

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW121719\
 Data File : VW014414.D
 Acq On : 18 Dec 2019 01:20
 Operator : SY/VA
 Sample : K6279-10
 Misc : 3.99G/10ML/MSVOA W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 BFEY4

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_W\METHOD\SOM2WLM120519S.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	2.160	38	42	47	rVB2	7984	11845	0.51%	0.238%
2	2.349	69	73	79	rVB	32949	50848	2.19%	1.020%
3	2.495	92	97	102	rVB	49009	78843	3.40%	1.582%
4	3.391	237	244	251	rVB3	6269	14585	0.63%	0.293%
5	3.526	257	266	293	rBV	977127	2320276	100.00%	46.561%
6	4.013	337	346	357	rBV2	52596	134652	5.80%	2.702%
7	4.202	371	377	388	rVB2	20687	59426	2.56%	1.193%
8	4.550	432	434	436	rVB2	899	808	0.03%	0.016%
9	4.574	436	438	439	rBV2	801	529	0.02%	0.011%
10	4.598	439	442	443	rBV3	762	559	0.02%	0.011%
11	4.794	473	474	475	rBV	871	607	0.03%	0.012%
12	4.848	479	483	484	rVV3	638	844	0.04%	0.017%
13	4.915	486	494	503	rVV7	5848	18668	0.80%	0.375%
14	5.007	507	509	510	rBV	1035	539	0.02%	0.011%
15	5.074	519	520	522	rVV2	806	553	0.02%	0.011%
16	5.226	543	545	548	rBV3	548	670	0.03%	0.013%
17	5.458	581	583	585	rBV2	821	674	0.03%	0.014%
18	5.495	585	589	592	rBV5	641	826	0.04%	0.017%
19	5.586	601	604	608	rBV	490	721	0.03%	0.014%
20	5.629	608	611	614	rBV3	532	822	0.04%	0.016%
21	5.714	623	625	626	rBV2	763	784	0.03%	0.016%
22	5.787	629	637	645	rVV4	6547	20160	0.87%	0.405%
23	6.117	690	691	693	rBV	777	652	0.03%	0.013%
24	6.171	698	700	702	rVB2	672	552	0.02%	0.011%
25	6.202	704	705	707	rBV	568	495	0.02%	0.010%
26	6.269	712	716	718	rBV2	688	843	0.04%	0.017%
27	6.287	718	719	723	rVB3	946	931	0.04%	0.019%
28	6.324	723	725	726	rBV	920	532	0.02%	0.011%
29	6.354	726	730	731	rBV2	483	604	0.03%	0.012%
30	6.397	731	737	745	rBV7	3445	10692	0.46%	0.215%
31	6.525	756	758	761	rBV4	810	1033	0.04%	0.021%
32	6.549	761	762	764	rVV2	852	485	0.02%	0.010%
33	6.586	766	768	770	rBV3	714	794	0.03%	0.016%
34	6.671	781	782	786	rVB2	533	609	0.03%	0.012%

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Integration Parameters: LSCINT.P

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Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM120519S.M
 Title : VOC Analysis

