

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX112120\
 Data File : VX019523.D
 Acq On : 21 Nov 2020 03:34
 Operator : JC/MD
 Sample : L4779-13
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 COMP-1654-1658

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\82X112120W.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	1.154	8	11	21	rBV	68440	66185	4.29%	0.348%
2	1.300	29	35	44	rBV3	20578	56049	3.63%	0.294%
3	1.380	44	48	53	rBV	65647	77465	5.02%	0.407%
4	1.471	60	63	68	rBV	41330	45145	2.93%	0.237%
5	1.575	75	80	91	rBV	8870	34288	2.22%	0.180%
6	2.123	165	170	176	rBV	83386	115886	7.51%	0.609%
7	2.331	200	204	214	rVB	17005	32889	2.13%	0.173%
8	2.428	214	220	235	rVB	362438	603604	39.11%	3.171%
9	2.611	242	250	259	rBV	83002	160993	10.43%	0.846%
10	3.044	306	321	332	rBV	60284	158461	10.27%	0.832%
11	4.306	518	528	546	rBV5	16489	67073	4.35%	0.352%
12	4.641	573	583	603	rVB	93956	272837	17.68%	1.433%
13	5.123	651	662	671	rBV2	13399	47791	3.10%	0.251%
14	5.464	704	718	732	rBV	172655	492724	31.93%	2.588%
15	5.623	732	744	760	rVV	312658	884174	57.29%	4.644%
16	6.031	797	811	819	rBV	189931	503384	32.62%	2.644%
17	6.434	867	877	887	rBV4	11504	36880	2.39%	0.194%
18	6.708	909	922	931	rBV	29790	80788	5.23%	0.424%
19	6.830	931	942	953	rBV	479255	1154253	74.79%	6.063%
20	7.513	1035	1054	1066	rBV	127016	280046	18.15%	1.471%
21	7.635	1068	1074	1080	rVV3	20276	48926	3.17%	0.257%
22	7.696	1080	1084	1095	rVB	21128	43369	2.81%	0.228%
23	7.958	1120	1127	1141	rVB2	63747	138191	8.95%	0.726%
24	8.622	1229	1236	1241	rBV	58532	93879	6.08%	0.493%
25	8.695	1241	1248	1254	rVV	980490	1543303	100.00%	8.107%
26	8.763	1254	1259	1270	rVV2	68138	136598	8.85%	0.718%
27	8.854	1270	1274	1278	rVV	21862	36947	2.39%	0.194%
28	9.073	1307	1310	1319	rVB	24814	39571	2.56%	0.208%
29	9.287	1339	1345	1349	rBV2	23400	40646	2.63%	0.214%
30	9.348	1349	1355	1363	rVB	38905	67258	4.36%	0.353%
31	9.476	1370	1376	1386	rBV	142831	210759	13.66%	1.107%
32	9.567	1386	1391	1402	rVB2	40421	69397	4.50%	0.365%
33	9.677	1402	1409	1417	rBV2	13655	31986	2.07%	0.168%
34	10.098	1468	1478	1489	rBV	914923	1329693	86.16%	6.985%

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Integration Parameters: RTEINT.P

Integrator: RTE
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35	10.244	1496	1502	1506	rBV3	27668	60854	3.94%	0.320%
36	10.281	1506	1508	1511	rVV	32278	40524	2.63%	0.213%
37	10.342	1514	1518	1525	rVV	37645	66389	4.30%	0.349%
38	10.427	1525	1532	1535	rVV	70208	111154	7.20%	0.584%
39	10.543	1545	1551	1555	rBV2	25441	33069	2.14%	0.174%
40	10.634	1561	1566	1570	rBV3	17516	31097	2.01%	0.163%
41	10.683	1570	1574	1577	rVV	29713	38530	2.50%	0.202%
42	10.719	1577	1580	1588	rVB	56288	71764	4.65%	0.377%
43	10.805	1588	1594	1600	rBV	259443	329564	21.35%	1.731%
44	10.896	1600	1609	1619	rVV7	50191	152691	9.89%	0.802%
45	11.067	1633	1637	1640	rBV	23552	32027	2.08%	0.168%
46	11.122	1640	1646	1653	rVV	795219	998317	64.69%	5.244%
47	11.207	1656	1660	1667	rVB3	21152	37329	2.42%	0.196%
48	11.335	1678	1681	1687	rVB3	35132	49382	3.20%	0.259%
49	11.414	1687	1694	1700	rBV2	22803	55919	3.62%	0.294%
50	11.469	1700	1703	1705	rVV	29155	36045	2.34%	0.189%
51	11.512	1706	1710	1720	rVB	584956	747028	48.40%	3.924%
52	11.597	1720	1724	1728	rBV	72451	106720	6.92%	0.561%
53	11.640	1728	1731	1737	rVV3	71467	129798	8.41%	0.682%
54	11.701	1737	1741	1746	rVB3	45217	72199	4.68%	0.379%
55	11.792	1746	1756	1760	rBV3	91490	152011	9.85%	0.798%
56	11.878	1765	1770	1775	rVB	164448	201191	13.04%	1.057%
57	11.933	1775	1779	1782	rBV2	29794	48498	3.14%	0.255%
58	12.061	1795	1800	1806	rVB	999613	1222709	79.23%	6.423%
59	12.140	1806	1813	1817	rBV2	77916	173534	11.24%	0.912%
60	12.256	1827	1832	1835	rBV	212712	296309	19.20%	1.556%
61	12.457	1861	1865	1868	rBV	137040	188318	12.20%	0.989%
62	12.487	1868	1870	1877	rVV3	85698	152528	9.88%	0.801%
63	12.555	1877	1881	1885	rVV	165942	255366	16.55%	1.341%
64	12.615	1885	1891	1895	rVV2	188540	325338	21.08%	1.709%
65	12.725	1904	1909	1914	rBV5	40271	87257	5.65%	0.458%
66	12.804	1918	1922	1925	rVV	116196	147619	9.57%	0.775%
67	12.835	1925	1927	1928	rVV2	49880	50364	3.26%	0.265%
68	12.853	1928	1930	1933	rVV	59460	71628	4.64%	0.376%
69	12.890	1933	1936	1941	rVV2	33863	53199	3.45%	0.279%
70	12.957	1944	1947	1950	rVV	35480	53355	3.46%	0.280%
71	13.000	1950	1954	1959	rVB	59410	91476	5.93%	0.481%

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72	13.201	1983	1987	1991	rVB2	134286	173929	11.27%	0.914%
73	13.329	2000	2008	2014	rVB4	163053	371181	24.05%	1.950%
74	13.420	2019	2023	2025	rBV3	23852	33072	2.14%	0.174%
75	13.463	2025	2030	2040	rVB	120163	185988	12.05%	0.977%
76	13.640	2052	2059	2062	rBV4	47772	92829	6.01%	0.488%
77	13.768	2078	2080	2084	rVB3	30131	32667	2.12%	0.172%
78	13.816	2084	2088	2094	rVB	203478	267088	17.31%	1.403%
79	13.890	2094	2100	2104	rBV5	40500	76204	4.94%	0.400%
80	13.938	2104	2108	2110	rBV2	24378	39374	2.55%	0.207%
81	14.079	2127	2131	2134	rBV	61699	69399	4.50%	0.365%
82	14.231	2152	2156	2158	rBV3	29421	40773	2.64%	0.214%
83	14.268	2159	2162	2166	rVB	59553	75206	4.87%	0.395%
84	14.414	2182	2186	2189	rBV2	38634	50054	3.24%	0.263%
85	14.518	2200	2203	2212	rVB3	58884	107324	6.95%	0.564%
86	14.603	2214	2217	2222	rVV2	38894	54940	3.56%	0.289%
87	14.676	2222	2229	2233	rVV2	148469	245859	15.93%	1.291%
88	14.725	2233	2237	2241	rVV3	45750	76146	4.93%	0.400%
89	14.792	2245	2248	2250	rVV	56703	63881	4.14%	0.336%
90	14.822	2250	2253	2257	rVB	84906	112414	7.28%	0.590%
91	14.957	2271	2275	2281	rBV3	59259	90783	5.88%	0.477%
92	15.255	2321	2324	2327	rVB	65903	72547	4.70%	0.381%
93	15.292	2327	2330	2339	rVB3	76040	114059	7.39%	0.599%
94	15.463	2354	2358	2364	rVB2	140235	229744	14.89%	1.207%
95	15.530	2365	2369	2375	rVV2	67692	116060	7.52%	0.610%
96	15.664	2387	2391	2394	rBV	61977	84550	5.48%	0.444%
97	15.700	2394	2397	2400	rVV	28956	39656	2.57%	0.208%
98	15.798	2409	2413	2419	rVB	93418	151488	9.82%	0.796%
99	16.042	2449	2453	2461	rVB6	59822	102897	6.67%	0.540%
100	16.773	2569	2573	2575	rBV2	62368	94856	6.15%	0.498%

