

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX082019\
 Data File : VX011678.D
 Acq On : 20 Aug 2019 12:41
 Operator : JC/SP
 Sample : K4391-04
 Misc : 5.06g/10mL/100uL/10mL/MSVOA_X/MEOH
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BOTTOM-1

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\82X081519W.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	1.696	62	89	90	rBV	59358	326910	17.36%	1.308%
2	5.488	700	711	725	rBV	175231	532258	28.27%	2.130%
3	5.647	726	737	750	rVV	344911	966750	51.35%	3.868%
4	6.049	794	803	814	rBV	186542	489234	25.98%	1.958%
5	6.439	860	867	874	rVB3	19986	52760	2.80%	0.211%
6	6.848	924	934	946	rBV	498391	1154879	61.34%	4.621%
7	7.451	1023	1033	1042	rVB	236374	552918	29.37%	2.212%
8	7.561	1046	1051	1057	rVV2	23278	47242	2.51%	0.189%
9	7.628	1057	1062	1068	rVB2	32520	62822	3.34%	0.251%
10	7.842	1087	1097	1105	rBV2	49140	107048	5.69%	0.428%
11	8.018	1120	1126	1133	rBV	45754	93485	4.97%	0.374%
12	8.250	1159	1164	1167	rBV	29594	47135	2.50%	0.189%
13	8.335	1173	1178	1181	rBV	49744	76728	4.08%	0.307%
14	8.378	1181	1185	1191	rVB3	36733	67148	3.57%	0.269%
15	8.494	1200	1204	1207	rVV2	46861	76719	4.07%	0.307%
16	8.524	1207	1209	1214	rVV	36046	59065	3.14%	0.236%
17	8.658	1224	1231	1234	rBV	311885	532183	28.27%	2.129%
18	8.707	1234	1239	1253	rVV	1022630	1775502	94.30%	7.104%
19	8.829	1253	1259	1264	rVV3	76878	131311	6.97%	0.525%
20	8.902	1268	1271	1276	rVB2	44161	70365	3.74%	0.282%
21	9.036	1287	1293	1299	rBV2	209333	335262	17.81%	1.341%
22	9.158	1304	1313	1319	rBV	116471	189354	10.06%	0.758%
23	9.286	1330	1334	1338	rVV	33571	47252	2.51%	0.189%
24	9.347	1338	1344	1349	rVB	88737	132523	7.04%	0.530%
25	9.439	1349	1359	1368	rBV	205744	342384	18.18%	1.370%
26	9.555	1373	1378	1384	rBV4	238296	479996	25.49%	1.921%
27	9.616	1384	1388	1393	rVB	279415	405875	21.56%	1.624%
28	9.677	1393	1398	1402	rBV	226033	368886	19.59%	1.476%
29	9.713	1402	1404	1408	rVB	40818	44419	2.36%	0.178%
30	9.811	1414	1420	1425	rBV3	52118	87662	4.66%	0.351%
31	9.884	1425	1432	1436	rBV2	204684	466778	24.79%	1.868%
32	9.957	1440	1444	1448	rVB	191484	272623	14.48%	1.091%
33	10.061	1454	1461	1465	rBV	229599	330359	17.55%	1.322%
34	10.109	1465	1469	1474	rBV	994697	1304056	69.26%	5.218%

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35	10.183	1474	1481	1484	rVB5	34366	78083	4.15%	0.312%
36	10.231	1484	1489	1496	rVB4	59996	115813	6.15%	0.463%
37	10.329	1496	1505	1508	rBV3	119995	236596	12.57%	0.947%
38	10.372	1508	1512	1515	rVB	228454	282279	14.99%	1.129%
39	10.408	1515	1518	1530	rVB2	151198	262216	13.93%	1.049%
40	10.506	1530	1534	1539	rBV	67297	100683	5.35%	0.403%
41	10.585	1543	1547	1549	rBV2	48518	65820	3.50%	0.263%
42	10.640	1549	1556	1563	rBV4	325655	519681	27.60%	2.079%
43	10.719	1563	1569	1574	rBV3	171327	270636	14.37%	1.083%
44	10.762	1574	1576	1579	rVB	71750	79003	4.20%	0.316%
45	10.835	1579	1588	1591	rBV2	413303	654986	34.79%	2.621%
46	10.878	1591	1595	1596	rBV2	64702	69628	3.70%	0.279%
47	10.908	1596	1600	1604	rBV2	282517	342583	18.20%	1.371%
48	10.945	1604	1606	1611	rVB	189345	225931	12.00%	0.904%
49	11.000	1611	1615	1620	rVB4	64170	111198	5.91%	0.445%
50	11.079	1620	1628	1632	rBV3	245825	455641	24.20%	1.823%
51	11.134	1632	1637	1647	rVB	1140423	1882798	100.00%	7.533%
52	11.213	1647	1650	1651	rBV	50688	46203	2.45%	0.185%
53	11.243	1651	1655	1657	rBV	194320	222533	11.82%	0.890%
54	11.274	1657	1660	1663	rVB3	112824	141961	7.54%	0.568%
55	11.353	1669	1673	1676	rBV	36003	51690	2.75%	0.207%
56	11.390	1676	1679	1683	rBV2	40847	57649	3.06%	0.231%
57	11.438	1683	1687	1691	rBV3	58736	86329	4.59%	0.345%
58	11.481	1691	1694	1697	rBV	152418	234652	12.46%	0.939%
59	11.682	1724	1727	1732	rVV3	100634	140403	7.46%	0.562%
60	11.731	1732	1735	1738	rVV2	88816	117832	6.26%	0.471%
61	11.768	1738	1741	1744	rVV	199938	243047	12.91%	0.972%
62	11.804	1744	1747	1757	rVB2	236994	416330	22.11%	1.666%
63	11.890	1757	1761	1763	rBV2	73622	104481	5.55%	0.418%
64	11.920	1763	1766	1767	rVV3	41180	52056	2.76%	0.208%
65	11.944	1767	1770	1774	rVB3	81204	113101	6.01%	0.453%
66	11.993	1774	1778	1781	rBV	128352	170582	9.06%	0.683%
67	12.024	1781	1783	1786	rVV3	44924	51787	2.75%	0.207%
68	12.079	1787	1792	1797	rVV	1181811	1581391	83.99%	6.327%
69	12.121	1797	1799	1801	rVV	106978	122561	6.51%	0.490%
70	12.146	1801	1803	1808	rVB3	148044	201690	10.71%	0.807%
71	12.225	1812	1816	1824	rVB2	120249	175969	9.35%	0.704%