35	6.775	796	799	800	rBV2	709	579	0.02%	0.012%
36	6.927	822	824	826	rVB2	733	550	0.02%	0.011%
37	7.000	828	836	837	rBV5	1201	2286	0.10%	0.046%
38	7.129	843	857	870	rBV3	24479	90452	3.90%	1.815%
39	7.269	878	880	882	rVB3	883	642	0.03%	0.013%
40	7.403	899	902	904	rBV2	823	1048	0.05%	0.021%
41	7.464	911	912	914	rVV	1058	657	0.03%	0.013%
42	7.647	932	942	954	rVV2	70228	172225	7.42%	3.456%
43	7.872	977	979	981	rVB2	632	509	0.02%	0.010%
44	7.897	981	983	984	rBV	805	619	0.03%	0.012%
45	8.049	1007	1008	1010	rBV	625	541	0.02%	0.011%
46	8.171	1026	1028	1032	rBV4	430	464	0.02%	0.009%
47	8.275	1032	1045	1058	rBV3	135012	406592	17.52%	8.159%
48	8.409	1065	1067	1069	rBV2	1268	1094	0.05%	0.022%
49	8.543	1084	1089	1091	rBV2	1439	2667	0.11%	0.054%
50	8.567	1091	1093	1095	rVB3	969	764	0.03%	0.015%
51	8.659	1106	1108	1111	rBV	512	494	0.02%	0.010%
52	8.707	1111	1116	1117	rBV3	3459	5356	0.23%	0.107%
53	8.842	1130	1138	1148	rVV	110726	224880	9.69%	4.513%
54	9.159	1189	1190	1193	rVB	827	546	0.02%	0.011%
55	9.274	1201	1209	1219	rBV	92628	198905	8.57%	3.991%
56	9.531	1249	1251	1253	rVB2	631	549	0.02%	0.011%
57	9.573	1255	1258	1264	rBV6	762	1119	0.05%	0.022%
58	9.634	1266	1268	1270	rBV2	621	532	0.02%	0.011%
59	9.848	1300	1303	1308	rVB	795	907	0.04%	0.018%
60	10.037	1323	1334	1342	rBV	31580	60780	2.62%	1.220%
61	10.128	1343	1349	1358	rVV	37955	73140	3.15%	1.468%
62	10.323	1374	1381	1389	rBV	145181	253967	10.95%	5.096%
63	10.433	1397	1399	1403	rVB4	1462	1519	0.07%	0.030%
64	10.488	1403	1408	1412	rBV5	1529	3436	0.15%	0.069%
65	10.579	1418	1423	1429	rVB	19839	34263	1.48%	0.688%
66	10.707	1438	1444	1448	rBV	16337	29920	1.29%	0.600%
67	10.793	1456	1458	1459	rVB	760	503	0.02%	0.010%
68	10.811	1459	1461	1462	rBV	572	493	0.02%	0.010%
69	10.933	1473	1481	1495	rBV2	19809	49598	2.14%	0.995%
70	11.061	1498	1502	1510	rVB2	10121	20404	0.88%	0.409%
71	11.244	1530	1532	1533	rBV	746	509	0.02%	0.010%

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 Title : VOC Analysis

72	11.311	1542	1543	1547	rVB2	809	822	0.04%	0.016%
73	11.439	1559	1564	1572	rVB4	5938	11075	0.48%	0.222%
74	11.494	1572	1573	1577	rBV3	944	570	0.02%	0.011%
75	11.628	1588	1595	1608	rVV	129298	232438	10.02%	4.664%
76	11.890	1635	1638	1640	rBV3	823	977	0.04%	0.020%
77	11.914	1640	1642	1645	rVB2	820	640	0.03%	0.013%
78	11.981	1651	1653	1656	rBV3	656	975	0.04%	0.020%
79	12.048	1658	1664	1666	rVV4	1187	1704	0.07%	0.034%
80	12.164	1681	1683	1685	rBV3	758	737	0.03%	0.015%
81	12.317	1706	1708	1711	rBV	677	821	0.04%	0.016%
82	12.554	1745	1747	1749	rBV	634	654	0.03%	0.013%
83	12.609	1749	1756	1763	rVV2	20040	35751	1.54%	0.717%
84	12.695	1763	1770	1780	rVB	62775	110056	4.74%	2.208%
85	12.768	1780	1782	1786	rBV3	983	1035	0.04%	0.021%
86	12.841	1791	1794	1796	rBV4	551	818	0.04%	0.016%
87	12.926	1805	1808	1814	rVB3	3407	5667	0.24%	0.114%
88	12.981	1814	1817	1821	rBV4	495	877	0.04%	0.018%
89	13.097	1834	1836	1838	rBV3	1140	1010	0.04%	0.020%
90	13.207	1849	1854	1859	rBV5	6212	11429	0.49%	0.229%
91	13.298	1864	1869	1873	rVB3	3334	5625	0.24%	0.113%
92	13.335	1873	1875	1877	rBV3	953	836	0.04%	0.017%
93	13.377	1880	1882	1884	rBV	507	478	0.02%	0.010%
94	13.408	1886	1887	1889	rVB2	1131	658	0.03%	0.013%
95	13.560	1906	1912	1922	rVB2	46815	81243	3.50%	1.630%
96	13.646	1924	1926	1929	rBV3	1074	978	0.04%	0.020%
97	13.768	1944	1946	1949	rVB2	888	647	0.03%	0.013%
98	13.853	1954	1960	1967	rBV	55710	93267	4.02%	1.872%
99	13.957	1975	1977	1979	rVB2	795	515	0.02%	0.010%
100	14.463	2058	2060	2062	rVB3	1407	1037	0.04%	0.021%