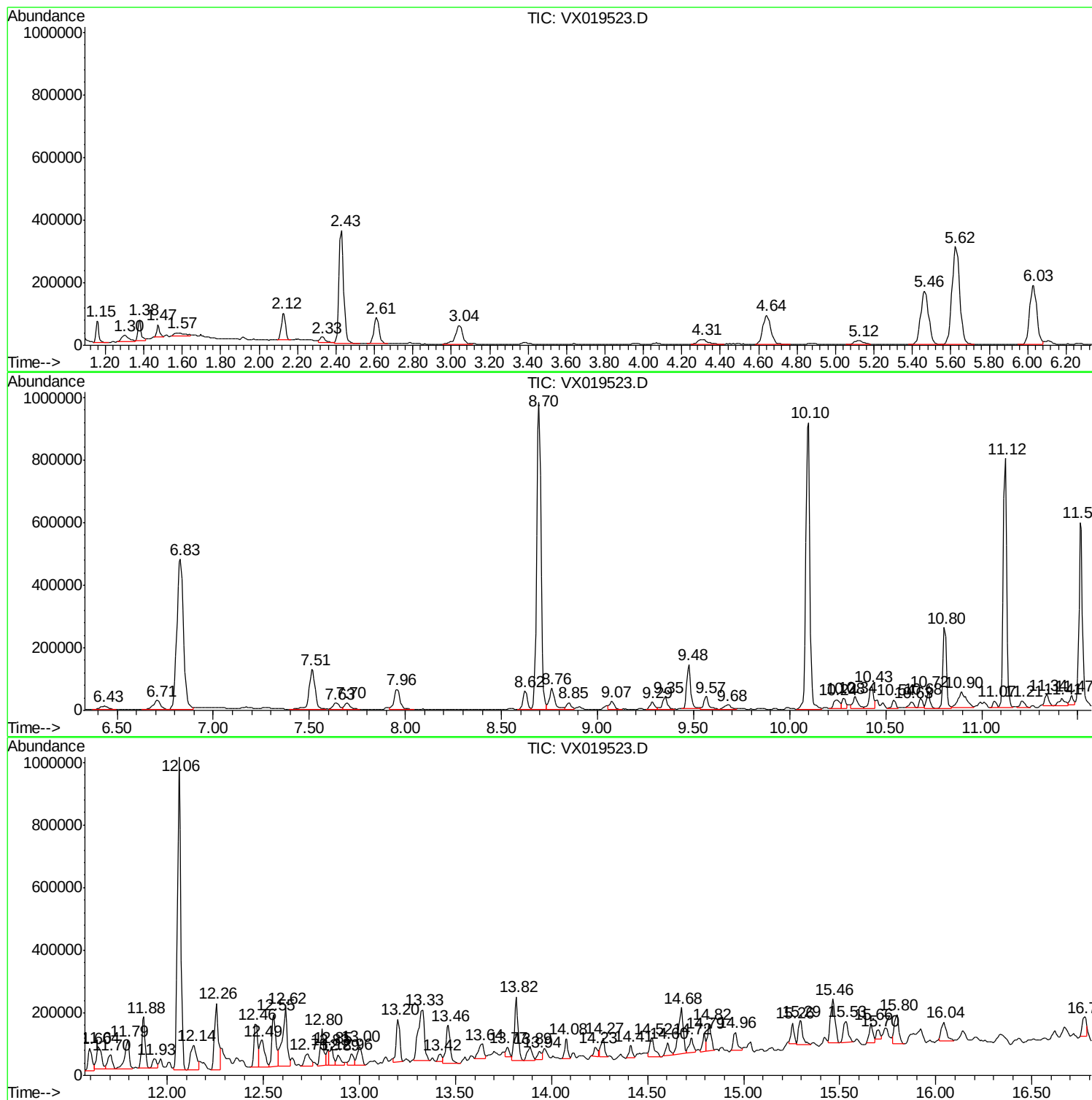
Sum of corrected areas: 19037506

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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112120W.M
Quant Title : SW846 8260

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



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ClientSampled :
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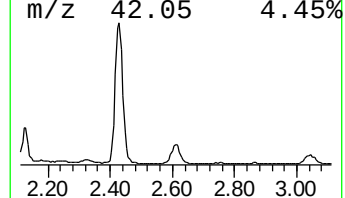
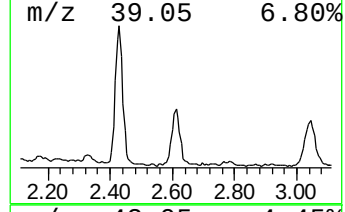
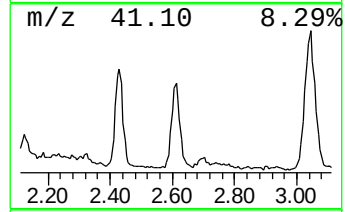
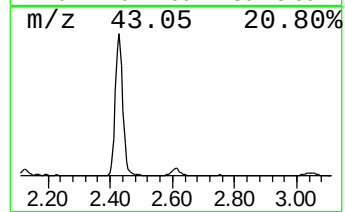
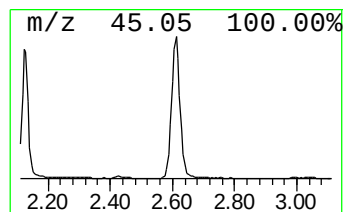
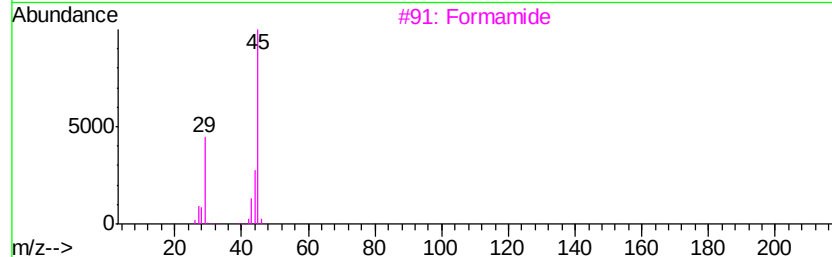
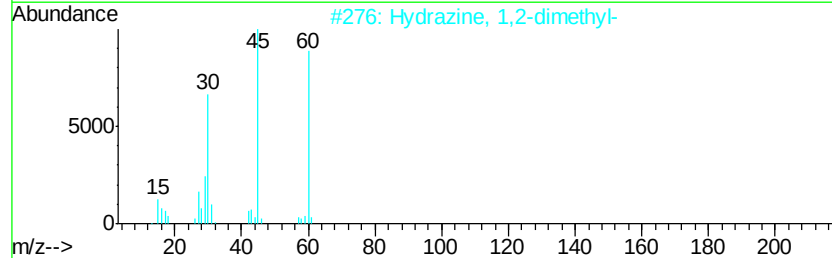
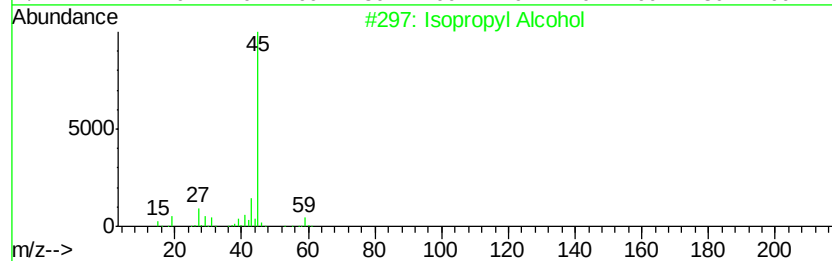
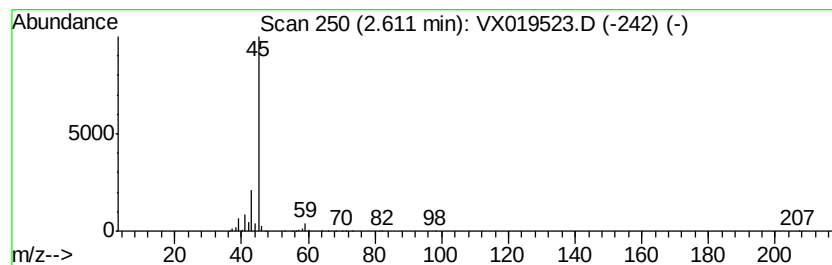
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112120W.M
Quant Title : SW846 8260

