

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV031620\
 Data File : VV014953.D
 Acq On : 16 Mar 2020 18:07
 Operator : SY/MD
 Sample : L1890-18
 Misc : 5.0mL/MSVOA V/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 C0DH5

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0 % 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_V\METHOD\SOMVLM031620WMA.M
 Title : VOC Analysis

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.222	36	41	47	rBV2	10868	10433	0.00%	0.001%
2	1.315	62	70	85	rVB	199783	215547	0.06%	0.023%
3	1.528	134	136	141	rBV5	1399	1543	0.00%	0.000%
4	1.579	144	152	167	rVV	132728	162138	0.04%	0.017%
5	2.119	310	320	339	rVB	265514	390972	0.10%	0.042%
6	2.309	370	379	387	rBV	6935	11690	0.00%	0.001%
7	2.524	437	446	457	rBV	19980	37335	0.01%	0.004%
8	2.592	459	467	485	rVB	38369	67645	0.02%	0.007%
9	2.698	497	500	502	rBV2	483	305	0.00%	0.000%
10	2.778	516	525	537	rVB10	5512	10660	0.00%	0.001%
11	2.878	554	556	561	rVB2	581	545	0.00%	0.000%
12	2.901	561	563	566	rBV2	394	296	0.00%	0.000%
13	3.010	591	597	601	rVB3	452	463	0.00%	0.000%
14	3.064	605	614	621	rBV5	2196	3610	0.00%	0.000%
15	3.129	630	634	638	rBV2	441	326	0.00%	0.000%
16	3.161	638	644	646	rVV2	328	312	0.00%	0.000%
17	3.187	649	652	653	rVV	559	299	0.00%	0.000%
18	3.196	653	655	657	rVV3	692	434	0.00%	0.000%
19	3.209	657	659	668	rVV2	529	567	0.00%	0.000%
20	3.290	678	684	686	rBV	409	304	0.00%	0.000%
21	3.331	693	697	701	rBV3	283	285	0.00%	0.000%
22	3.569	768	771	776	rVV2	370	382	0.00%	0.000%
23	3.653	792	797	800	rVB3	361	349	0.00%	0.000%
24	3.794	821	841	860	rVV2	87317	226430	0.06%	0.024%
25	3.939	870	886	909	rBV2	67394	255999	0.07%	0.028%
26	4.380	1005	1023	1056	rBV3	228512	671506	0.17%	0.072%
27	4.572	1081	1083	1089	rVB2	690	330	0.00%	0.000%
28	4.708	1108	1125	1144	rVV2	52435	138928	0.04%	0.015%
29	4.804	1150	1155	1160	rBV4	346	401	0.00%	0.000%
30	4.827	1160	1162	1170	rVB3	689	358	0.00%	0.000%
31	4.878	1174	1178	1181	rVB3	657	404	0.00%	0.000%
32	4.910	1181	1188	1194	rBV2	772	744	0.00%	0.000%
33	4.978	1206	1209	1210	rBV3	533	357	0.00%	0.000%
34	5.209	1222	1281	1322	rBV3	75397213	389727645	100.00%	42.010%

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 Title : VOC Analysis

35	5.650	1406	1418	1440	rVB	460915	895961	0.23%	0.097%
36	5.942	1499	1509	1522	rBV2	23785	46763	0.01%	0.005%
37	6.100	1544	1558	1573	rBV	291523	600274	0.15%	0.065%
38	6.318	1624	1626	1629	rBV3	584	284	0.00%	0.000%
39	6.367	1639	1641	1642	rBV	562	249	0.00%	0.000%
40	6.383	1642	1646	1647	rBV2	1935	1354	0.00%	0.000%
41	6.425	1657	1659	1661	rBV3	664	394	0.00%	0.000%
42	6.466	1667	1672	1673	rBV2	693	476	0.00%	0.000%
43	6.502	1679	1683	1686	rVB3	700	552	0.00%	0.000%
44	6.569	1701	1704	1706	rBV2	689	473	0.00%	0.000%
45	6.633	1719	1724	1727	rBV3	905	855	0.00%	0.000%
46	6.659	1727	1732	1734	rBV3	885	742	0.00%	0.000%
47	6.717	1748	1750	1752	rBV2	726	397	0.00%	0.000%
48	6.823	1773	1783	1784	rBV5	5711	6640	0.00%	0.001%
49	7.013	1828	1842	1862	rBV	170314	310020	0.08%	0.033%
50	7.135	1878	1880	1881	rBV2	624	315	0.00%	0.000%
51	7.251	1911	1916	1917	rBV2	448	393	0.00%	0.000%
52	7.341	1930	1944	1959	rBV	681238	1217091	0.31%	0.131%
53	7.415	1960	1967	1988	rVB	70771	129309	0.03%	0.014%
54	7.592	2020	2022	2028	rBV6	486	439	0.00%	0.000%
55	7.646	2028	2039	2056	rBV	118442	207606	0.05%	0.022%
56	7.749	2068	2071	2076	rVB5	870	631	0.00%	0.000%
57	7.791	2082	2084	2086	rBV2	610	244	0.00%	0.000%
58	7.894	2114	2116	2118	rVB2	566	249	0.00%	0.000%
59	7.920	2119	2124	2127	rBV4	582	439	0.00%	0.000%
60	7.942	2127	2131	2132	rBV2	601	470	0.00%	0.000%
61	7.961	2133	2137	2139	rVV3	2071	1551	0.00%	0.000%
62	8.003	2139	2150	2164	rVB3	44913	79034	0.02%	0.009%
63	8.087	2170	2176	2177	rBV3	731	692	0.00%	0.000%
64	8.132	2178	2190	2225	rVV2	300146	638160	0.16%	0.069%
65	8.312	2243	2246	2253	rVB7	770	617	0.00%	0.000%
66	8.376	2264	2266	2268	rVB3	654	280	0.00%	0.000%
67	8.444	2285	2287	2293	rVB5	924	733	0.00%	0.000%
68	8.495	2300	2303	2305	rBV2	566	279	0.00%	0.000%
69	8.646	2348	2350	2352	rBV2	504	277	0.00%	0.000%
70	8.727	2370	2375	2378	rBV2	308	265	0.00%	0.000%
71	10.241	2833	2846	2870	rBV2	568678	904446	0.23%	0.097%

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72	10.405	2894	2897	2900	rBV2	2152	1357	0.00%	0.000%
73	10.450	2902	2911	2926	rVB	63494	100485	0.03%	0.011%
74	10.566	2938	2947	2962	rVB	41669	68616	0.02%	0.007%
75	10.723	2993	2996	3000	rBV3	905	777	0.00%	0.000%
76	10.871	3040	3042	3046	rBV4	729	274	0.00%	0.000%
77	10.913	3053	3055	3057	rBV3	505	285	0.00%	0.000%
78	10.942	3060	3064	3066	rBV	1319	1195	0.00%	0.000%
79	11.010	3082	3085	3087	rBV3	790	476	0.00%	0.000%
80	11.071	3102	3104	3106	rBV3	550	332	0.00%	0.000%
81	11.209	3131	3147	3161	rBV	9304433	16584985	4.26%	1.788%
82	11.337	3161	3187	3201	rVB5	108168793	292756995	75.12%	31.557%
83	11.701	3274	3300	3334	rBV4	86525779	198212283	50.86%	21.366%
84	12.135	3430	3435	3448	rVB8	5480	8523	0.00%	0.001%
85	12.418	3513	3523	3536	rBV	86008	130915	0.03%	0.014%
86	12.601	3572	3580	3594	rVB3	13620	29366	0.01%	0.003%
87	12.669	3597	3601	3605	rVB5	2474	2406	0.00%	0.000%
88	12.772	3622	3633	3643	rBV2	36336	57867	0.01%	0.006%
89	13.026	3703	3712	3720	rVB8	3069	5302	0.00%	0.001%
90	13.286	3776	3793	3818	rBV	14320733	21749031	5.58%	2.344%
91	13.527	3860	3868	3879	rVB	14974	23321	0.01%	0.003%
92	13.765	3930	3942	3962	rVB	625628	968017	0.25%	0.104%
93	13.878	3975	3977	3981	rVB4	1262	732	0.00%	0.000%
94	13.926	3988	3992	3993	rBV2	704	573	0.00%	0.000%
95	14.009	4016	4018	4019	rBV2	902	329	0.00%	0.000%
96	14.099	4043	4046	4050	rBV3	959	637	0.00%	0.000%
97	14.382	4132	4134	4136	rVB2	729	326	0.00%	0.000%
98	14.440	4149	4152	4155	rBV4	1265	760	0.00%	0.000%
99	14.456	4155	4157	4162	rVB4	895	474	0.00%	0.000%
100	14.881	4287	4289	4291	rBV3	773	403	0.00%	0.000%