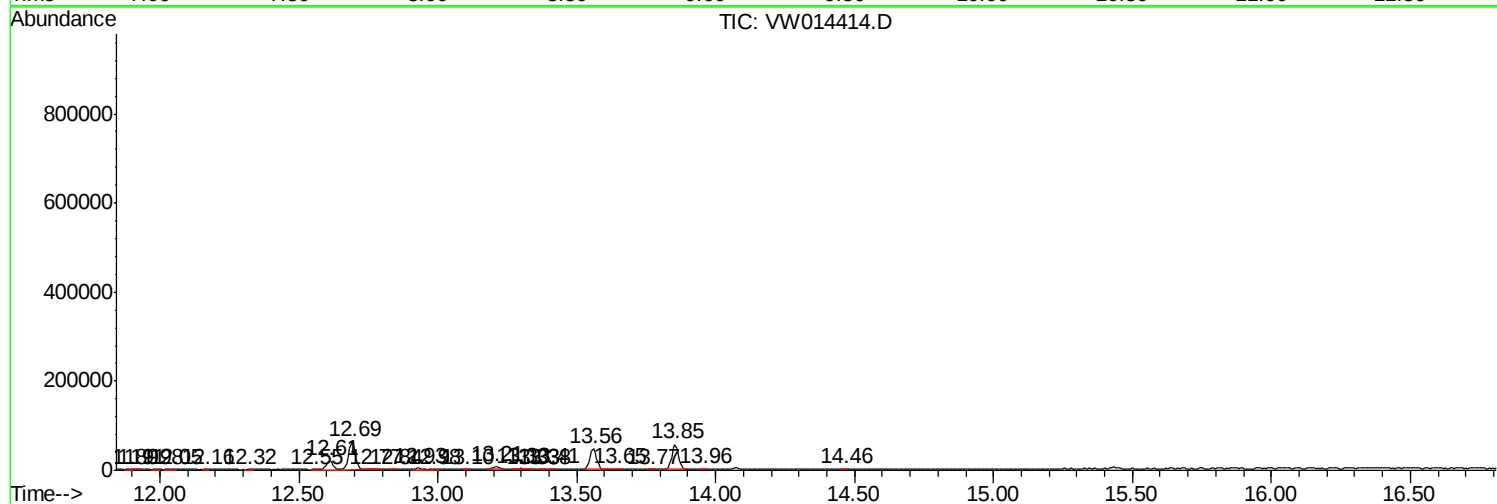
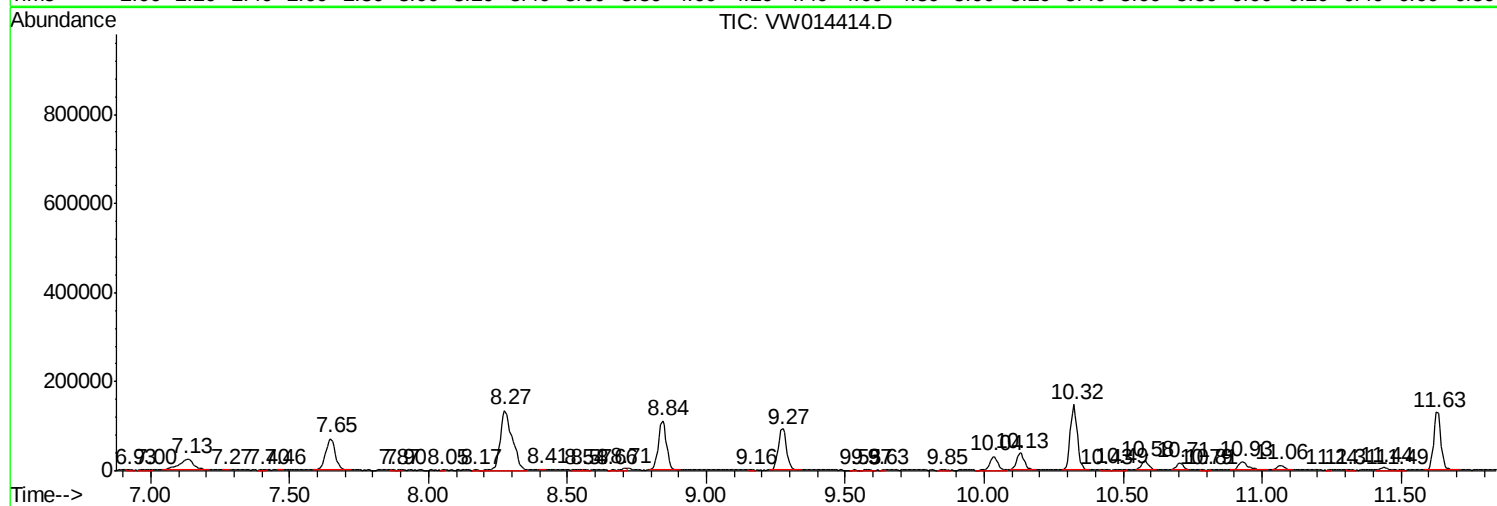
