL-0582

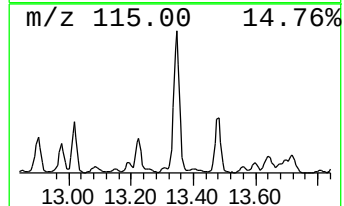
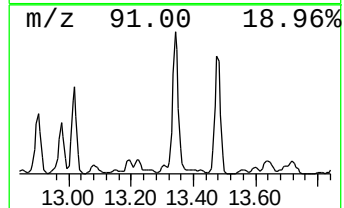
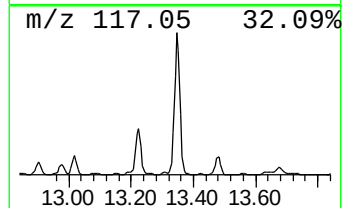
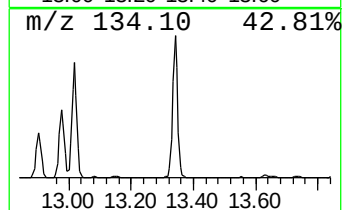
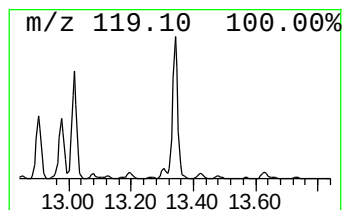
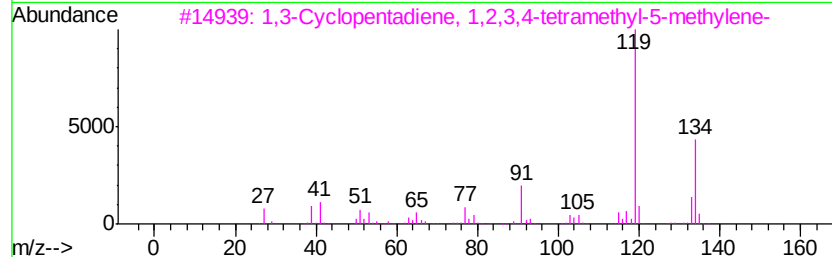
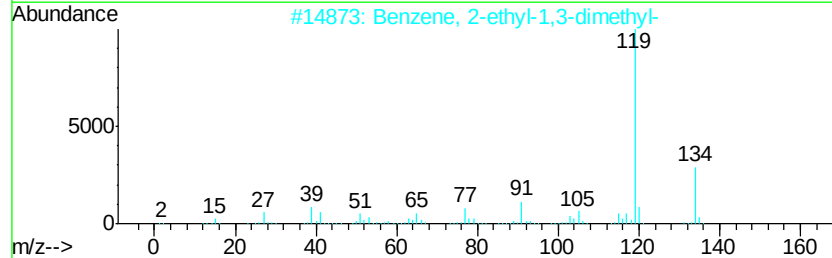
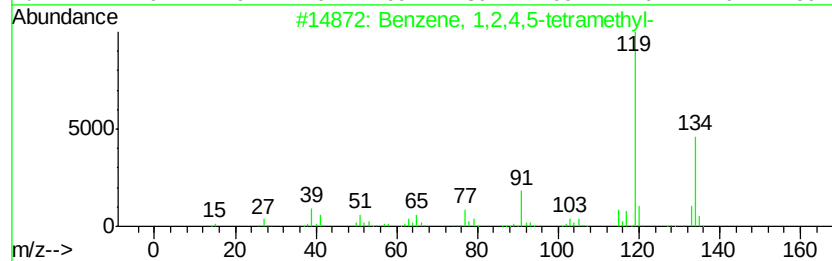
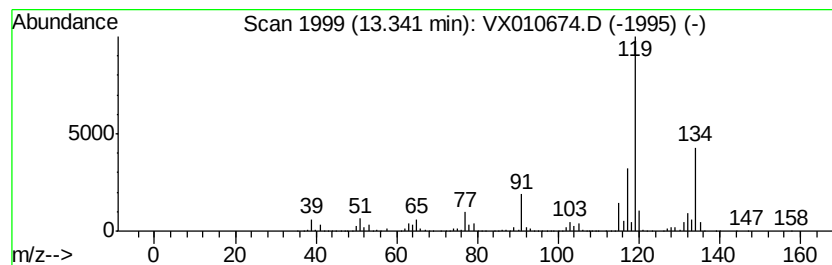
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.34	143.23 ug/l	3471410	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	89
2		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	81
3		1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	81
4		o-Cymene	134	C10H14	000527-84-4	81
5		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	81



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX070319\
Data File : VX010674.D
Acq On : 04 Jul 2019 07:51
Operator : JC/SP
Sample : K3664-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 45 Sample Multiplier: 1