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72	12.298	1824	1828	1834	rBV6	33768	75359	4.00%	0.302%
73	12.365	1834	1839	1846	rVB3	116374	193930	10.30%	0.776%
74	12.469	1853	1856	1861	rBV	52988	70720	3.76%	0.283%
75	12.518	1861	1864	1870	rVB4	63034	103240	5.48%	0.413%
76	12.609	1872	1879	1881	rBV2	59163	124383	6.61%	0.498%
77	12.737	1896	1900	1903	rBV	42484	61127	3.25%	0.245%
78	12.767	1903	1905	1911	rVB4	27182	47955	2.55%	0.192%
79	12.871	1919	1922	1926	rVB2	44258	49194	2.61%	0.197%
80	12.944	1929	1934	1941	rVV4	85080	216453	11.50%	0.866%
81	13.024	1943	1947	1953	rVV4	58596	124451	6.61%	0.498%
82	13.091	1954	1958	1961	rVV3	55847	73066	3.88%	0.292%
83	13.347	1996	2000	2005	rVB2	91404	139647	7.42%	0.559%
84	13.432	2011	2014	2018	rVV2	72660	106681	5.67%	0.427%
85	13.475	2018	2021	2030	rVB7	56095	139849	7.43%	0.560%
86	13.633	2044	2047	2054	rBV5	23717	49548	2.63%	0.198%
87	13.731	2058	2063	2065	rBV4	30839	46543	2.47%	0.186%
88	13.792	2069	2073	2076	rVB3	38403	46873	2.49%	0.188%
89	13.895	2085	2090	2096	rVB2	75139	115095	6.11%	0.461%
90	14.267	2148	2151	2163	rVB5	38053	70403	3.74%	0.282%
91	14.529	2190	2194	2199	rVV5	25445	61474	3.27%	0.246%
92	14.663	2213	2216	2218	rBV2	89793	107142	5.69%	0.429%
93	14.694	2218	2221	2225	rVB	121527	140902	7.48%	0.564%
94	14.840	2241	2245	2252	rBV2	71354	99703	5.30%	0.399%
95	15.273	2313	2316	2319	rVV	80735	100799	5.35%	0.403%
96	15.438	2339	2343	2347	rBV	45748	68199	3.62%	0.273%
97	15.541	2356	2360	2365	rVB	88541	136425	7.25%	0.546%
98	15.682	2379	2383	2387	rBV	103912	165125	8.77%	0.661%
99	15.718	2387	2389	2398	rVB	63682	100450	5.34%	0.402%
100	16.438	2503	2507	2514	rVB	22213	42007	2.23%	0.168%

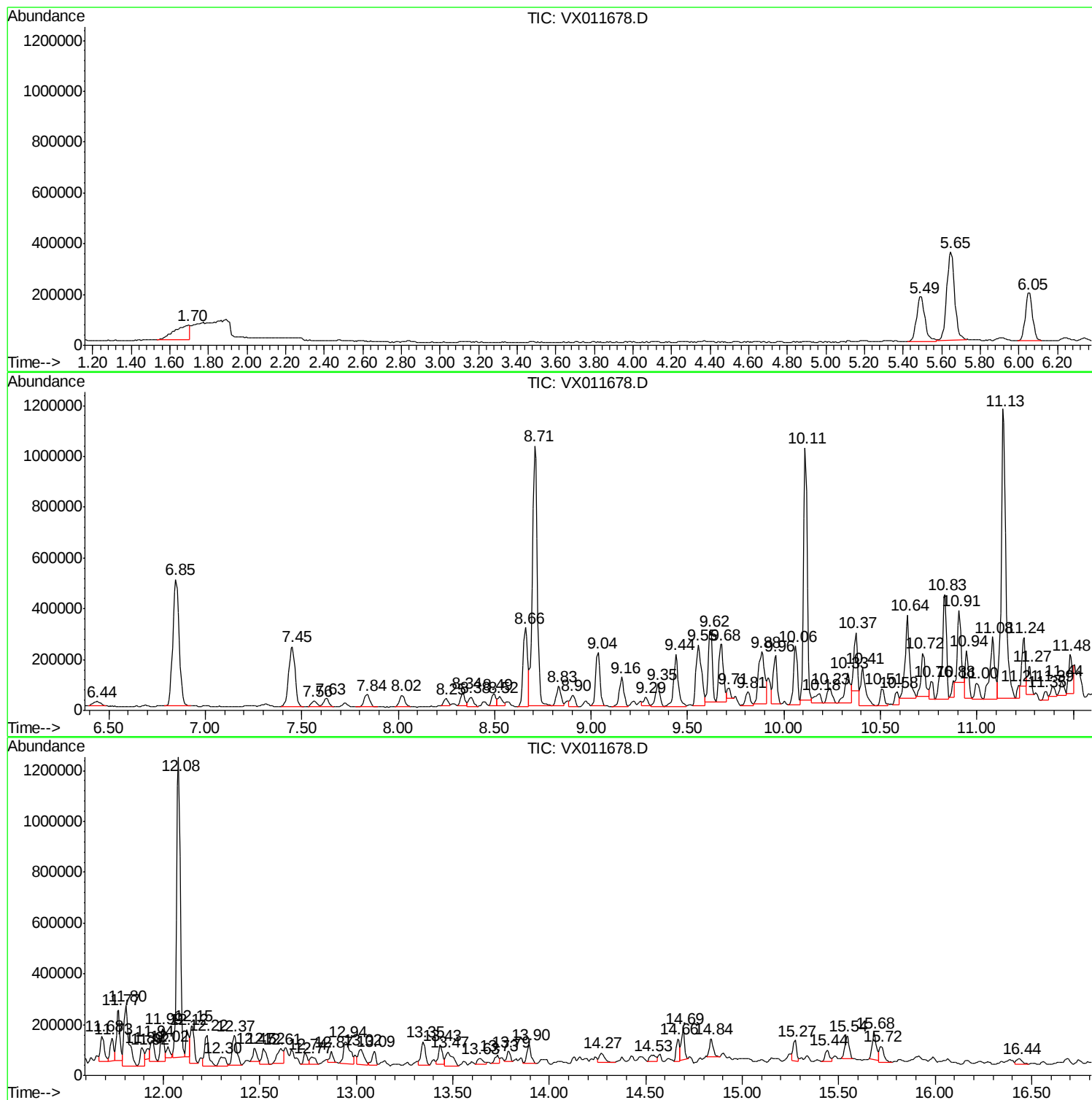
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Instrument :
MSVOA_X
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BOTTOM-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X081519W.M
Quant Title : SW846 8260