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

Peak Number 1 Isopropyl Alcohol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.61	9.10 ug/l	160993	Pentafluorobenzene	5.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Isopropyl Alcohol	60	C3H8O	000067-63-0	80
2		Hydrazine, 1,2-dimethyl-	60	C2H8N2	000540-73-8	9
3		Formamide	45	CH3NO	000075-12-7	5
4		Ethanol, 2-nitro-	91	C2H5NO3	000625-48-9	4
5		4-Penten-2-ol	86	C5H10O	000625-31-0	4



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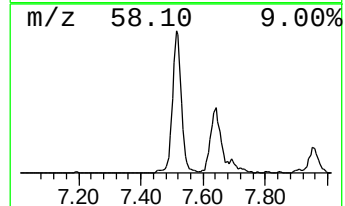
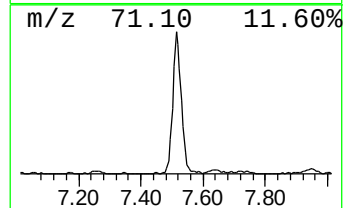
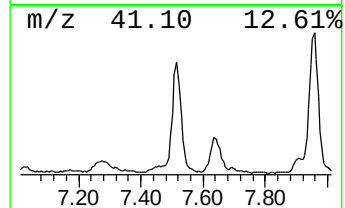
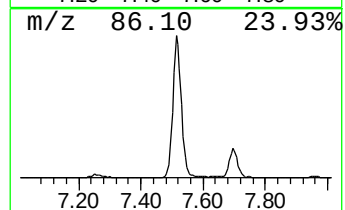
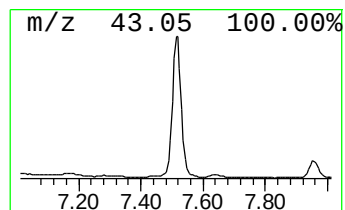
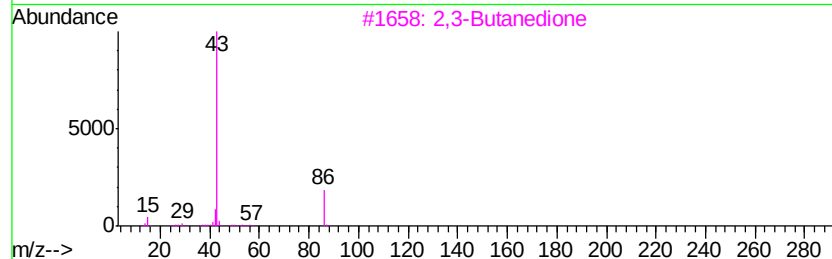
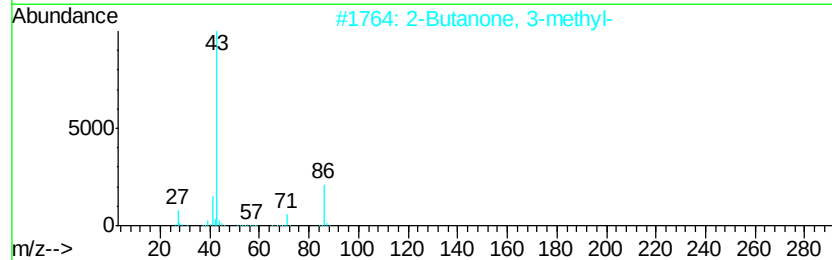
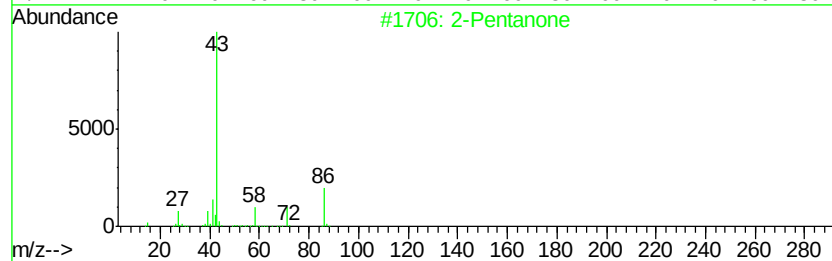
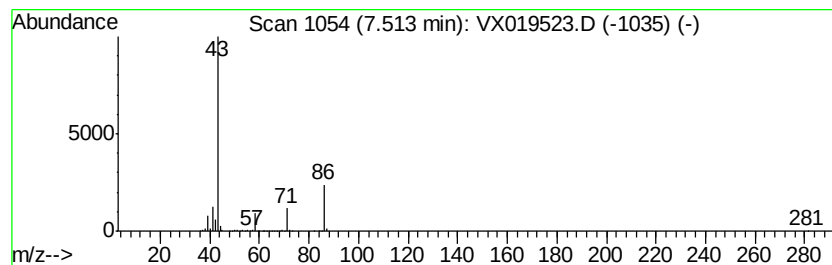
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112120W.M
Quant Title : SW846 8260

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

Peak Number 2 2-Pentanone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.51	12.13 ug/l	280046	1,4-Difluorobenzene	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone	86	C5H10O	000107-87-9	91
2		2-Butanone, 3-methyl-	86	C5H10O	000563-80-4	59
3		2,3-Butanedione	86	C4H6O2	000431-03-8	52
4		2,3-Hexanedione	114	C6H10O2	003848-24-6	12
5		Butane	58	C4H10	000106-97-8	9