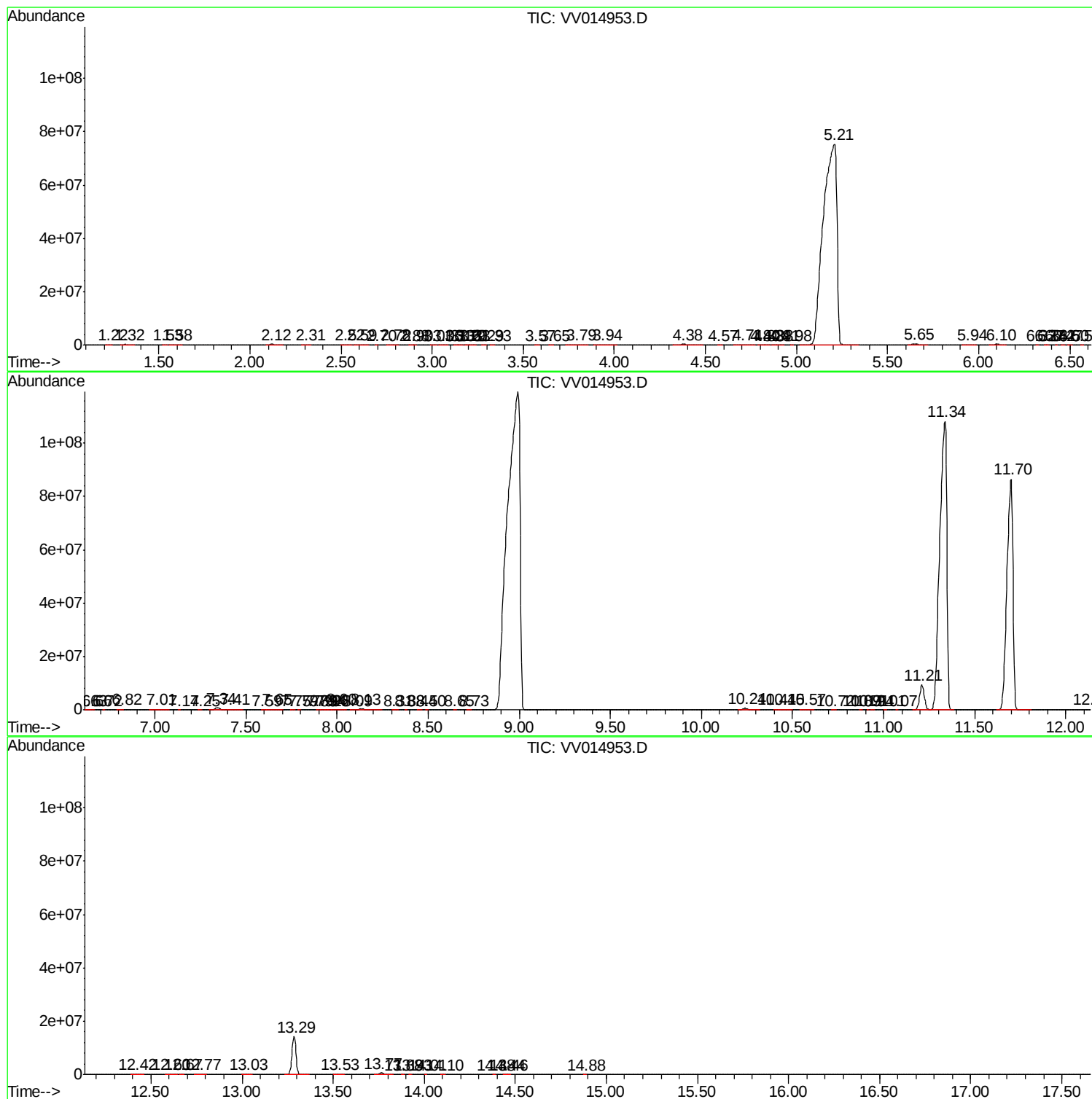
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM031620WMA.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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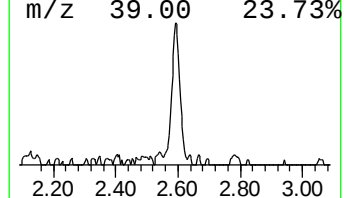
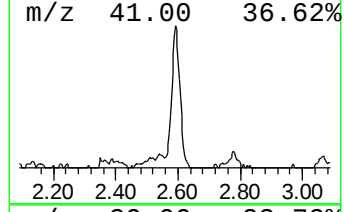
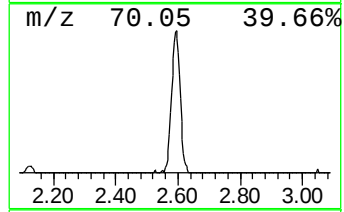
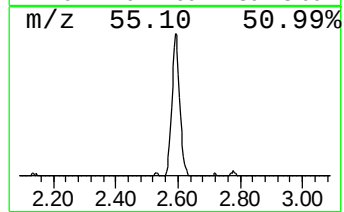
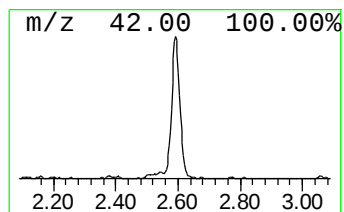
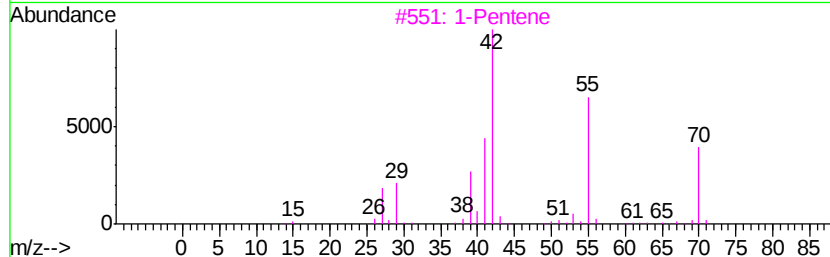
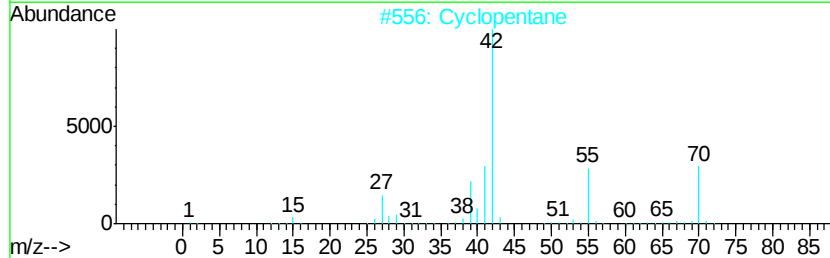
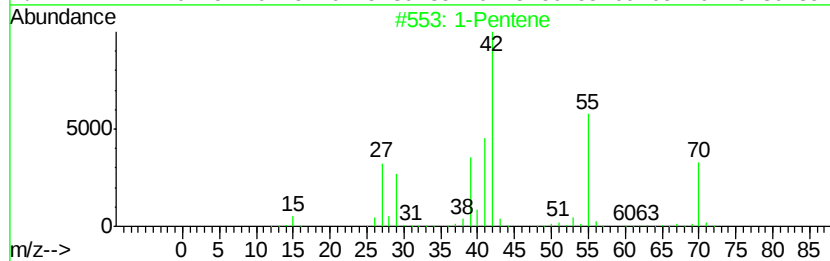
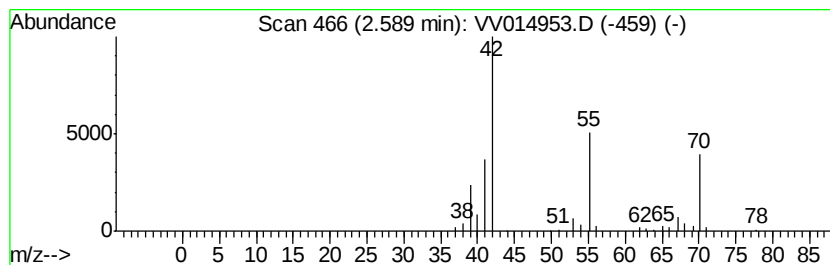
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM031620WMA.M
Quant Title : VOC Analysis

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

Peak Number 1 (DEL) Alkane: Cyclic2.59 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.59	3.78 ug/L	67645	1,4-Difluorobenzene	5.65

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Pentene	70	C5H10	000109-67-1	72
2			Cyclopentane	70	C5H10	000287-92-3	72
3			1-Pentene	70	C5H10	000109-67-1	64
4			Cyclopentane	70	C5H10	000287-92-3	56
5			Cyclopropane, ethyl-	70	C5H10	001191-96-4	32