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



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Instrument :
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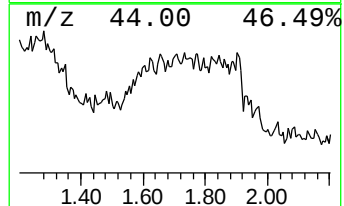
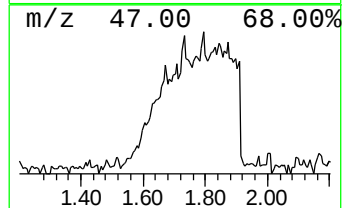
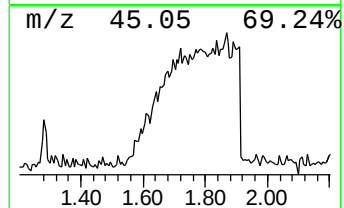
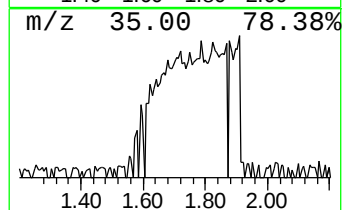
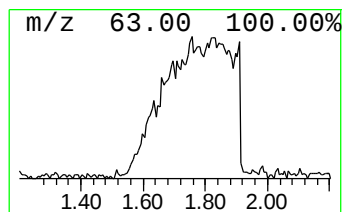
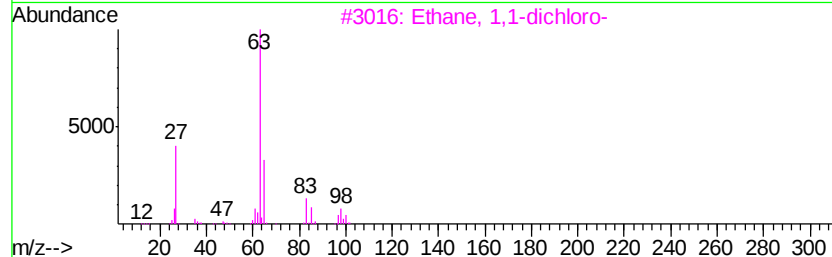
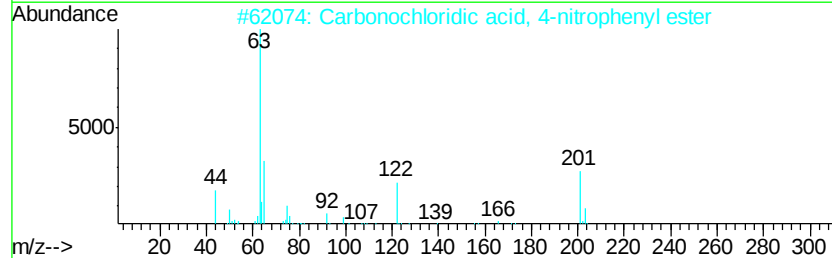
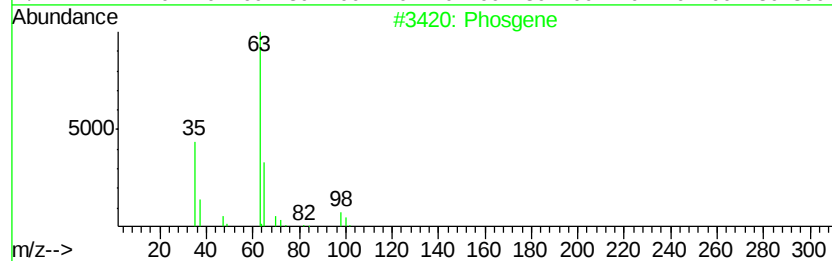
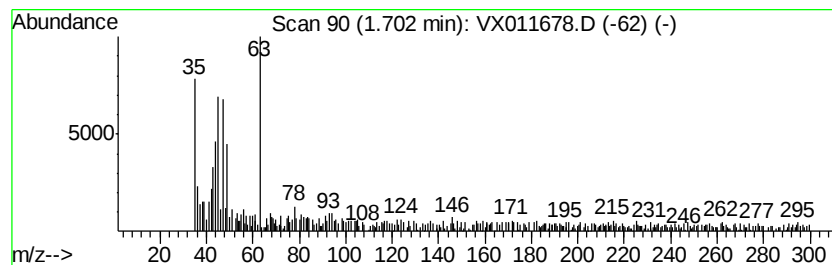
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X081519W.M
Quant Title : SW846 8260

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

Peak Number 1 unknown1.70 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.70	16.91 ug/l	326910	Pentafluorobenzene	5.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phosgene	98	CCl2O	000075-44-5	40
2		Carbonochloridic acid, 4-nitroph...	201	C7H4ClN04	007693-46-1	32
3		Ethane, 1,1-dichloro-	98	C2H4Cl2	000075-34-3	27
4		1-Bromo-1-chloroethane	142	C2H4BrCl	000593-96-4	27
5		Bis(2-(2-chloroethoxy)ethyl)ether	230	C8H16Cl2O3	000638-56-2	25



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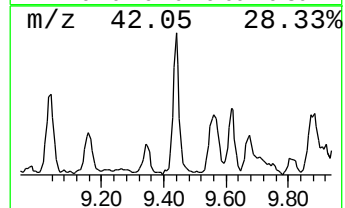
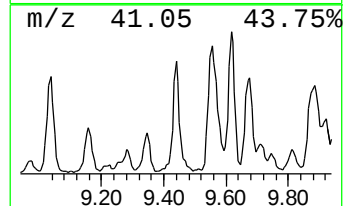
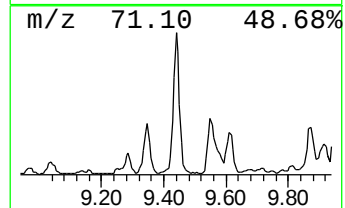
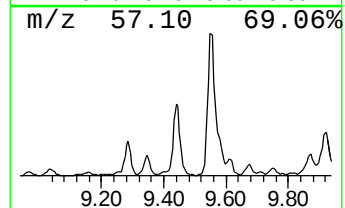
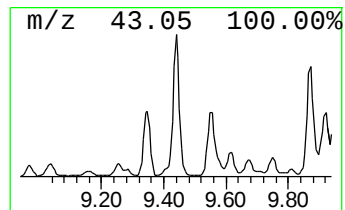
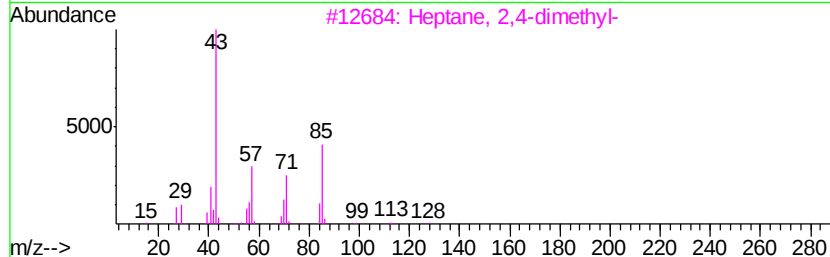
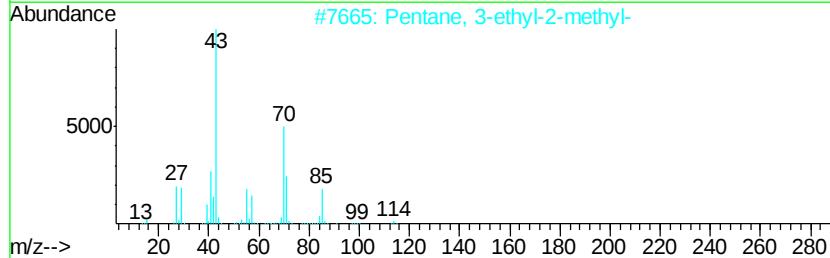
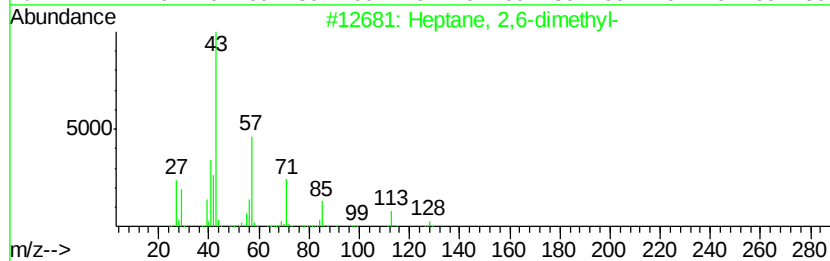
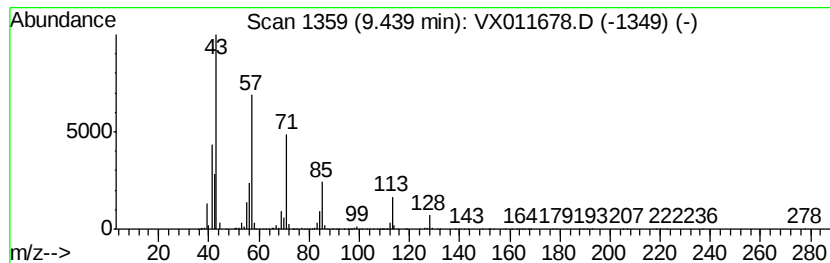
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TIC Integration Parameters: LSCINT.P