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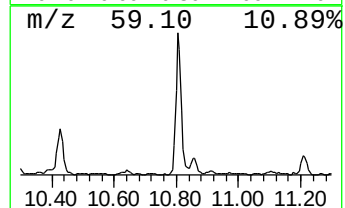
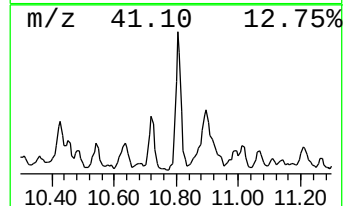
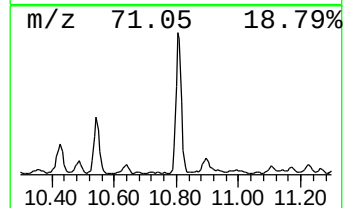
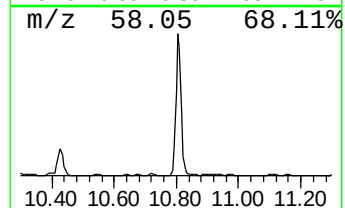
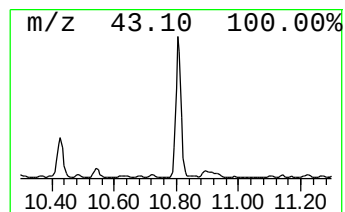
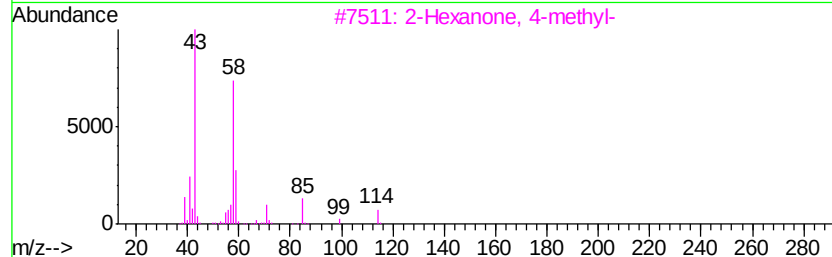
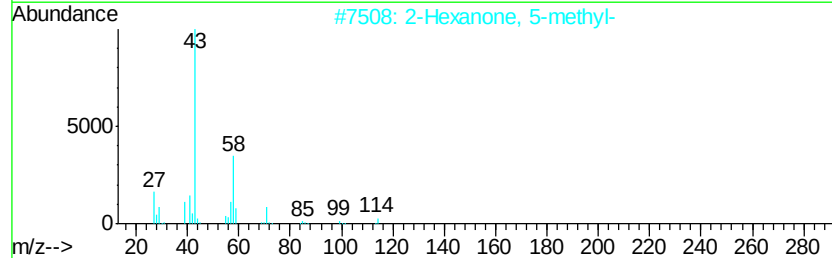
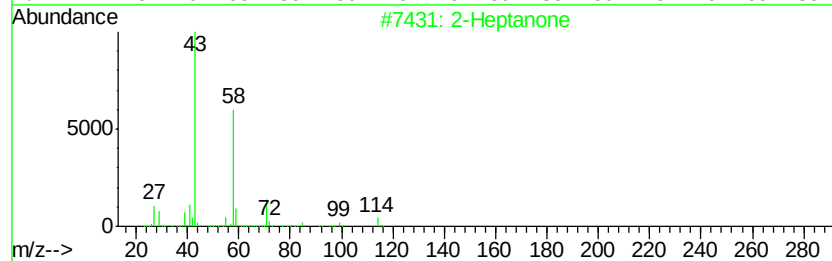
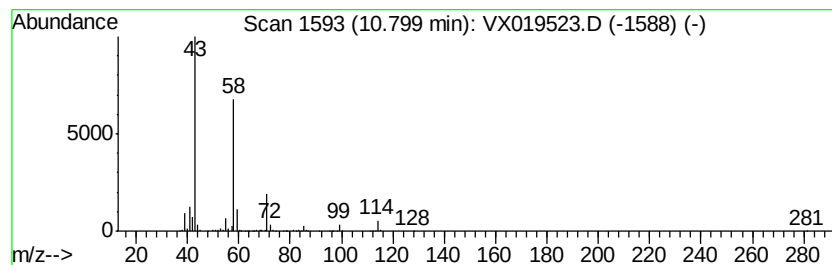
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TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Heptanone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.80	12.39 ug/l	329564	Chlorobenzene-d5	10.10	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Heptanone	114	C7H14O	000110-43-0	91
2	2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	83
3	2-Hexanone, 4-methyl-	114	C7H14O	000105-42-0	59
4	2-Pentanone, 4,4-dimethyl-	114	C7H14O	000590-50-1	47
5	2,3-Pentanedione, 4-methyl-	114	C6H10O2	007493-58-5	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX112120\
Data File : VX019523.D
Acq On : 21 Nov 2020 03:34
Operator : JC/MD
Sample : L4779-13
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 42 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
COMP-1654-1658

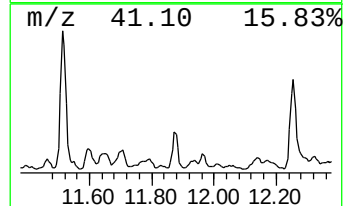
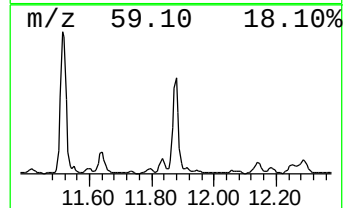
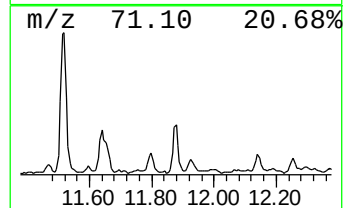
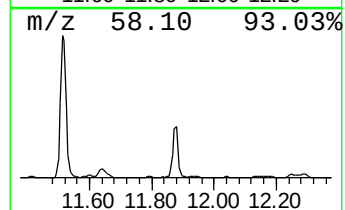
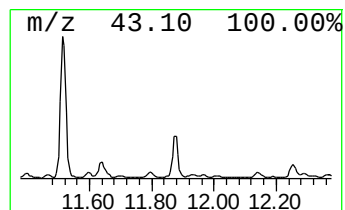
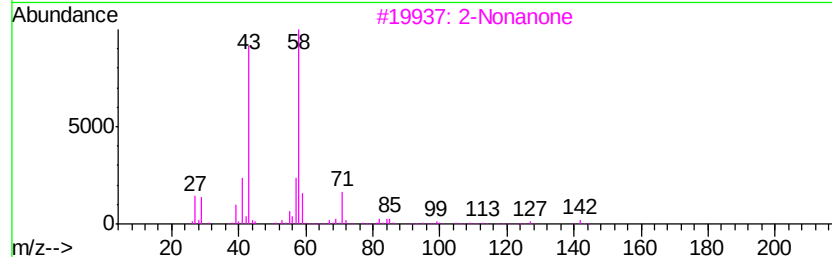
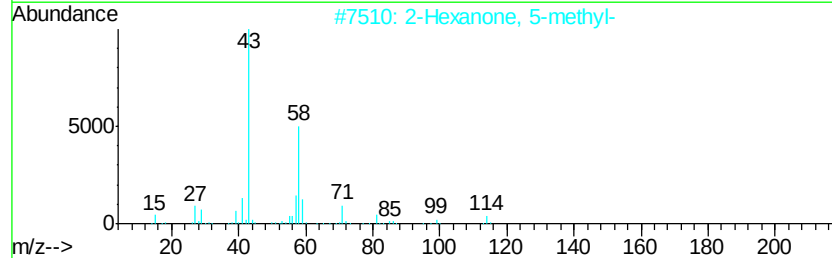
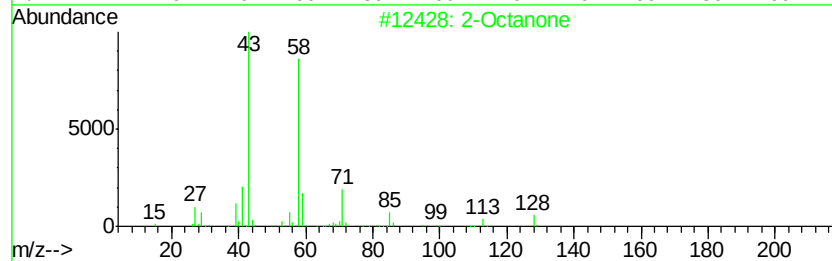
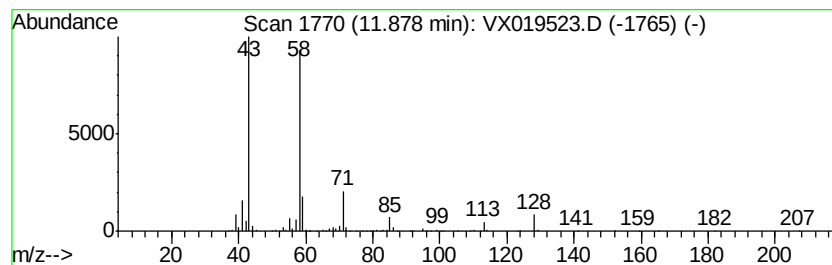
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Quant Title : SW846 8260