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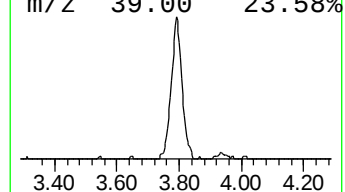
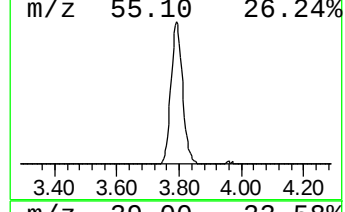
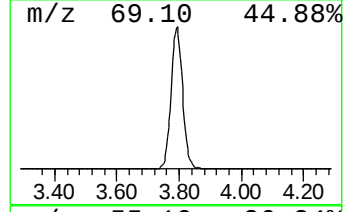
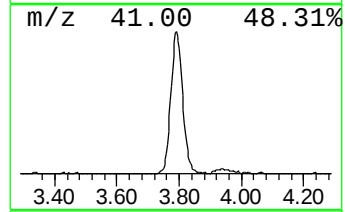
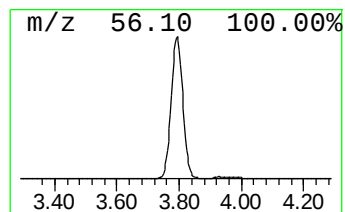
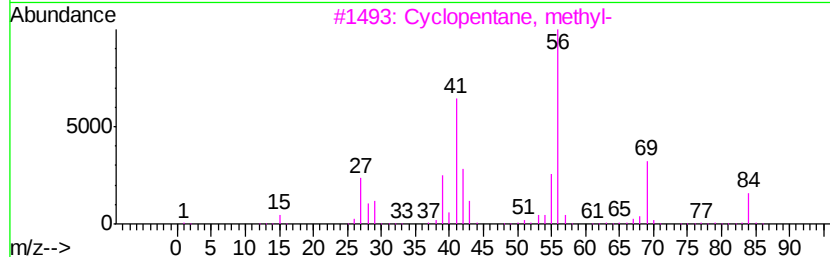
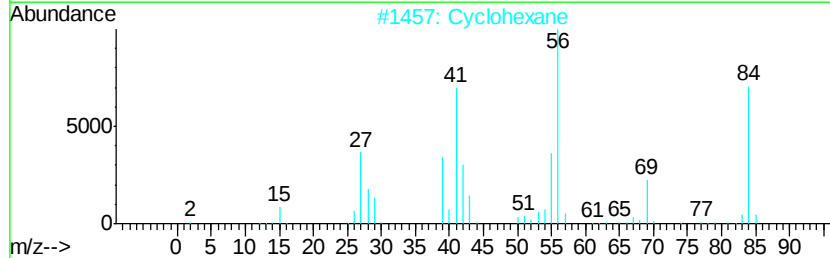
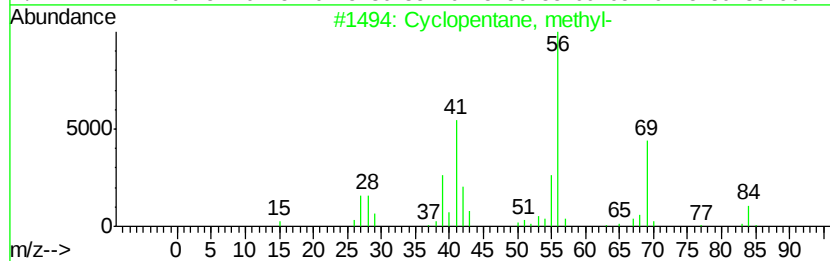
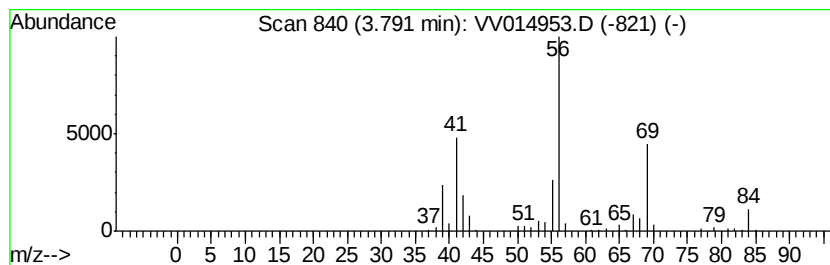
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM031620WMA.M
Quant Title : VOC Analysis

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

Peak Number 2 (DEL) Alkane: Cyclic3.79 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.79	12.64 ug/L	226430	1,4-Difluorobenzene	5.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	90
2		Cyclohexane	84	C6H12	000110-82-7	83
3		Cyclopentane, methyl-	84	C6H12	000096-37-7	80
4		Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
5		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	59



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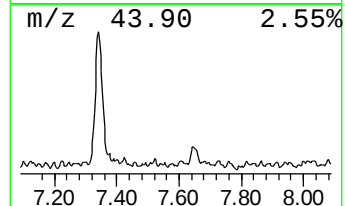
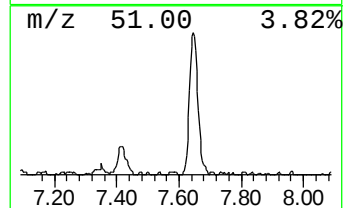
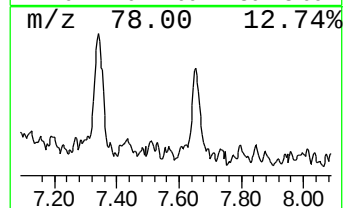
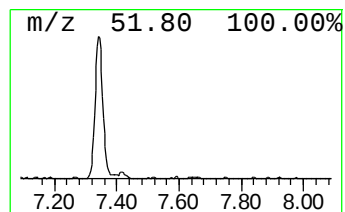
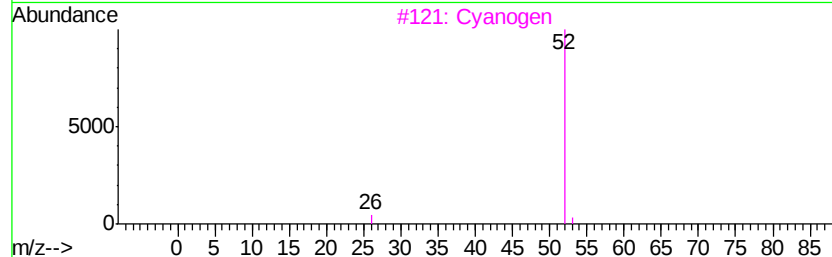
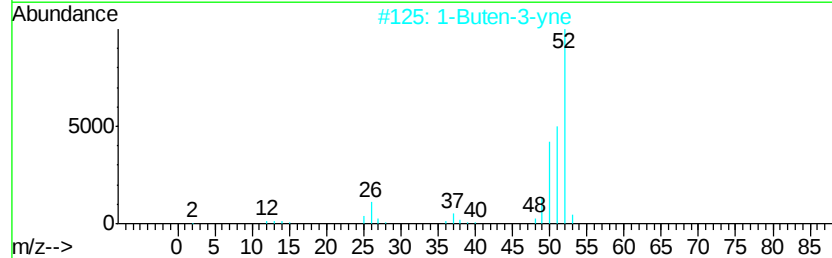
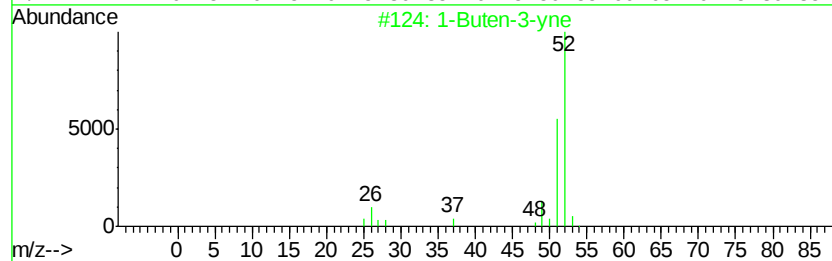
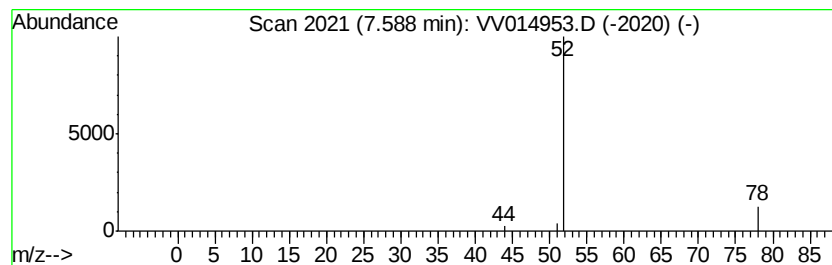
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Quant Title : VOC Analysis

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

Peak Number 5 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	82.83 ug/L	439	Chlorobenzene-d5	8.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Buten-3-yne	52	C4H4	000689-97-4	3
2		1-Buten-3-yne	52	C4H4	000689-97-4	3
3		Cyanoden	52	C2N2	000460-19-5	3
4		1-Buten-3-yne	52	C4H4	000689-97-4	3
5		1-Buten-3-yne	52	C4H4	000689-97-4	3



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ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0DH5