Peak Number 2 Heptane, 2,6-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.44	13.13 ug/l	342384	Chlorobenzene-d5	10.11

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 2,6-dimethyl-	128	C9H20	001072-05-5	90	
2	Pentane, 3-ethyl-2-methyl-	114	C8H18	000609-26-7	64	
3	Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	59	
4	Dodecane, 5-methyl-	184	C13H28	017453-93-9	53	
5	Nonane	128	C9H20	000111-84-2	52	



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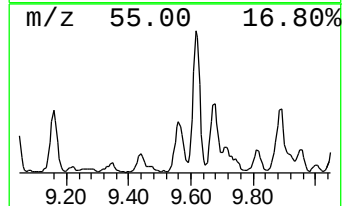
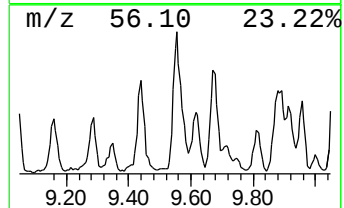
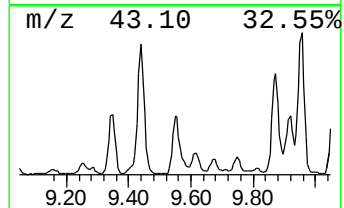
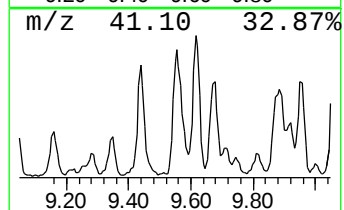
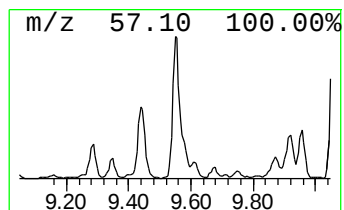
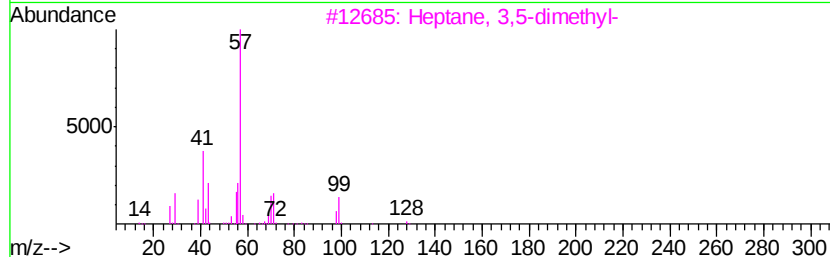
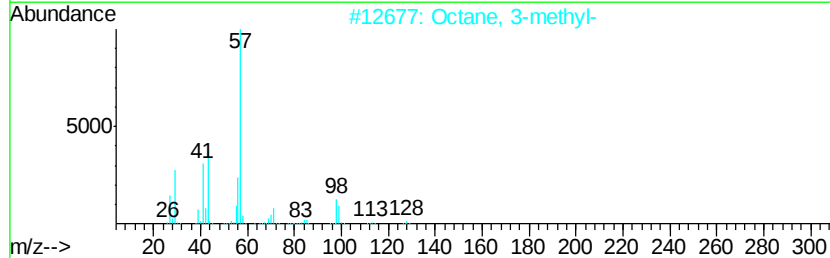
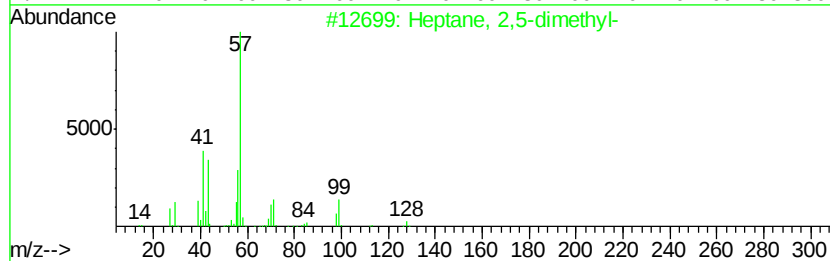
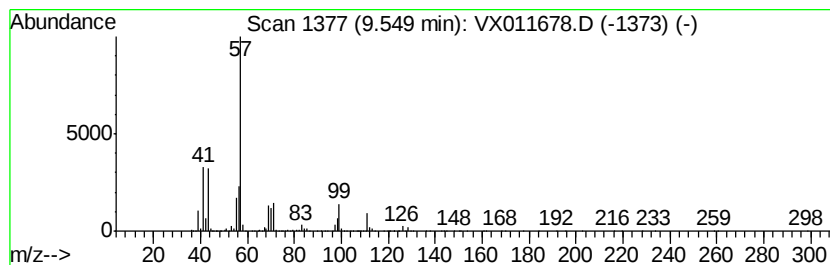
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X081519W.M
Quant Title : SW846 8260

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

Peak Number 3 Heptane, 2,5-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.55	18.40 ug/l	479996	Chlorobenzene-d5	10.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptane, 2,5-dimethyl-	128	C9H20	002216-30-0	76
2		Octane, 3-methyl-	128	C9H20	002216-33-3	52
3		Heptane, 3,5-dimethyl-	128	C9H20	000926-82-9	52
4		Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	50
5		Hexane, 3-ethyl-4-methyl-	128	C9H20	003074-77-9	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX082019\
Data File : VX011678.D
Acq On : 20 Aug 2019 12:41
Operator : JC/SP
Sample : K4391-04
Misc : 5.06g/10mL/100uL/10mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BOTTOM-1