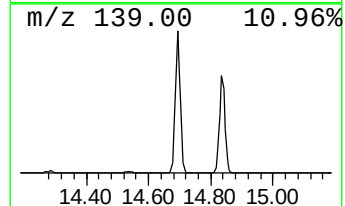
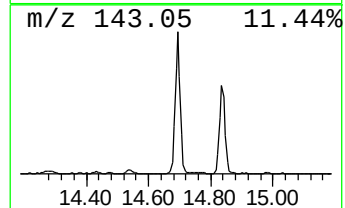
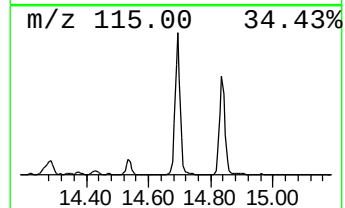
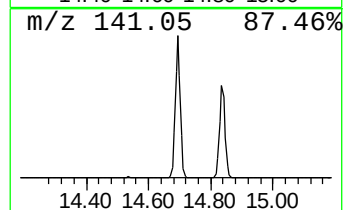
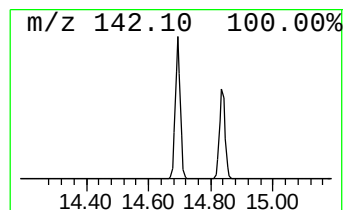
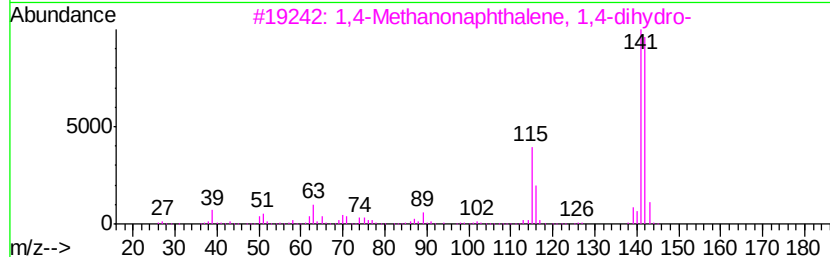
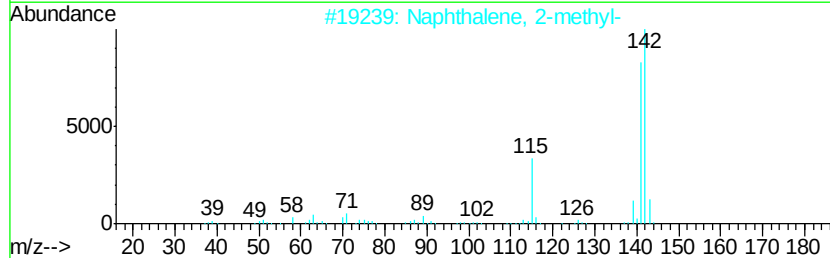
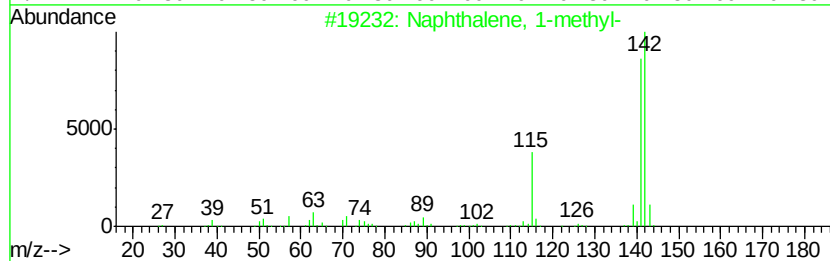
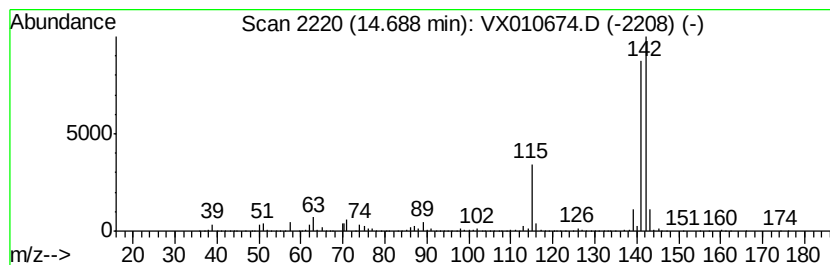
Instrument :
MSVOA_X
ClientSampleId :
FL-0582

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	88.50 ug/l	2144950	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1-methyl-	142 C11H10	000090-12-0 96
2		Naphthalene, 2-methyl-	142 C11H10	000091-57-6 96
3		1,4-Methanonaphthalene, 1,4-dihv...	142 C11H10	004453-90-1 91
4		Benzocycloheptatriene	142 C11H10	000264-09-5 91
5		1H-Indene, 1-ethylidene-	142 C11H10	002471-83-2 64



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070319\
Data File : VX010674.D
Acq On : 04 Jul 2019 07:51
Operator : JC/SP
Sample : K3664-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0582

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Pentane, 3-methyl-	3.18	83.0	ug/l	1088210	1	5.66	655635	50.0
Cyclopentane, met...	4.40	199.4	ug/l	2614270	1	5.66	655635	50.0
Benzene, 1,2,3-tr...	12.14	321.8	ug/l	7800030	4	12.08	1211820	50.0
Benzene, 1-propenyl-	12.30	84.5	ug/l	2047150	4	12.08	1211820	50.0
o-Cymene	12.66	66.9	ug/l	1622110	4	12.08	1211820	50.0
Indan, 1-methyl-	12.76	64.8	ug/l	1569710	4	12.08	1211820	50.0
Benzene, 1,2,3,4-...	13.02	75.9	ug/l	1840510	4	12.08	1211820	50.0
Benzene, 1,2,4,5-...	13.34	143.2	ug/l	3471410	4	12.08	1211820	50.0
Naphthalene, 1-me...	14.69	88.5	ug/l	2144950	4	12.08	1211820	50.0