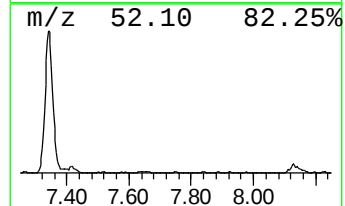
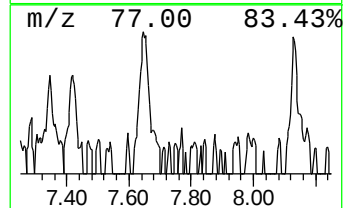
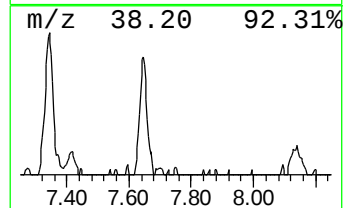
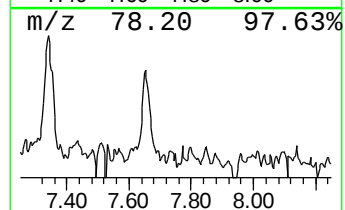
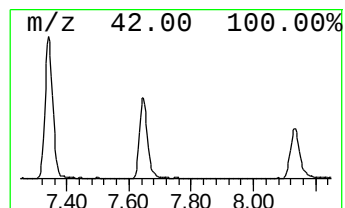
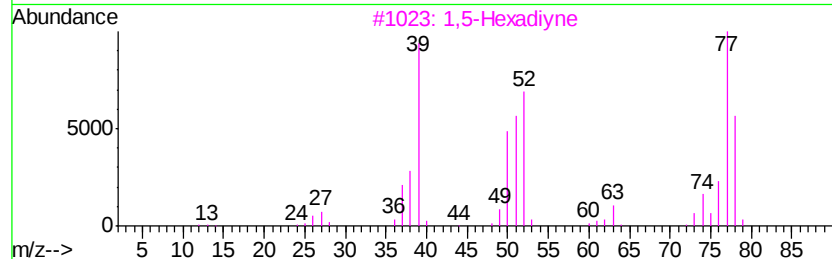
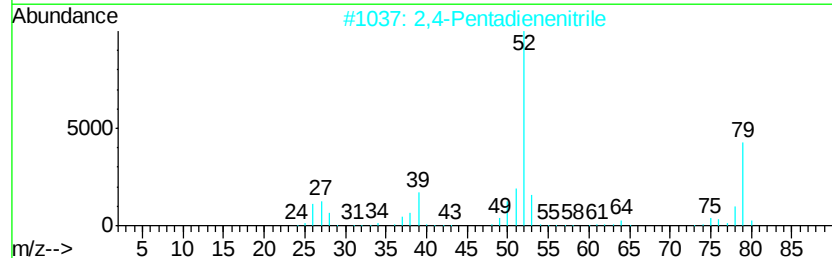
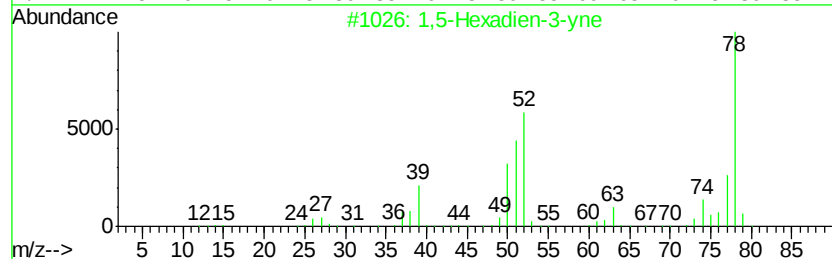
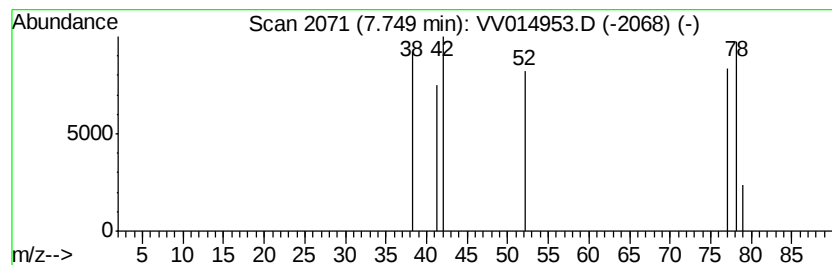
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Quant Title : VOC Analysis

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

Peak Number 6 unknown-02 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.75	119.06 ug/L	631	Chlorobenzene-d5	8.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,5-Hexadien-3-yne	78	C6H6	000821-08-9	7
2		2,4-Pentadienenitrile	79	C5H5N	001615-70-9	5
3		1,5-Hexadiyne	78	C6H6	000628-16-0	5
4		2,4-Pentadienenitrile	79	C5H5N	001615-70-9	5
5		Propanedinitrile, methylene-	78	C4H2N2	000922-64-5	3



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV031620\
Data File : VV014953.D
Acq On : 16 Mar 2020 18:07
Operator : SY/MD
Sample : L1890-18
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0DH5

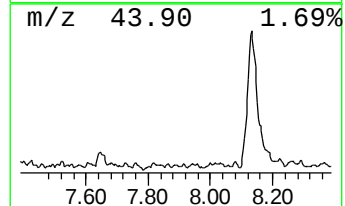
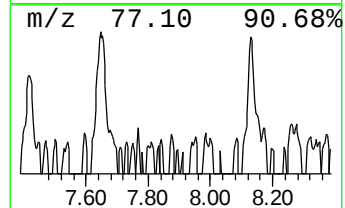
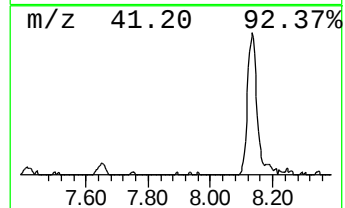
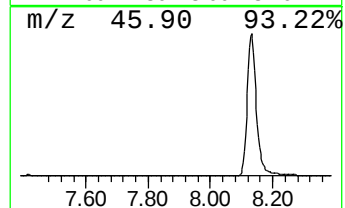
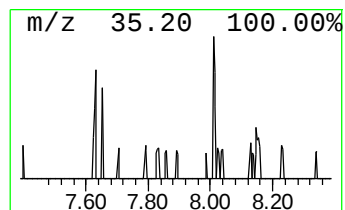
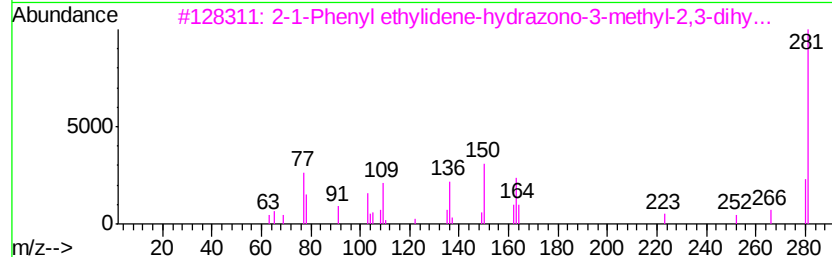
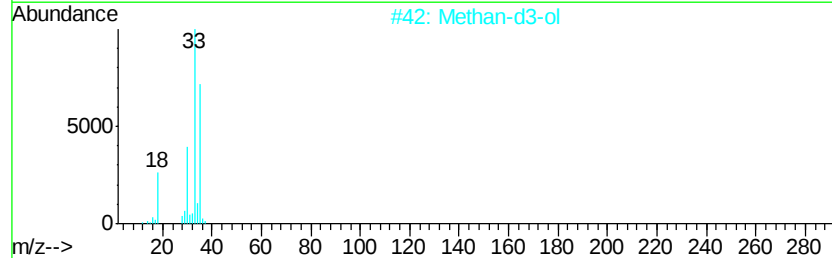
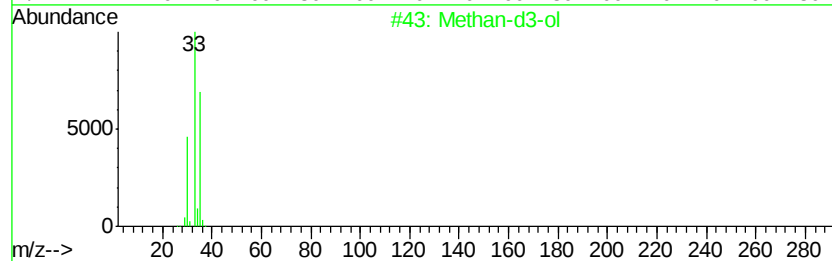
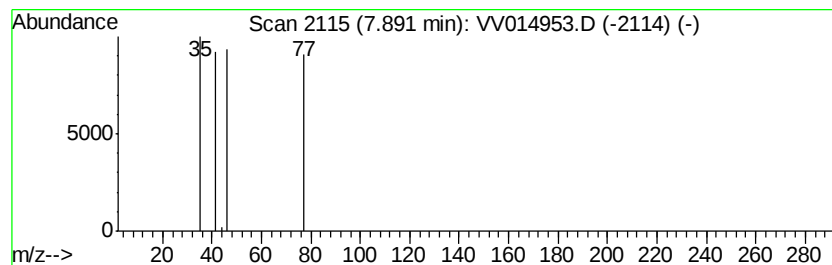
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Quant Title : VOC Analysis