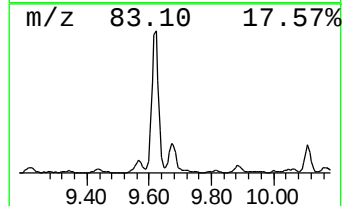
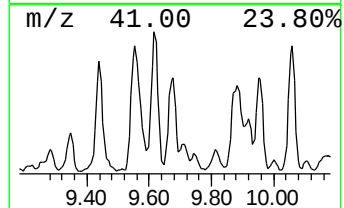
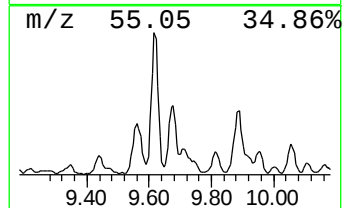
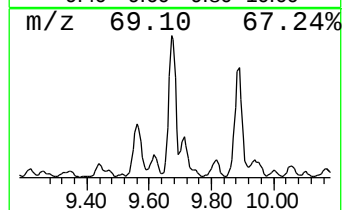
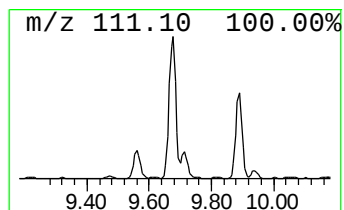
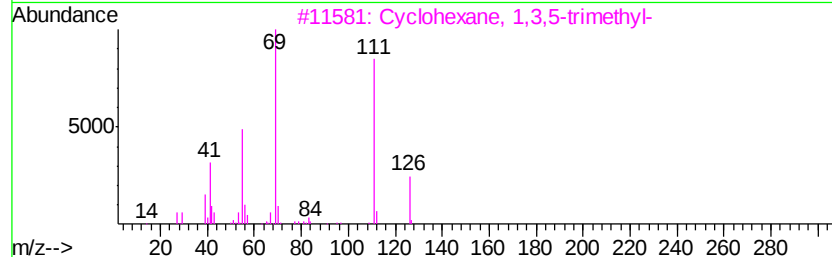
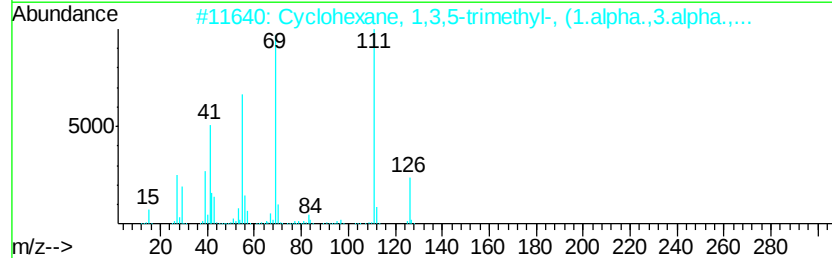
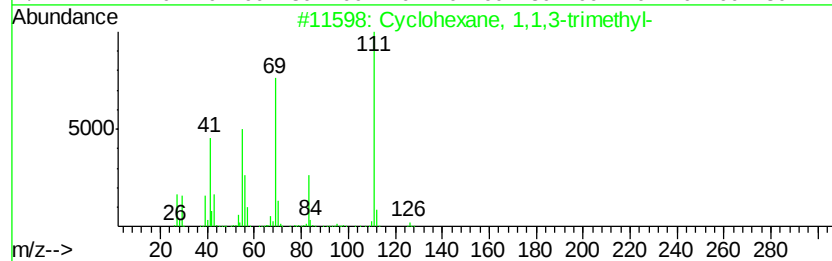
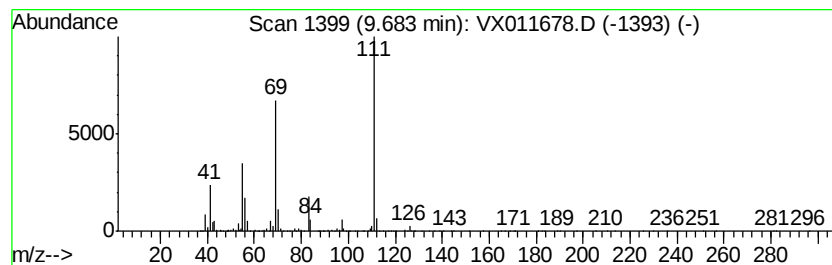
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X081519W.M
Quant Title : SW846 8260

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

Peak Number 5 Cyclohexane, 1,1,3-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.68	14.14 ug/l	368886	Chlorobenzene-d5	10.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	92
2		Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-27-3	78
3		Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	72
4		1,1,4-Trimethylcyclohexane	126	C9H18	007094-27-1	72
5		Cyclooctane, (1-methylpropyl)-	168	C12H24	016538-89-9	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX082019\
Data File : VX011678.D
Acq On : 20 Aug 2019 12:41
Operator : JC/SP
Sample : K4391-04
Misc : 5.06g/10mL/100uL/10mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BOTTOM-1

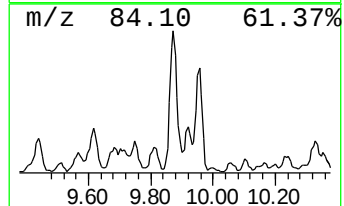
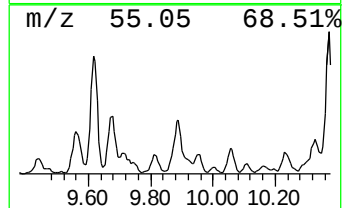
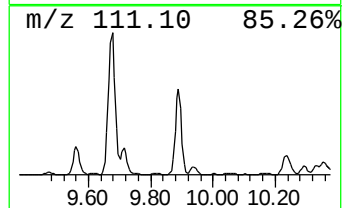
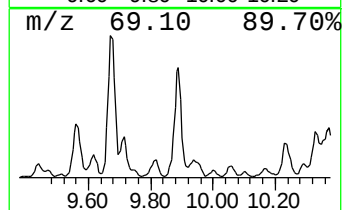
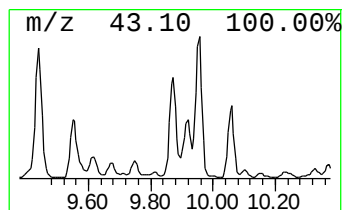
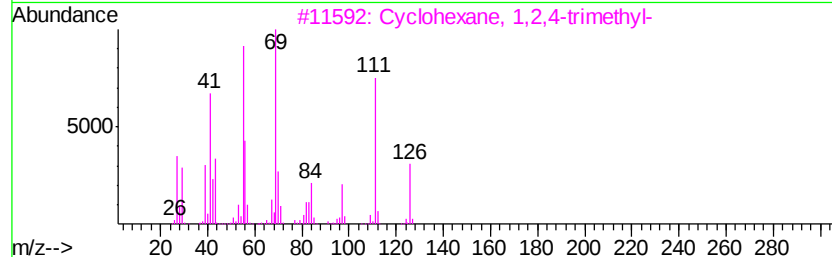
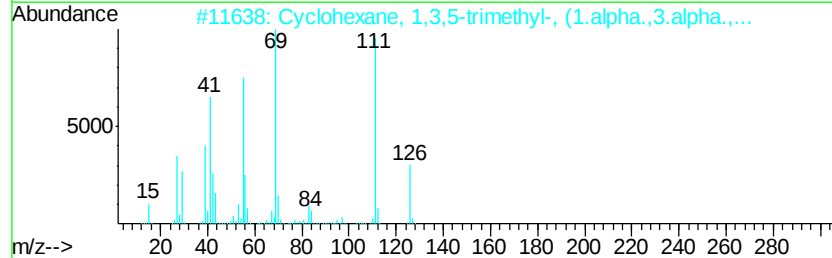
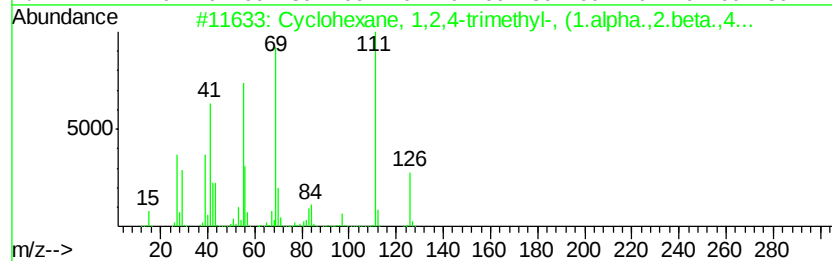
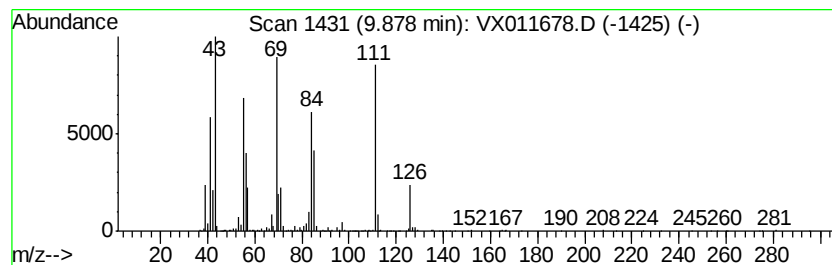
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Quant Title : SW846 8260