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

Peak Number 4 2-Octanone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.88	8.23 ug/l	201191	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Octanone	128	C8H16O	000111-13-7	94
2		2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	72
3		2-Nonanone	142	C9H18O	000821-55-6	72
4		2,4(3H,5H)-Furandione, 5-hexyl-5...	198	C11H18O3	054852-79-8	64
5		Butanimidamide	86	C4H10N2	000107-90-4	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX112120\
Data File : VX019523.D
Acq On : 21 Nov 2020 03:34
Operator : JC/MD
Sample : L4779-13
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 42 Sample Multiplier: 1

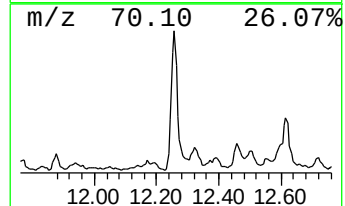
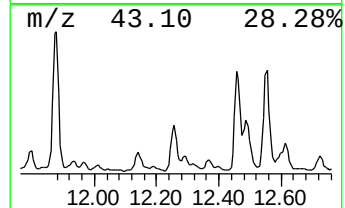
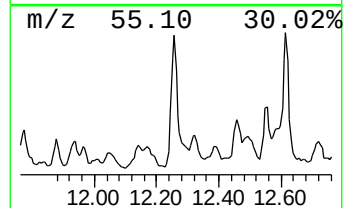
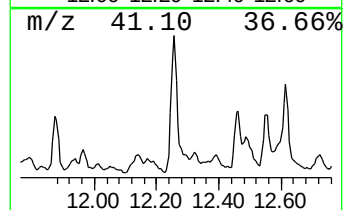
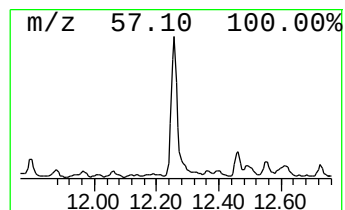
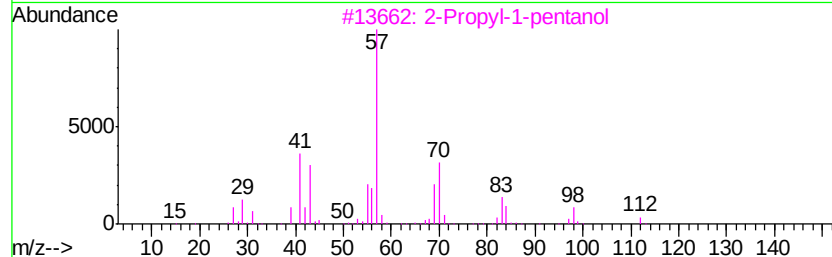
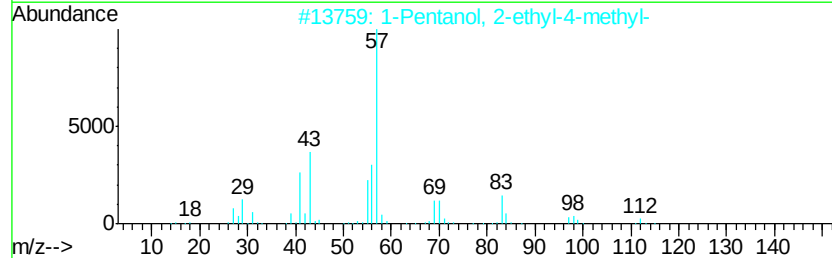
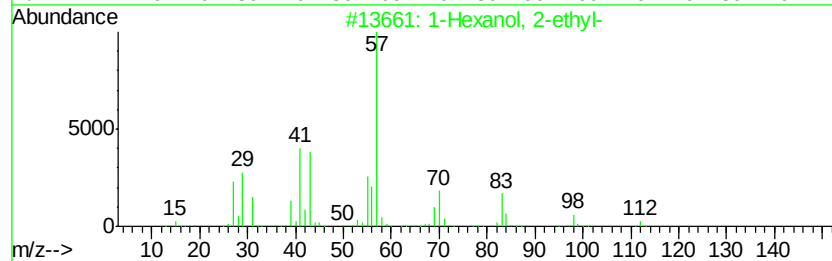
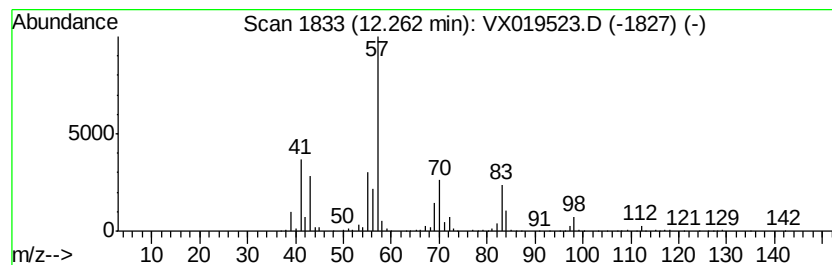
Instrument :
MSVOA_X
ClientSampleId :
COMP-1654-1658

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112120W.M
Quant Title : SW846 8260

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

Peak Number 5 1-Hexanol, 2-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.26	12.12 ug/l	296309	1,4-Dichlorobenzene-d4	12.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1-Hexanol, 2-ethyl-	130 C8H18O	000104-76-7 90
2		1-Pentanol, 2-ethyl-4-methyl-	130 C8H18O	000106-67-2 56
3		2-Propyl-1-pentanol	130 C8H18O	058175-57-8 50
4		Heptane, 1,1'-oxybis-	214 C14H30O	000629-64-1 50
5		2-Ethyl-1-hexanol, methyl ether	144 C9H20O	1000365-10-2 47



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX112120\
Data File : VX019523.D
Acq On : 21 Nov 2020 03:34
Operator : JC/MD
Sample : L4779-13
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 42 Sample Multiplier: 1

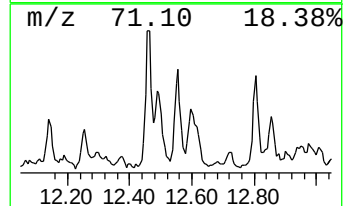
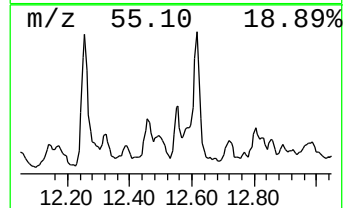
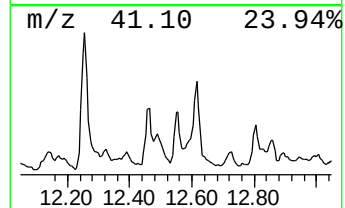
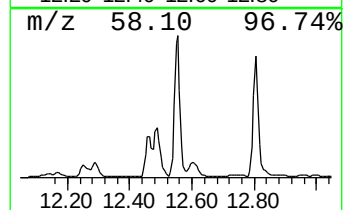
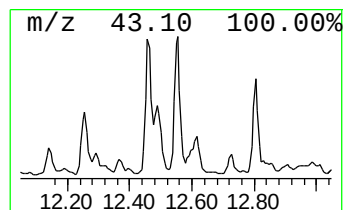
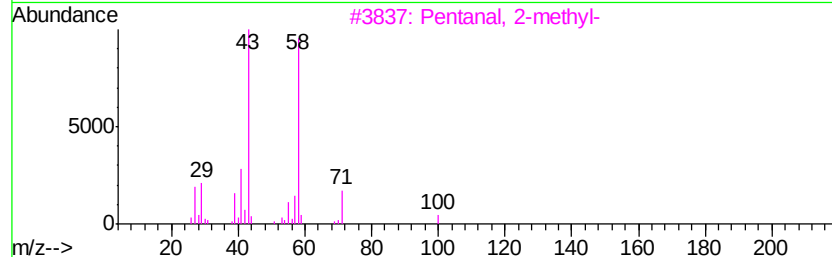
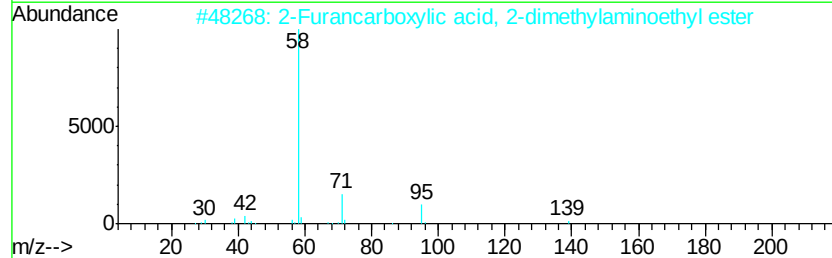
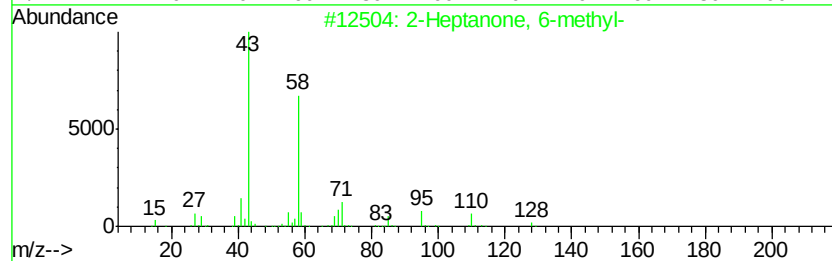
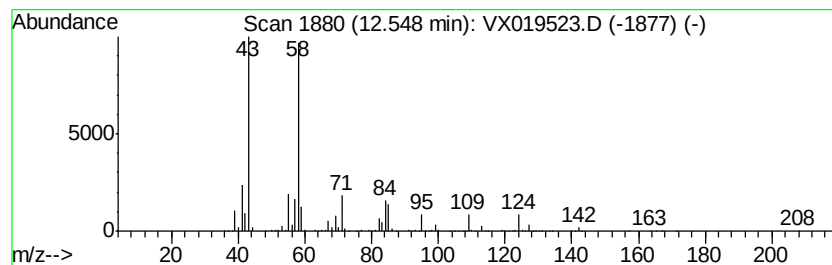
Instrument :
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ClientSampleId :
COMP-1654-1658