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

Peak Number 8 unknown-03 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.89	46.98 ug/L	249	Chlorobenzene-d5	8.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methan-d3-ol	35	CHD3O	001849-29-2	3
2		Methan-d3-ol	35	CHD3O	001849-29-2	3
3		2-1-Phenyl ethylidene-hydrazono-...	281	C16H15N3S	1000148-55-3	3
4		4-Bromo-2,6-difluorobenzyl alcoh...	294	C10H13BrF2OSi	1000364-60-0	2
5		6,8-Dioxa-3-thia-bicyclo(3,2,1)o...	132	C5H8O2S	1000143-27-3	2



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV031620\
Data File : VV014953.D
Acq On : 16 Mar 2020 18:07
Operator : SY/MD
Sample : L1890-18
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0DH5

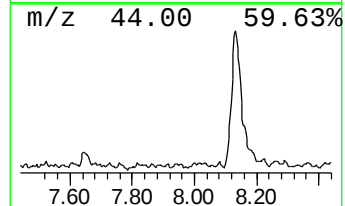
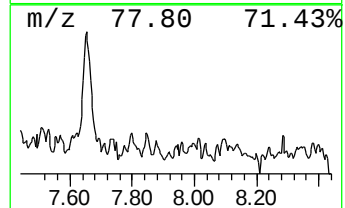
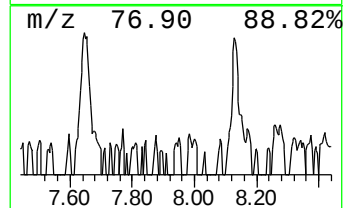
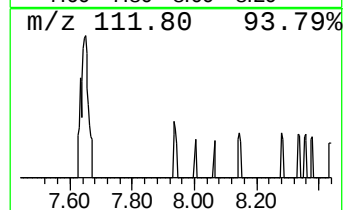
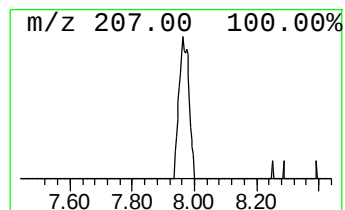
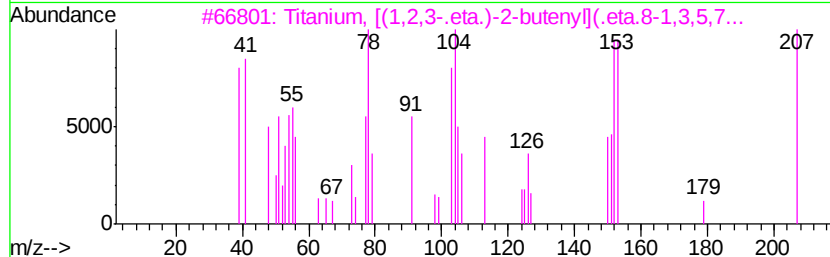
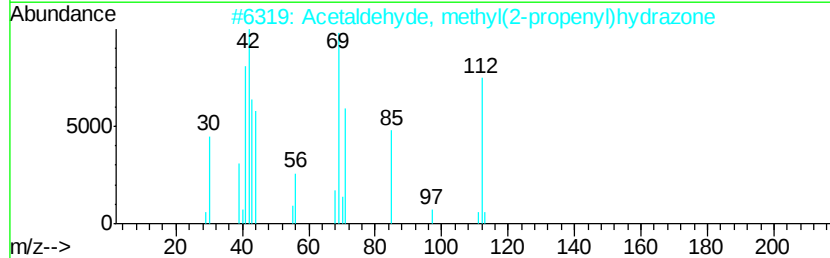
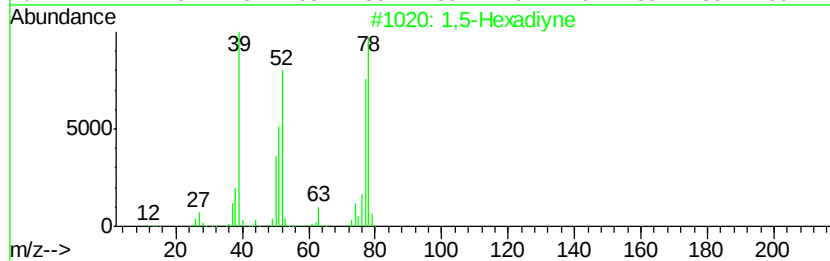
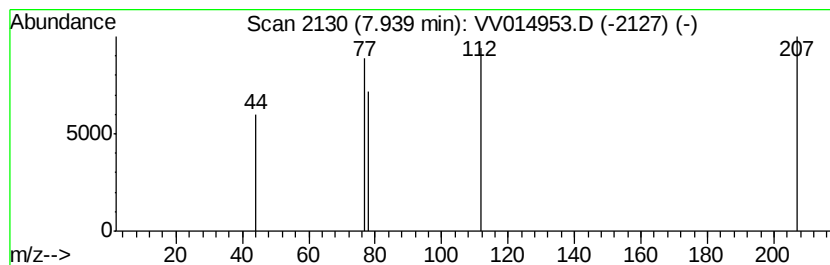
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Quant Title : VOC Analysis

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

Peak Number 10 unknown-04 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.94	88.68 ug/L	470	Chlorobenzene-d5	8.94

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,5-Hexadiyne	78	C6H6	000628-16-0	5
2			Acetaldehyde, methyl(2-propenyl)...	112	C6H12N2	066075-10-3	5
3			Titanium, [(1,2,3-.eta.)-2-buten...	207	C12H15Ti	039015-02-6	5
4			5-Amino-3,4-dimethyl-isoxazole	112	C5H8N2O	019947-75-2	4
5			Hydrazine, 1,1-di-2-propenyl-	112	C6H12N2	005164-11-4	4



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV031620\
Data File : VV014953.D
Acq On : 16 Mar 2020 18:07
Operator : SY/MD
Sample : L1890-18
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0DH5

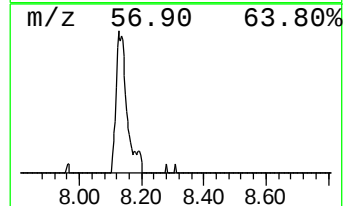
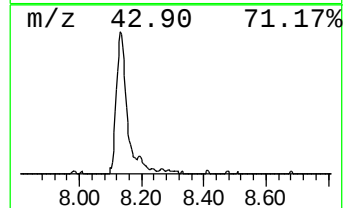
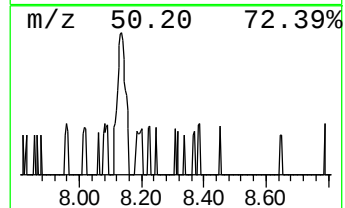
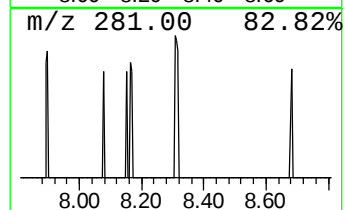
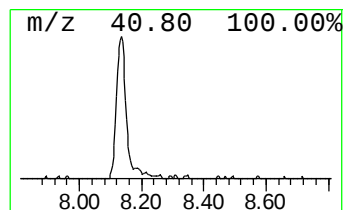
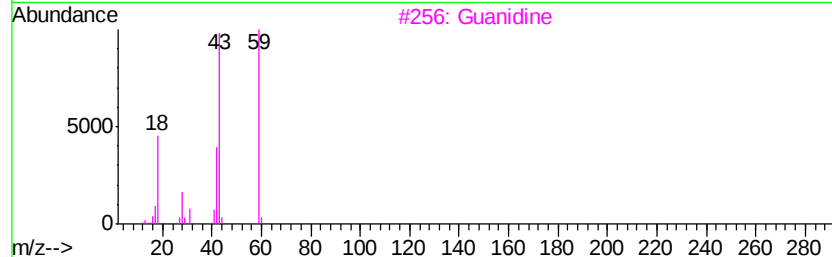
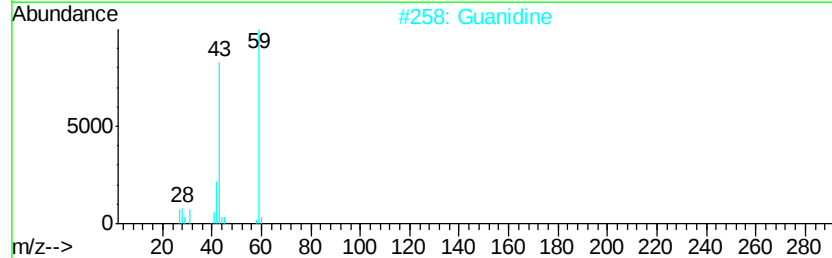
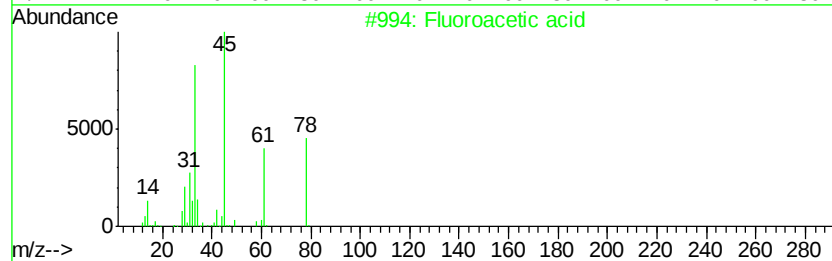
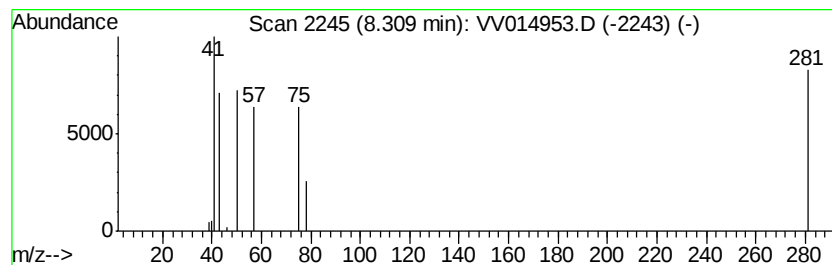
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Quant Title : VOC Analysis