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

Peak Number 6 Cyclohexane, 1,2,4-trimethyl-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.88	17.90 ug/l	466778	Chlorobenzene-d5	10.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1,2,4-trimethyl-, (...	126	C9H18	007667-60-9	93
2		Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-26-2	87
3		Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	87
4		Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	83
5		Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-27-3	74



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX082019\
Data File : VX011678.D
Acq On : 20 Aug 2019 12:41
Operator : JC/SP
Sample : K4391-04
Misc : 5.06g/10mL/100uL/10mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BOTTOM-1

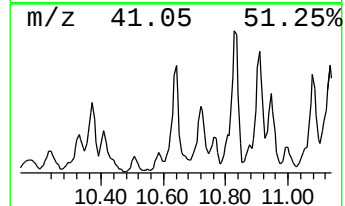
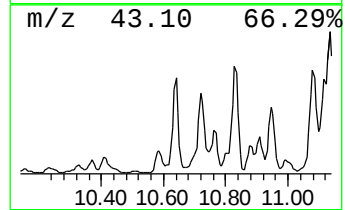
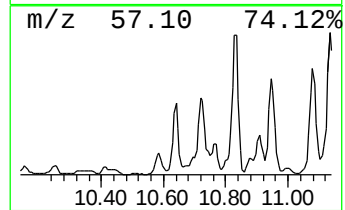
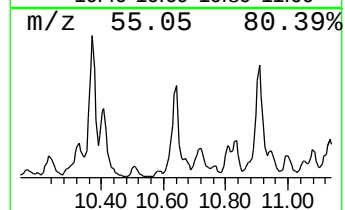
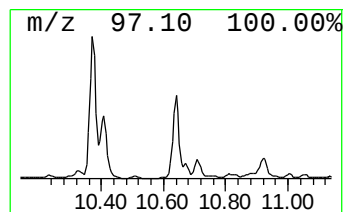
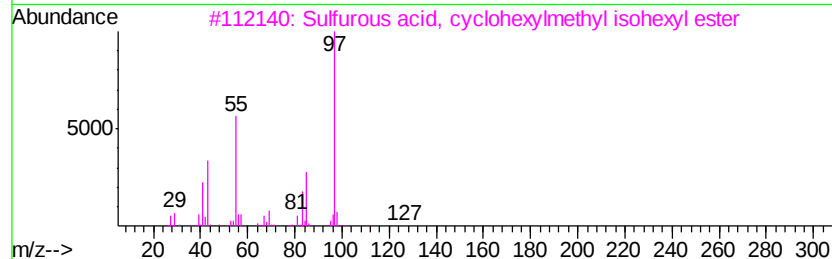
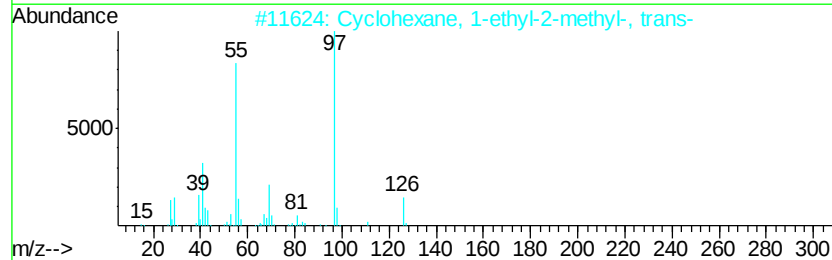
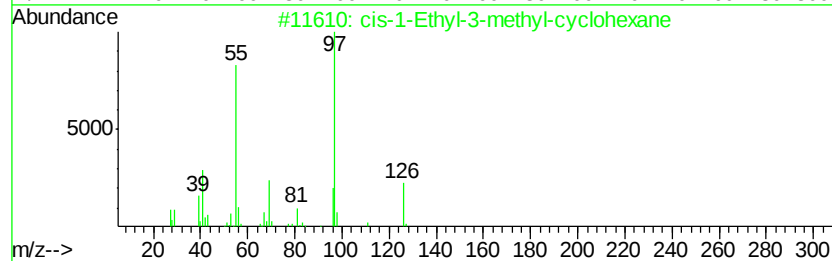
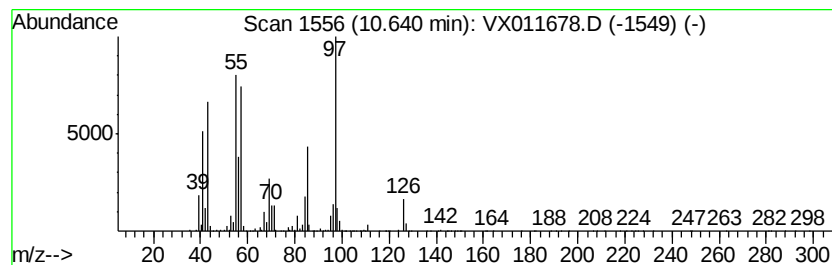
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X081519W.M
Quant Title : SW846 8260

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

Peak Number 7 cis-1-Ethyl-3-methyl-cyclohexane... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.64	19.93 ug/l	519681	Chlorobenzene-d5	10.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	cis-1-Ethyl-3-methyl-cyclohexane	126	C9H18	019489-10-2	55
2		Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	55
3		Sulfurous acid, cyclohexylmethyl...	262	C13H26O3S	1000309-21-5	53
4		Cyclohexane, 1-ethyl-2-methyl-	126	C9H18	003728-54-9	49
5		1-Ethyl-3-methylcyclohexane (c,t)	126	C9H18	003728-55-0	46



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX082019\
Data File : VX011678.D
Acq On : 20 Aug 2019 12:41
Operator : JC/SP
Sample : K4391-04
Misc : 5.06g/10mL/100uL/10mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
BOTTOM-1