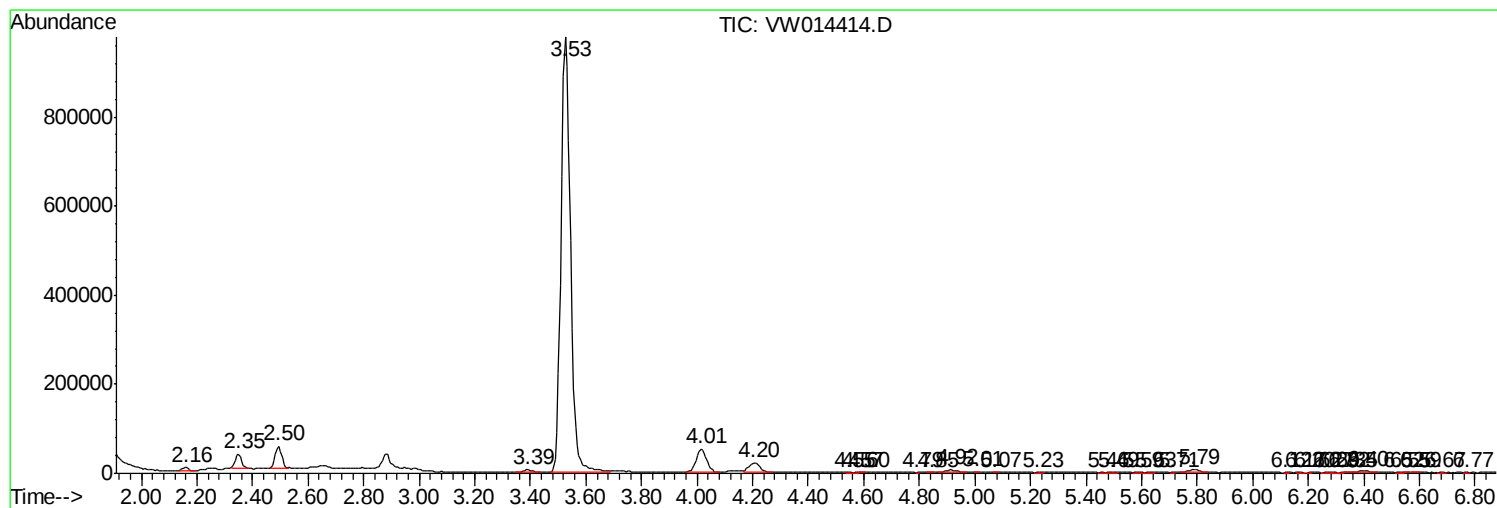
Sum of corrected areas: 4983310