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

Peak Number 11 unknown-05 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.31	116.42 ug/L	617	Chlorobenzene-d5	8.94

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Fluoroacetic acid	78	C2H3FO2	000144-49-0	7
2		Guanidine	59	CH5N3	000113-00-8	5
3		Guanidine	59	CH5N3	000113-00-8	5
4		Guanidine	59	CH5N3	000113-00-8	5
5		Acetaldoxime	59	C2H5NO	000107-29-9	4



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV031620\
Data File : VV014953.D
Acq On : 16 Mar 2020 18:07
Operator : SY/MD
Sample : L1890-18
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0DH5

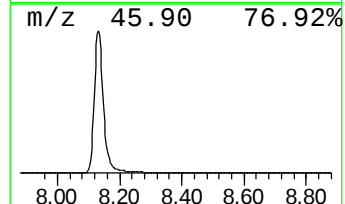
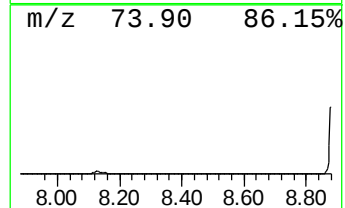
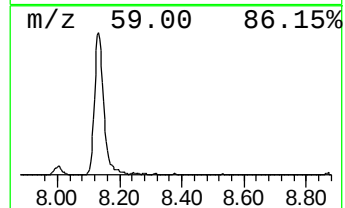
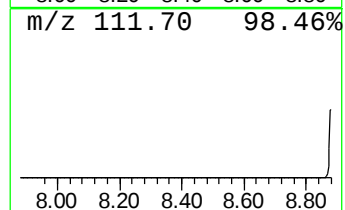
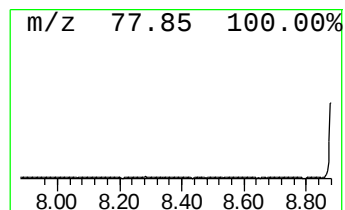
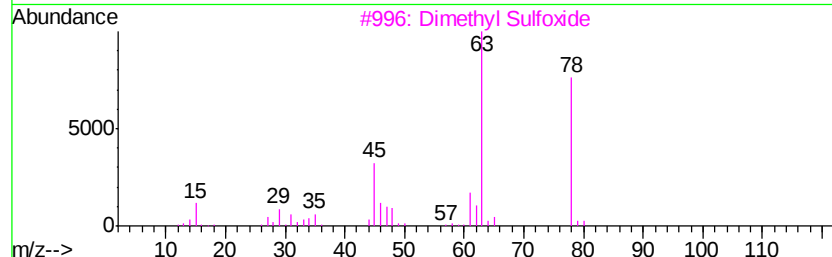
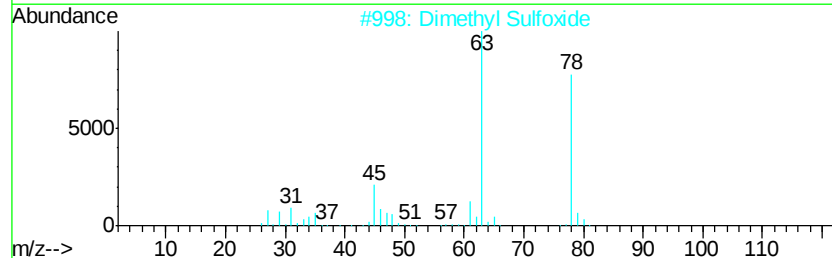
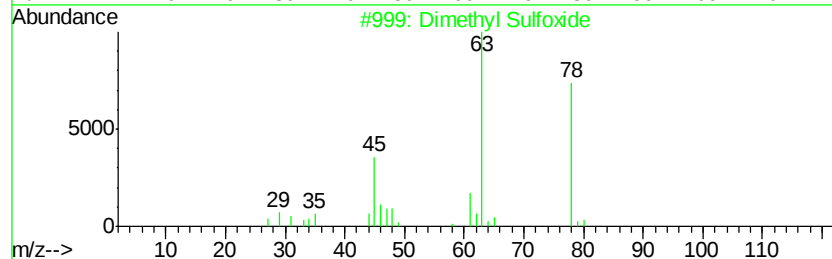
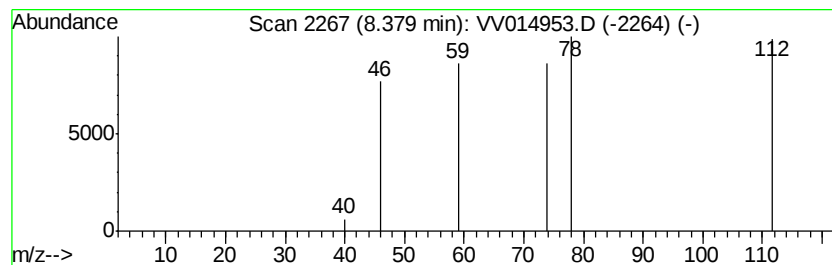
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Quant Title : VOC Analysis

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

Peak Number 12 unknown-06 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.38	52.83 ug/L	280	Chlorobenzene-d5	8.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dimethyl Sulfoxide	78	C2H6OS	000067-68-5	7
2		Dimethyl Sulfoxide	78	C2H6OS	000067-68-5	7
3		Dimethyl Sulfoxide	78	C2H6OS	000067-68-5	7
4		Dimethyl Sulfoxide	78	C2H6OS	000067-68-5	5
5		Thiirane, methyl-	74	C3H6S	001072-43-1	5



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV031620\
Data File : VV014953.D
Acq On : 16 Mar 2020 18:07
Operator : SY/MD
Sample : L1890-18
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0DH5