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112120W.M
Quant Title : SW846 8260

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

Peak Number 6 2-Heptanone, 6-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.55	10.44 ug/l	255366	1,4-Dichlorobenzene-d4	12.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2-Heptanone, 6-methyl-	128 C8H16O	000928-68-7	59
2	2-Furancarboxylic acid, 2-dimeth...	183 C9H13NO3	1000331-10-1	50
3	Pentanal, 2-methyl-	100 C6H12O	000123-15-9	47
4	2-Nonanone	142 C9H18O	000821-55-6	47
5	2-Octanone	128 C8H16O	000111-13-7	40



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX112120\
Data File : VX019523.D
Acq On : 21 Nov 2020 03:34
Operator : JC/MD
Sample : L4779-13
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 42 Sample Multiplier: 1

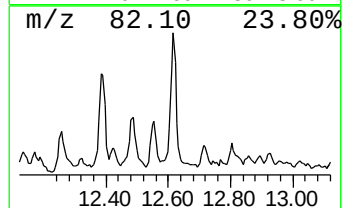
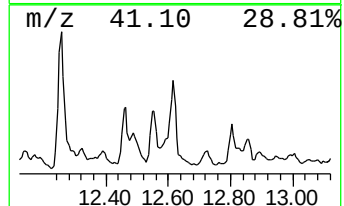
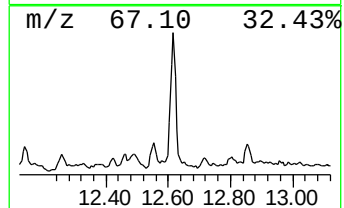
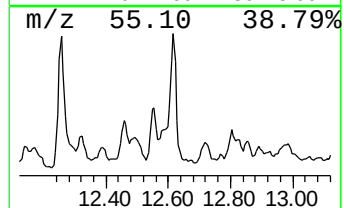
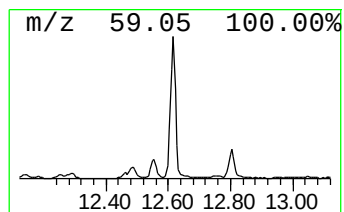
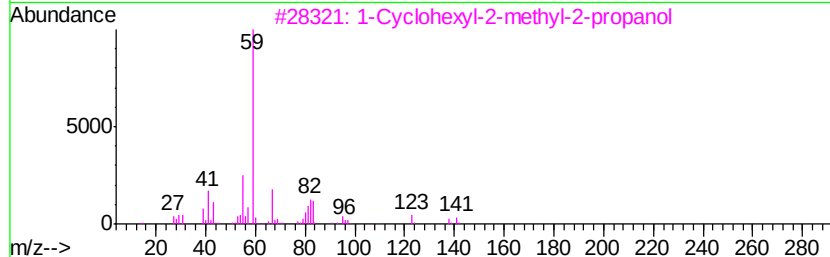
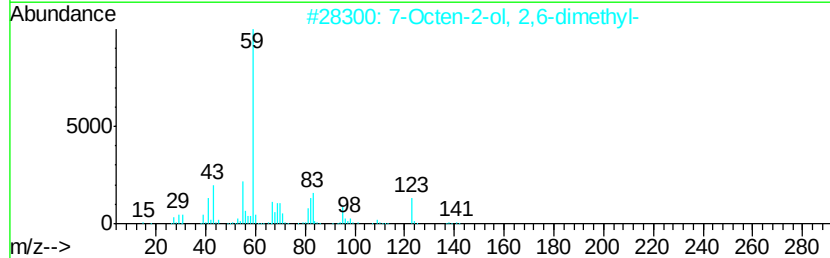
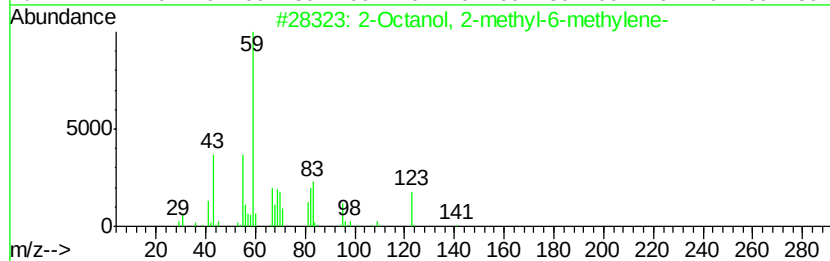
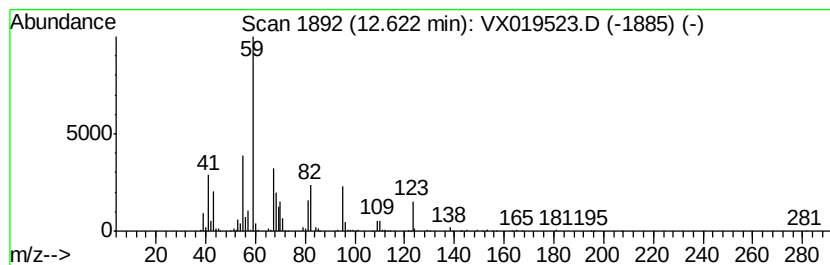
Instrument :
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ClientSampleId :
COMP-1654-1658

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112120W.M
Quant Title : SW846 8260

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

Peak Number 7 unknown12.62 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.62	13.30 ug/l	325338	1,4-Dichlorobenzene-d4	12.06	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Octanol, 2-methyl-6-methylene-	156	C10H20O	018479-59-9	38
2	7-Octen-2-ol, 2,6-dimethyl-	156	C10H20O	018479-58-8	38
3	1-Cyclohexyl-2-methyl-2-propanol	156	C10H20O	005531-30-6	36
4	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	000498-81-7	28
5	1,7-Octanediol, 3,7-dimethyl-	174	C10H22O2	000107-74-4	28