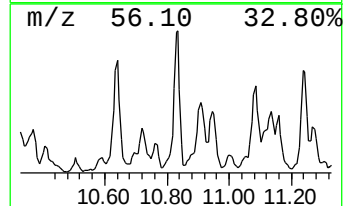
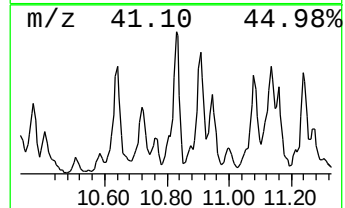
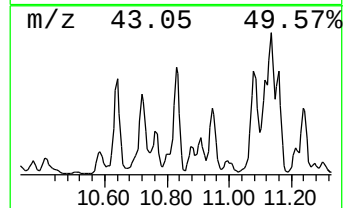
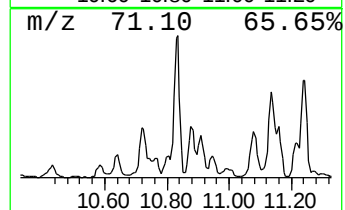
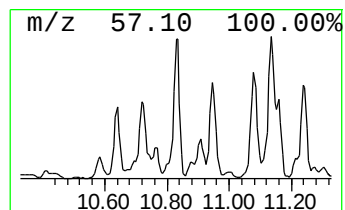
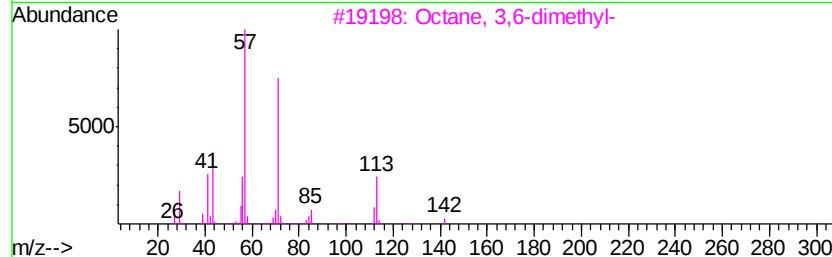
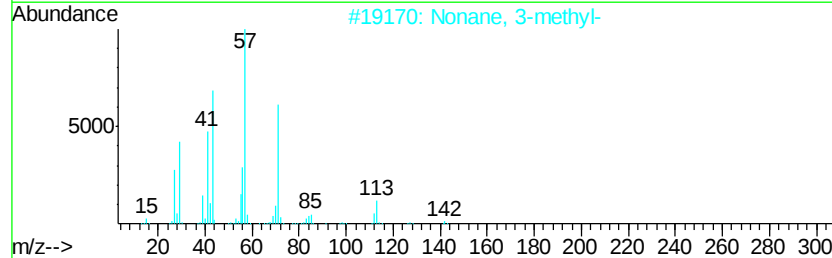
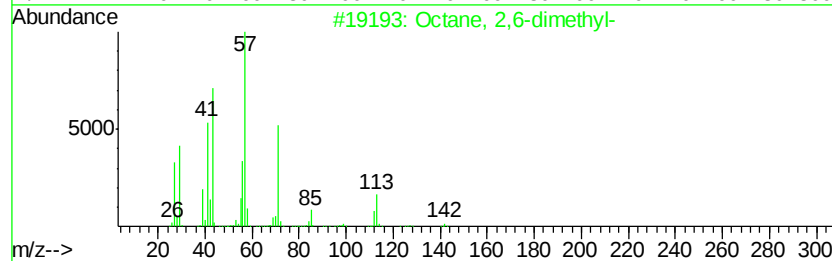
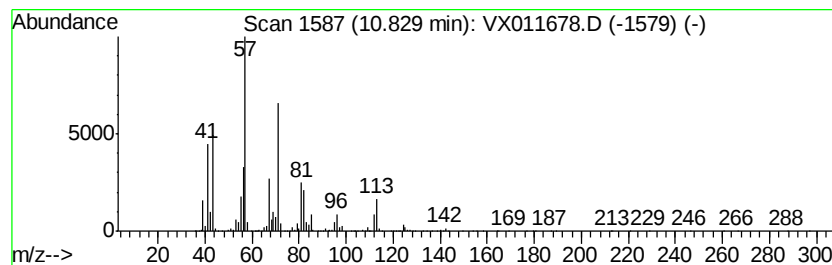
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X081519W.M
Quant Title : SW846 8260

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

Peak Number 8 Octane, 2,6-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.83	25.11 ug/l	654986	Chlorobenzene-d5	10.11

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	64
2		Nonane, 3-methyl-	142	C10H22	005911-04-6	60
3		Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	55
4		Sulfurous acid, di(2-ethylhexyl)...	306	C16H34O3S	1000309-19-1	43
5		Undecane, 6,6-dimethyl-	184	C13H28	017312-76-4	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX082019\
Data File : VX011678.D
Acq On : 20 Aug 2019 12:41
Operator : JC/SP
Sample : K4391-04
Misc : 5.06g/10mL/100uL/10mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BOTTOM-1

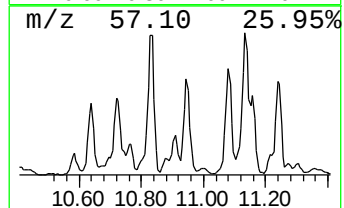
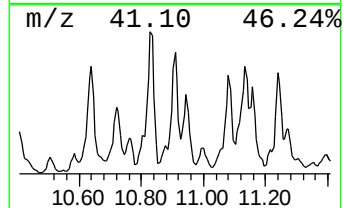
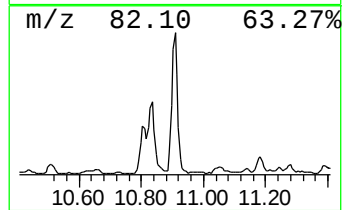
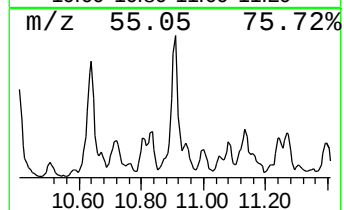
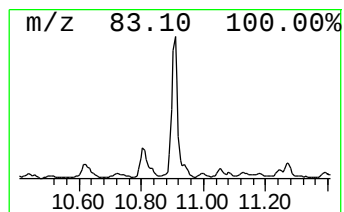
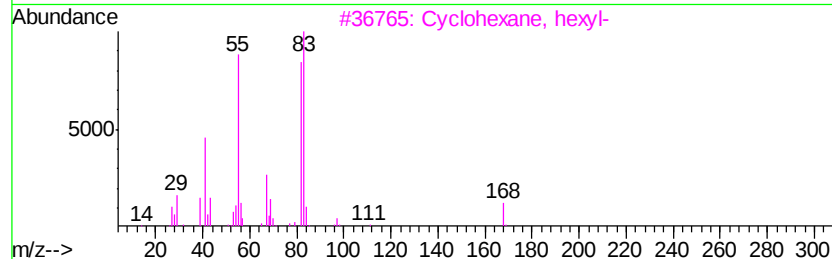
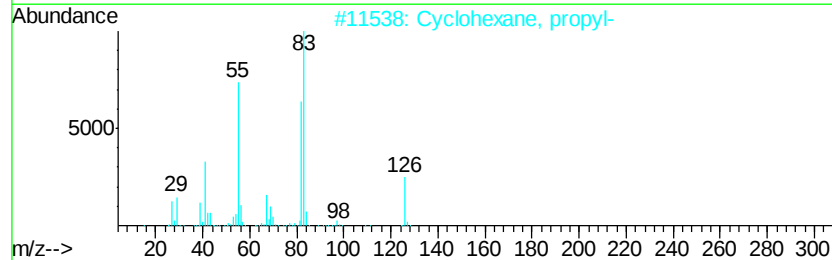
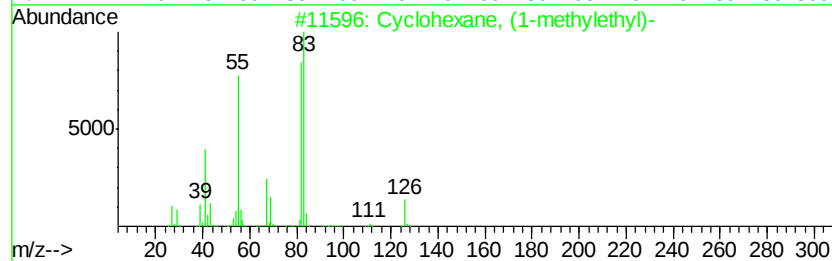
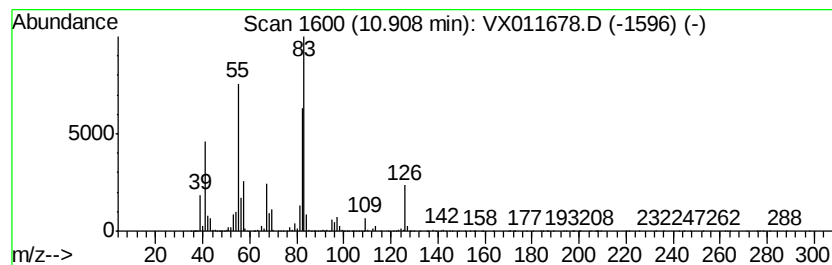
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X081519W.M
Quant Title : SW846 8260