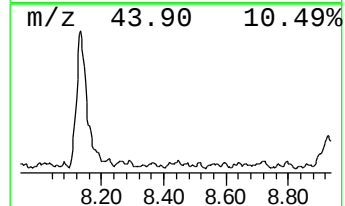
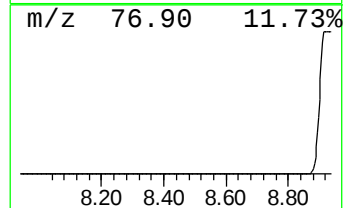
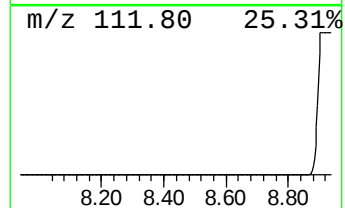
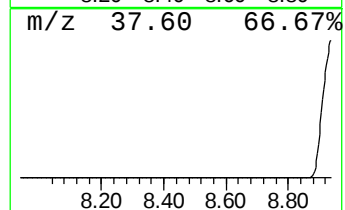
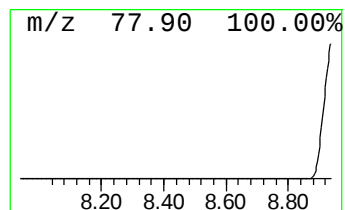
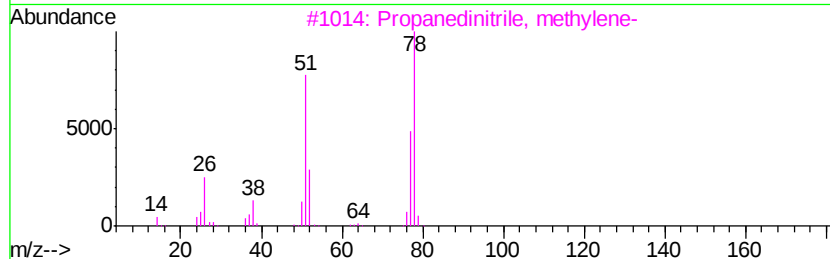
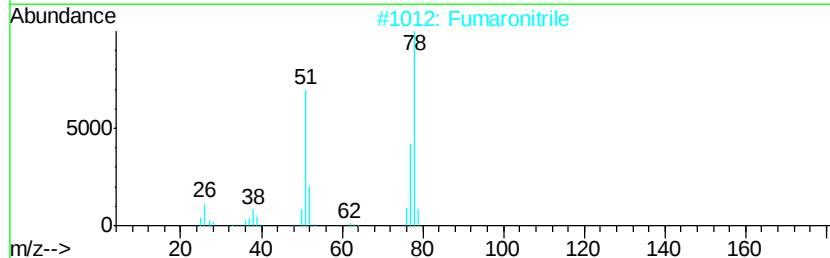
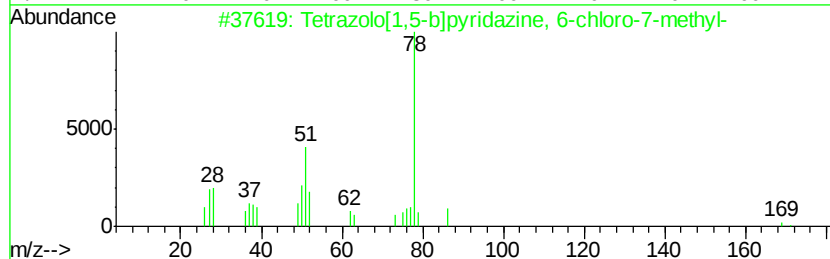
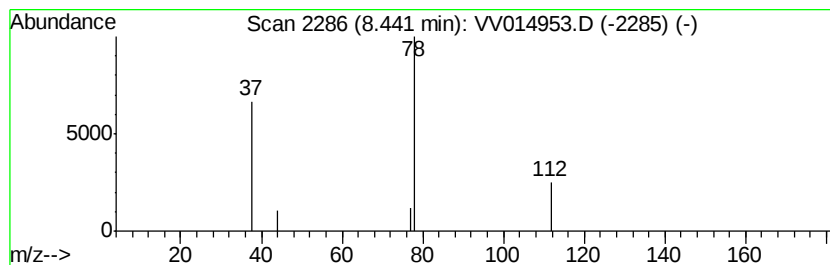
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Quant Title : VOC Analysis

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

Peak Number 13 unknown-07 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.44	138.30 ug/L	733	Chlorobenzene-d5	8.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetrazolo[1,5-b]pyridazine, 6-ch...	169	C5H4ClN5	028691-22-7	9
2		Fumaronitrile	78	C4H2N2	000764-42-1	5
3		Propanedinitrile, methylene-	78	C4H2N2	000922-64-5	5
4		Fumaronitrile	78	C4H2N2	000764-42-1	5
5		1,5-Hexadiyne	78	C6H6	000628-16-0	4



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV031620\
Data File : VV014953.D
Acq On : 16 Mar 2020 18:07
Operator : SY/MD
Sample : L1890-18
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0DH5

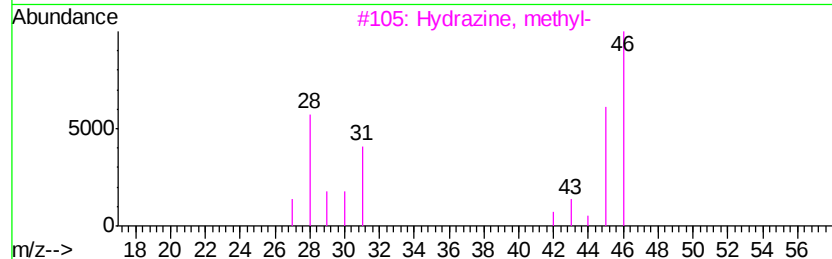
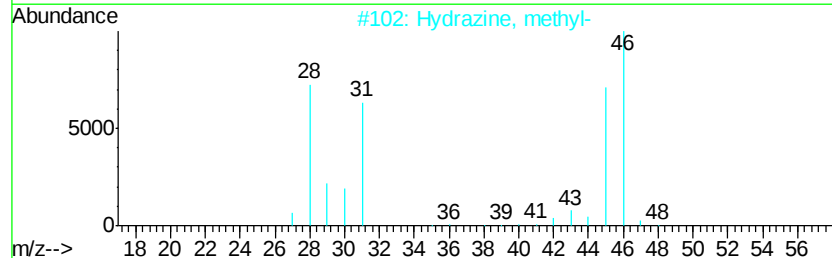
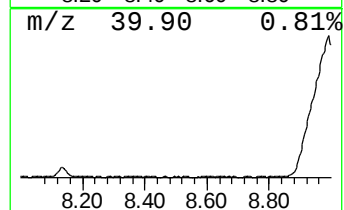
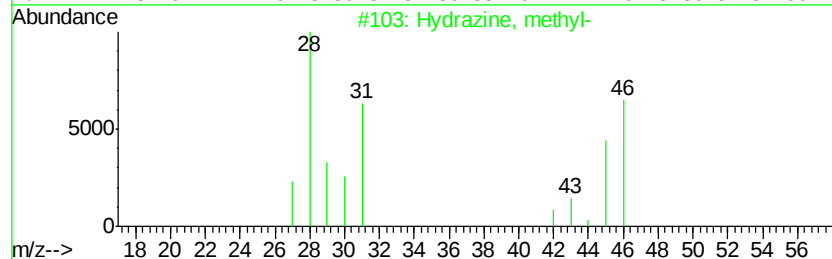
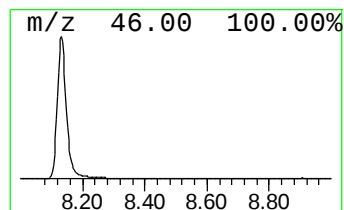
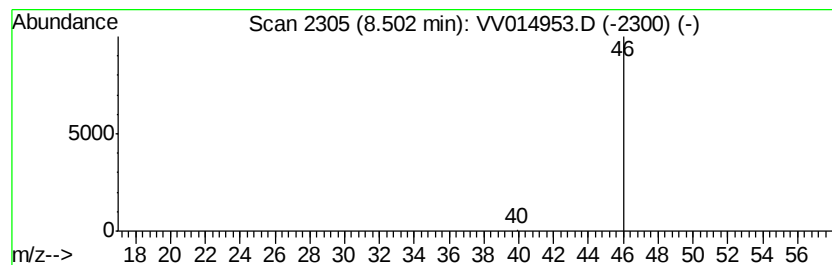
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Quant Title : VOC Analysis

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

Peak Number 14 unknown-08 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.50	52.64 ug/L	279	Chlorobenzene-d5	8.94

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hvdrazine, methyl-	46	CH6N2	000060-34-4	4
2			Hvdrazine, methyl-	46	CH6N2	000060-34-4	4
3			Hvdrazine, methyl-	46	CH6N2	000060-34-4	4
4			Mercaptamine	77	C2H7NS	000060-23-1	3
5			Silanamine, N-silyl-	77	H7NSi2	005702-11-4	3



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV031620\
Data File : VV014953.D
Acq On : 16 Mar 2020 18:07
Operator : SY/MD
Sample : L1890-18
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0DH5