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX112120\
Data File : VX019523.D
Acq On : 21 Nov 2020 03:34
Operator : JC/MD
Sample : L4779-13
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 42 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
COMP-1654-1658

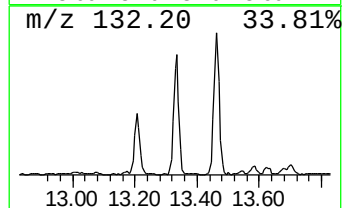
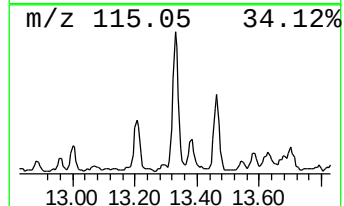
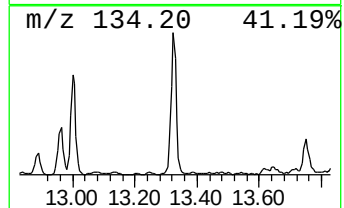
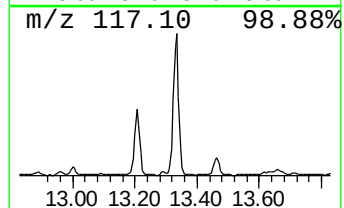
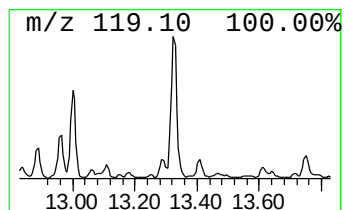
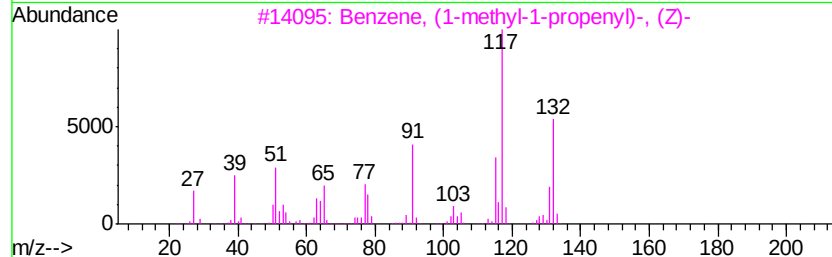
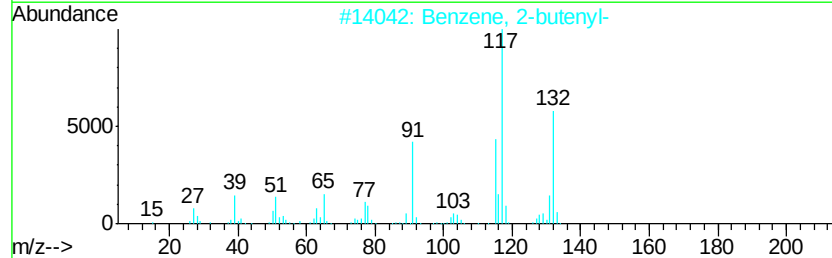
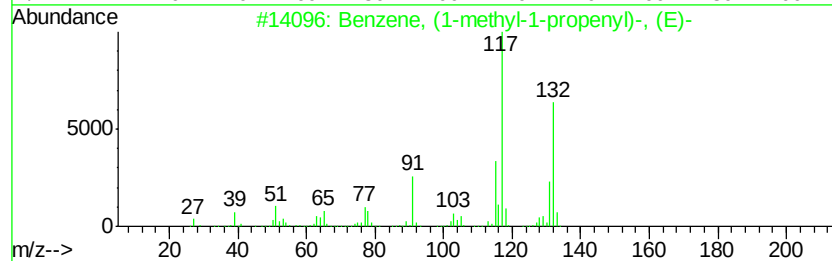
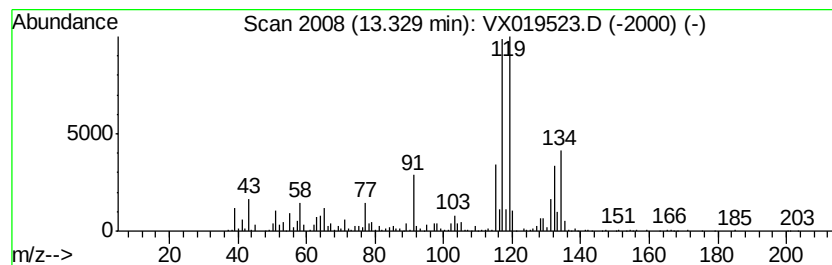
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, (1-methyl-1-propen... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.33	15.18 ug/l	371181	1,4-Dichlorobenzene-d4	12.06

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	64
2			Benzene, 2-butenyl-	132	C10H12	001560-06-1	64
3			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	50
4			Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	50
5			o-Cymene	134	C10H14	000527-84-4	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX112120\
Data File : VX019523.D
Acq On : 21 Nov 2020 03:34
Operator : JC/MD
Sample : L4779-13
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 42 Sample Multiplier: 1

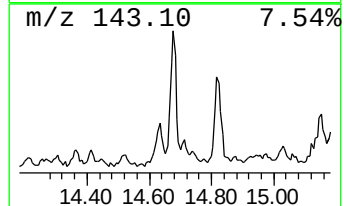
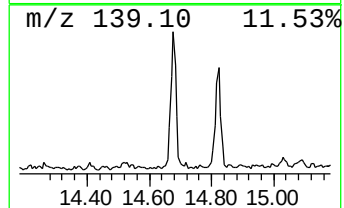
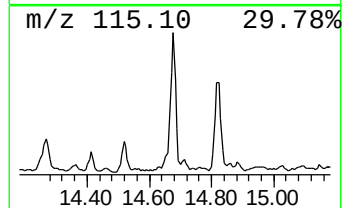
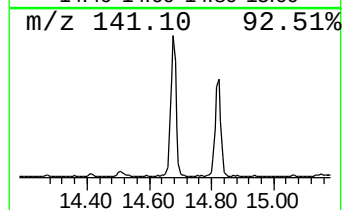
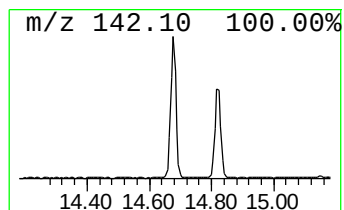
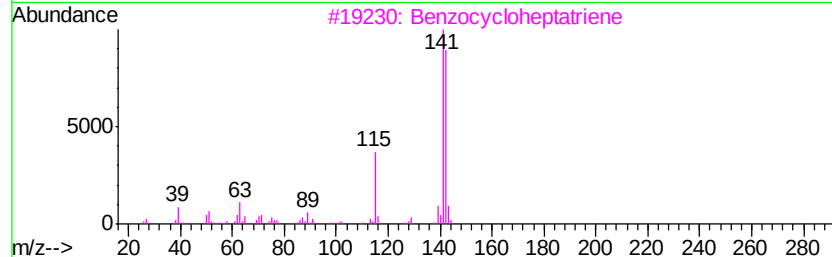
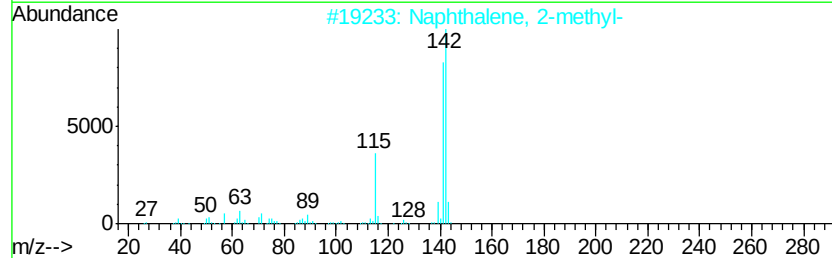
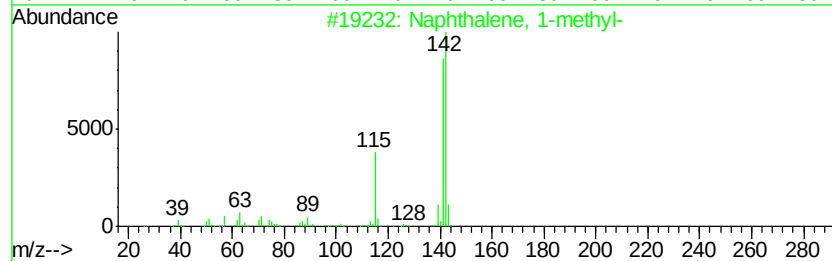
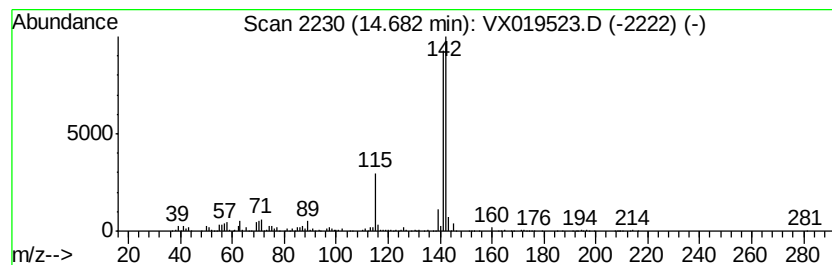
Instrument :
MSVOA_X
ClientSampleId :
COMP-1654-1658