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

Peak Number 9 Cyclohexane, (1-methylethyl)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.91	13.14 ug/l	342583	Chlorobenzene-d5	10.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	90
2		Cyclohexane, propyl-	126	C9H18	001678-92-8	87
3		Cyclohexane, hexyl-	168	C12H24	004292-75-5	72
4		Cyclohexane, octyl-	196	C14H28	001795-15-9	72
5		Ethanone, 1-(1-methylcyclopentyl)-	126	C8H14O	013388-93-7	64



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX082019\
Data File : VX011678.D
Acq On : 20 Aug 2019 12:41
Operator : JC/SP
Sample : K4391-04
Misc : 5.06g/10mL/100uL/10mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BOTTOM-1

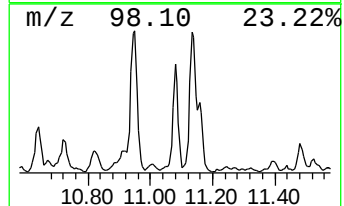
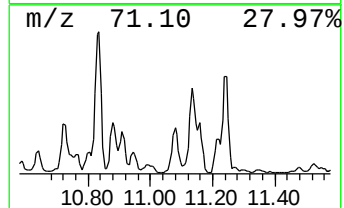
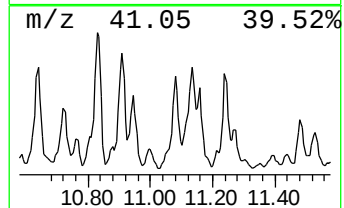
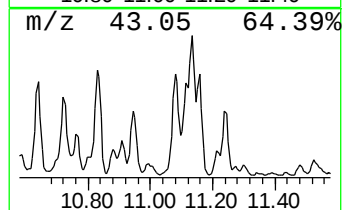
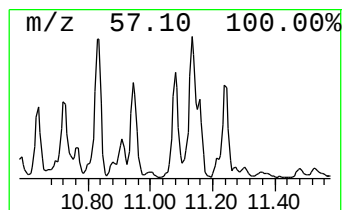
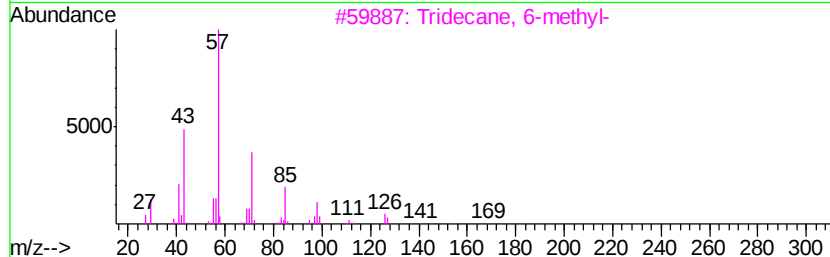
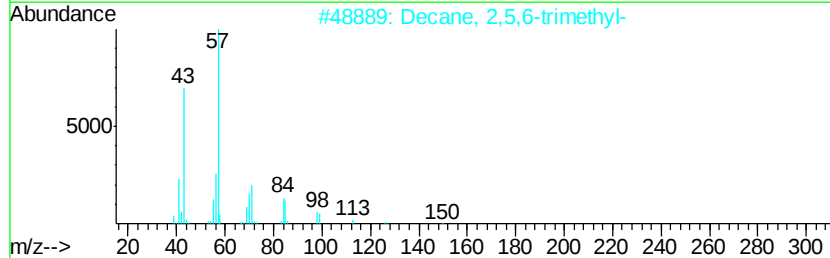
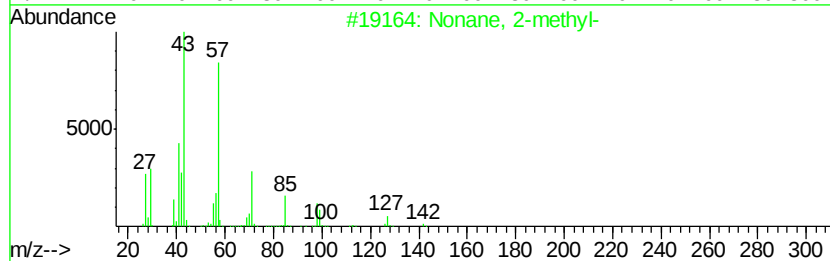
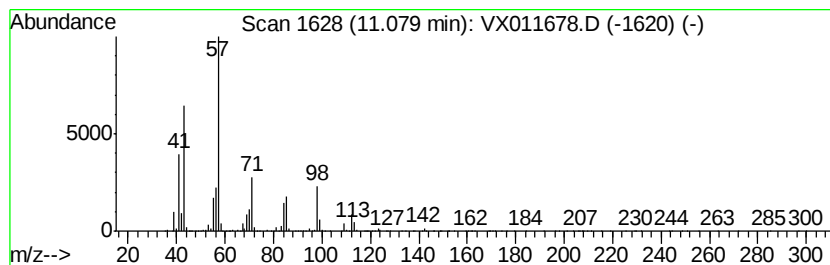
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Quant Title : SW846 8260

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

Peak Number 10 Nonane, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.08	17.47 ug/l	455641	Chlorobenzene-d5	10.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonane, 2-methyl-	142	C10H22	000871-83-0	52
2		Decane, 2,5,6-trimethyl-	184	C13H28	062108-23-0	50
3		Tridecane, 6-methyl-	198	C14H30	013287-21-3	50
4		Oxalic acid, isobutyl nonyl ester	272	C15H28O4	1000309-37-4	47
5		Octane, 2,3-dimethyl-	142	C10H22	007146-60-3	46



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX082019\
Data File : VX011678.D
Acq On : 20 Aug 2019 12:41
Operator : JC/SP
Sample : K4391-04
Misc : 5.06g/10mL/100uL/10mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BOTTOM-1

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X081519W.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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Heptane, 2,6-dime...	9.44	13.1	ug/l	342384	3	10.11	1304060	50.0
Heptane, 2,5-dime...	9.55	18.4	ug/l	479996	3	10.11	1304060	50.0
Cyclohexane, 1,1,...	9.68	14.1	ug/l	368886	3	10.11	1304060	50.0
Cyclohexane, 1,2,...	9.88	17.9	ug/l	466778	3	10.11	1304060	50.0
cis-1-Ethyl-3-met...	10.64	19.9	ug/l	519681	3	10.11	1304060	50.0
Octane, 2,6-dimet...	10.83	25.1	ug/l	654986	3	10.11	1304060	50.0
Cyclohexane, (1-m...	10.91	13.1	ug/l	342583	3	10.11	1304060	50.0
Nonane, 2-methyl-	11.08	17.5	ug/l	455641	3	10.11	1304060	50.0