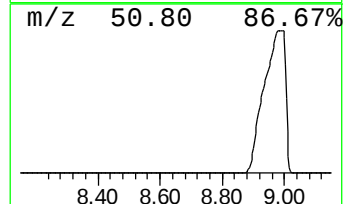
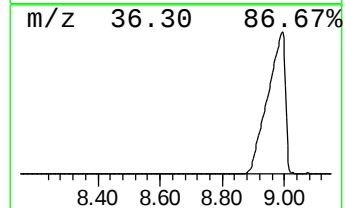
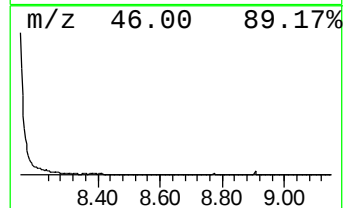
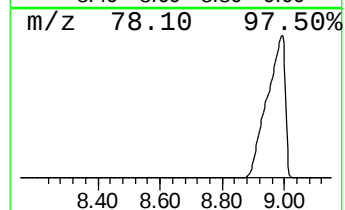
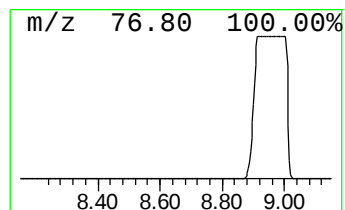
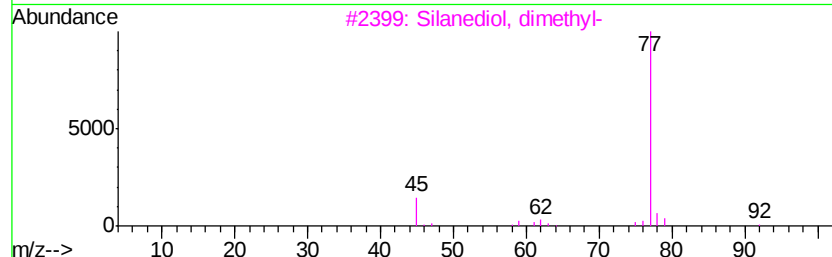
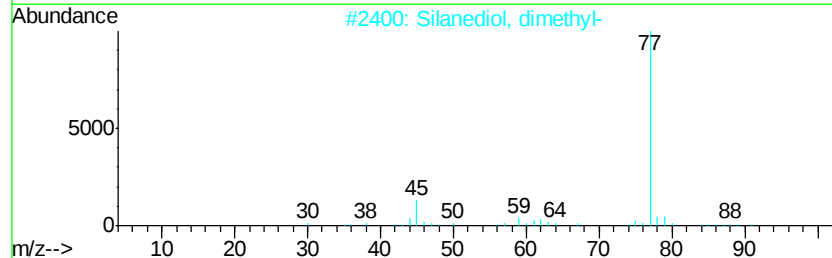
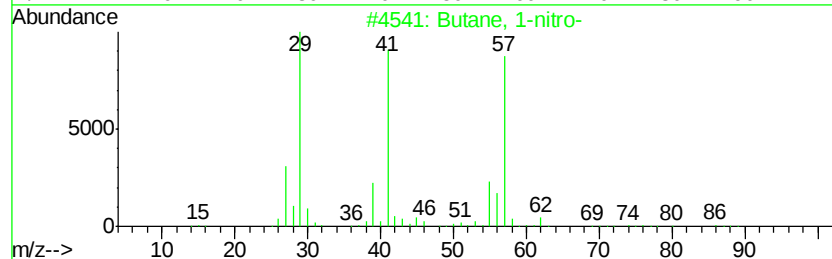
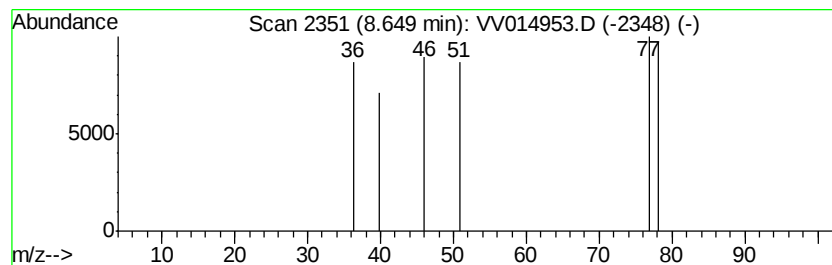
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Quant Title : VOC Analysis

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

Peak Number 15 unknown-09 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.65	52.26 ug/L	277	Chlorobenzene-d5	8.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 1-nitro-	103	C4H9NO2	000627-05-4	1
2		Silanediol, dimethyl-	92	C2H8O2Si	001066-42-8	1
3		Silanediol, dimethyl-	92	C2H8O2Si	001066-42-8	1
4		1,5-Hexadiene, 3,3,4,4-tetrafluoro-	154	C6H6F4	001763-21-9	1
5		2-Chloroethylamine	79	C2H6ClN	000689-98-5	1



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV031620\
Data File : VV014953.D
Acq On : 16 Mar 2020 18:07
Operator : SY/MD
Sample : L1890-18
Misc : 5.0mL/MSVOA V/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0DH5

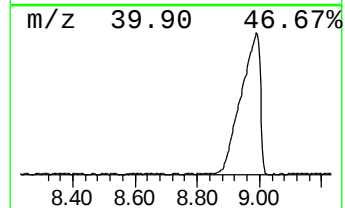
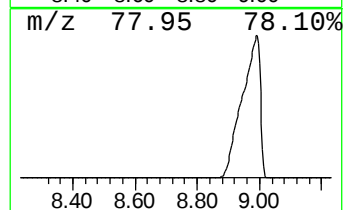
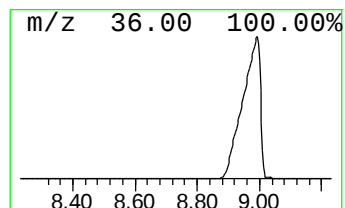
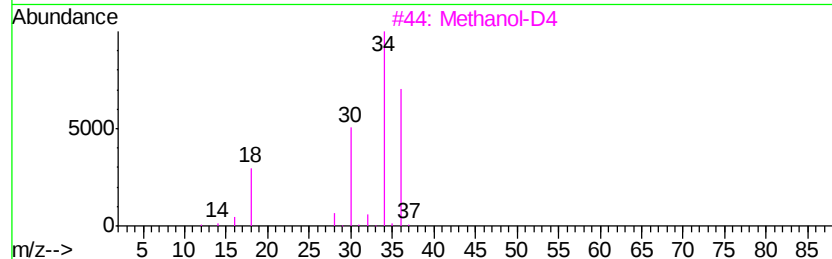
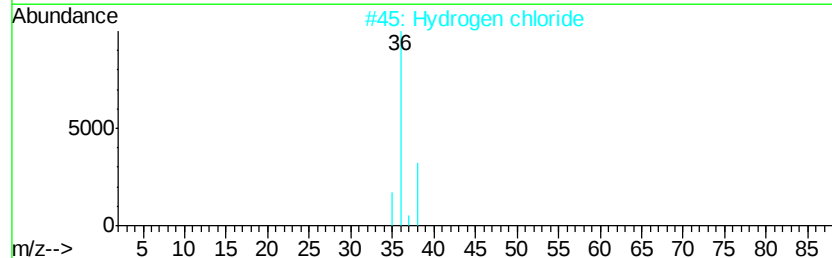
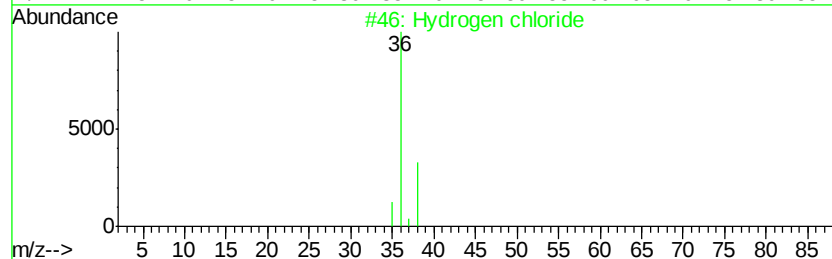
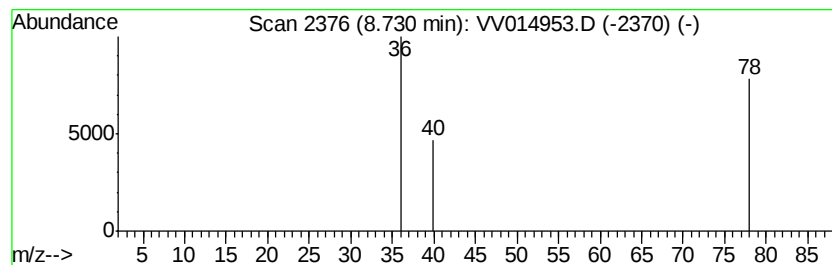
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM031620WMA.M
Quant Title : VOC Analysis

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

Peak Number 16 unknown-10 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.73	50.00 ug/L	265	Chlorobenzene-d5	8.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hydrogen chloride	36	ClH	007647-01-0	2
2		Hydrogen chloride	36	ClH	007647-01-0	2
3		Methanol-D4	36	CD4O	000811-98-3	2
4		Ammonium Chloride	53	ClH4N	012125-02-9	1
5		2-Chloroethylamine	79	C2H6ClN	000689-98-5	1



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV031620\
Data File : VV014953.D
Acq On : 16 Mar 2020 18:07
Operator : SY/MD
Sample : L1890-18
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0DH5

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM031620WMA.M
Quant Title : VOC Analysis

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: Cyc...	2.59	3.8	ug/L	67645	1	5.65	895961	50.0
(DEL) Alkane: Cyc...	3.79	12.6	ug/L	226430	1	5.65	895961	50.0
unknown-01	7.59	82.8	ug/L	439	2	8.94	265	50.0
unknown-02	7.75	119.1	ug/L	631	2	8.94	265	50.0
unknown-03	7.89	47.0	ug/L	249	2	8.94	265	50.0
unknown-04	7.94	88.7	ug/L	470	2	8.94	265	50.0
unknown-05	8.31	116.4	ug/L	617	2	8.94	265	50.0
unknown-06	8.38	52.8	ug/L	280	2	8.94	265	50.0
unknown-07	8.44	138.3	ug/L	733	2	8.94	265	50.0
unknown-08	8.50	52.6	ug/L	279	2	8.94	265	50.0
unknown-09	8.65	52.3	ug/L	277	2	8.94	265	50.0
unknown-10	8.73	50.0	ug/L	265	2	8.94	265	50.0