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112120W.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.68	10.05 ug/l	245859	1,4-Dichlorobenzene-d4	12.06
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 94
3		Benzocycloheptatriene	142 C11H10	000264-09-5 91
4		1,4-Methanonaphthalene, 1,4-dihy...	142 C11H10	004453-90-1 90
5		Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142 C11H10	002443-46-1 87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX112120\
Data File : VX019523.D
Acq On : 21 Nov 2020 03:34
Operator : JC/MD
Sample : L4779-13
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ALS Vial : 42 Sample Multiplier: 1

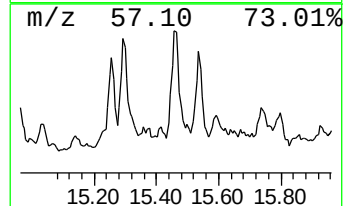
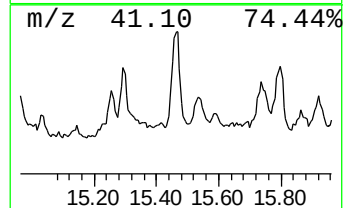
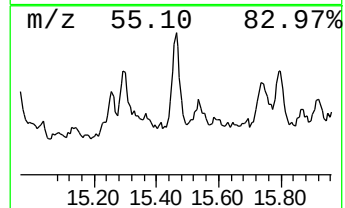
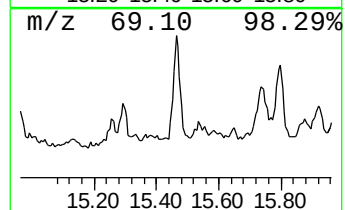
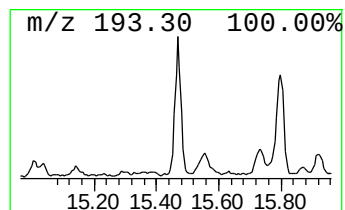
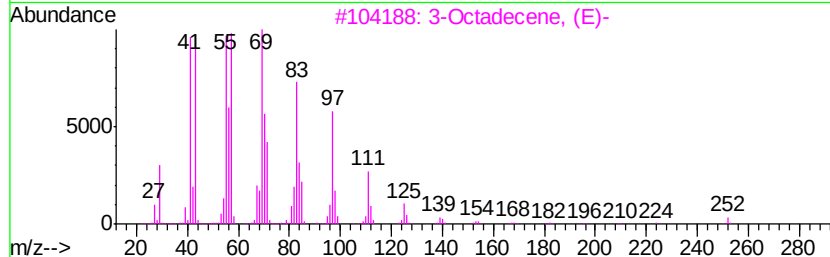
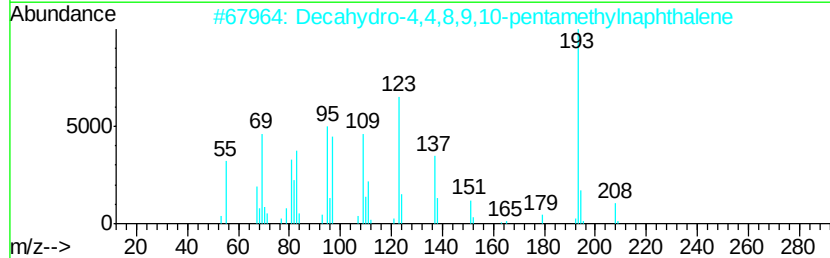
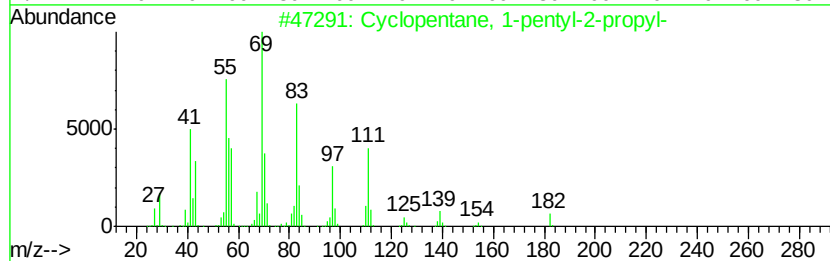
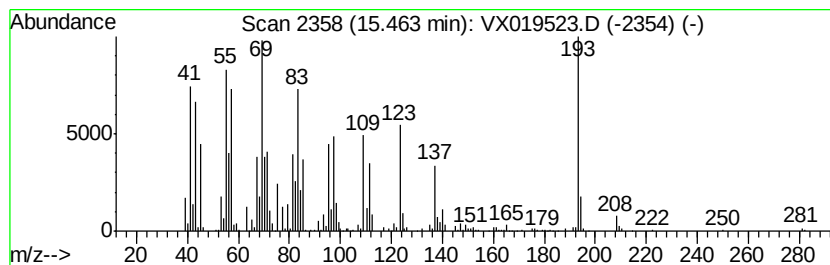
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COMP-1654-1658

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X112120W.M
Quant Title : SW846 8260

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

Peak Number 10 unknown15.46 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.46	9.39 ug/l	229744	1,4-Dichlorobenzene-d4	12.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclopentane, 1-pentyl-2-propyl-	182 C13H26	062199-51-3 43
2		Decahydro-4,4,8,9,10-pentamethyl...	208 C15H28	080655-44-3 38
3		3-Octadecene, (E)-	252 C18H36	007206-19-1 30
4		3-Eicosene, (E)-	280 C20H40	074685-33-9 30
5		3-Chloropropionic acid, heptadec...	346 C20H39ClO2	1000283-05-1 25



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX112120\
Data File : VX019523.D
Acq On : 21 Nov 2020 03:34
Operator : JC/MD
Sample : L4779-13
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 42 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
COMP-1654-1658

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112120W.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
Isopropyl Alcohol	2.61	9.1	ug/l	160993	1	5.62	884174	50.0
2-Pentanone	7.51	12.1	ug/l	280046	2	6.83	1154250	50.0
2-Heptanone	10.80	12.4	ug/l	329564	3	10.10	1329690	50.0
2-Octanone	11.88	8.2	ug/l	201191	4	12.06	1222710	50.0
1-Hexanol, 2-ethyl-	12.26	12.1	ug/l	296309	4	12.06	1222710	50.0
2-Heptanone, 6-me...	12.55	10.4	ug/l	255366	4	12.06	1222710	50.0
unknown12.62	12.62	13.3	ug/l	325338	4	12.06	1222710	50.0
Benzene, (1-methy...	13.33	15.2	ug/l	371181	4	12.06	1222710	50.0
Naphthalene, 1-me...	14.68	10.1	ug/l	245859	4	12.06	1222710	50.0
unknown15.46	15.46	9.4	ug/l	229744	4	12.06	1222710	50.0