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ALS Vial : 1 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM120519S.M
Quant Title : VOC Analysis

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



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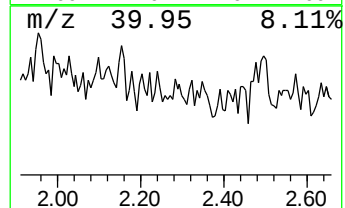
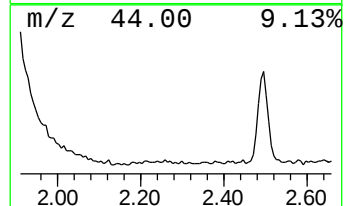
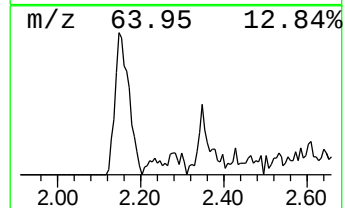
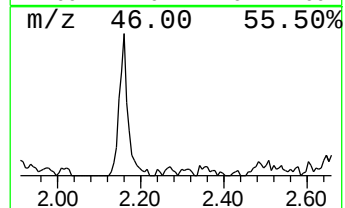
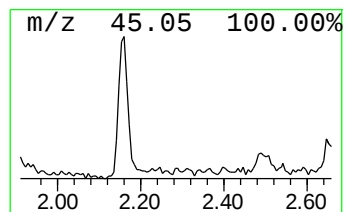
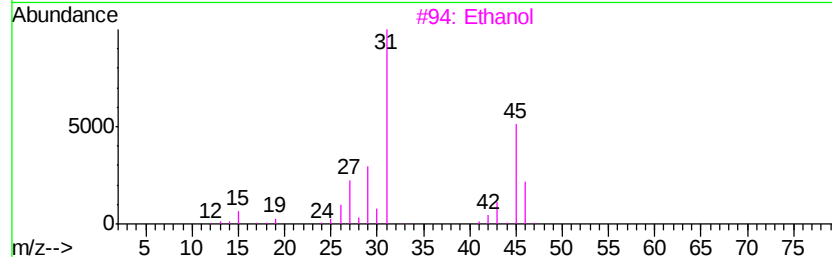
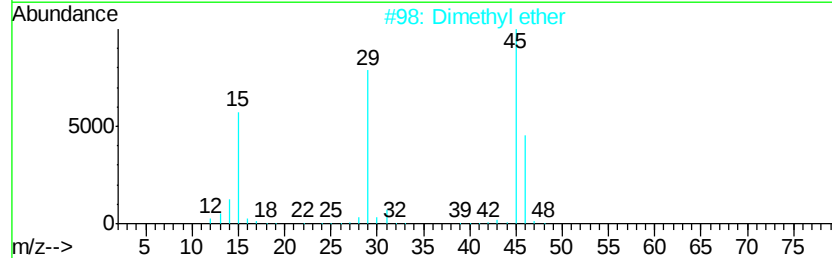
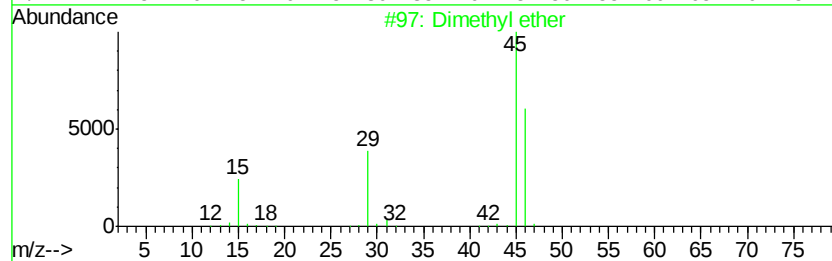
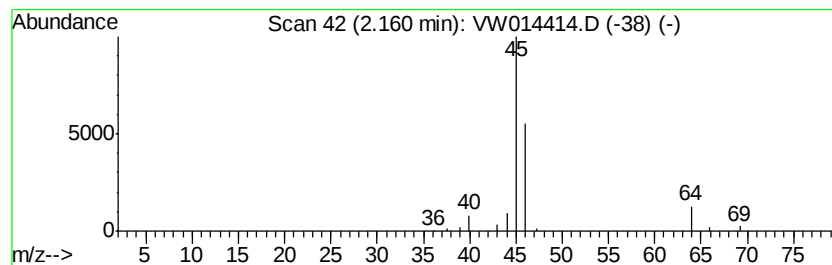
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM120519S.M
Quant Title : VOC Analysis

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

Peak Number 1 unknown-01 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.16	1.32 ug/L	11845	1,4-Difluorobenzene	8.84

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dimethyl ether	46	C2H6O	000115-10-6	9
2			Dimethyl ether	46	C2H6O	000115-10-6	9
3			Ethanol	46	C2H6O	000064-17-5	7
4			Acetone-D6	64	C3D6O	000666-52-4	7
5			Ethanol	46	C2H6O	000064-17-5	7



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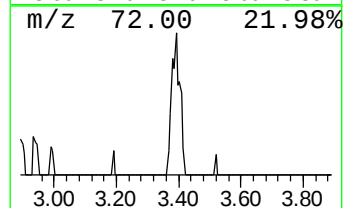
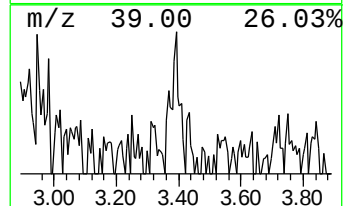
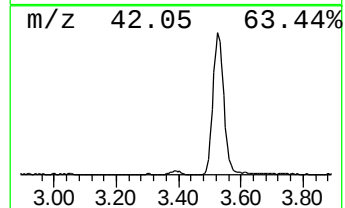
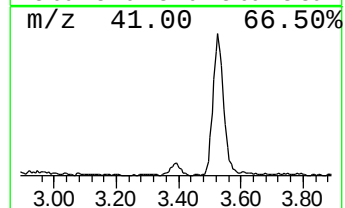
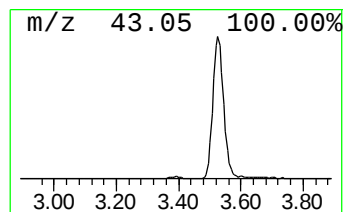
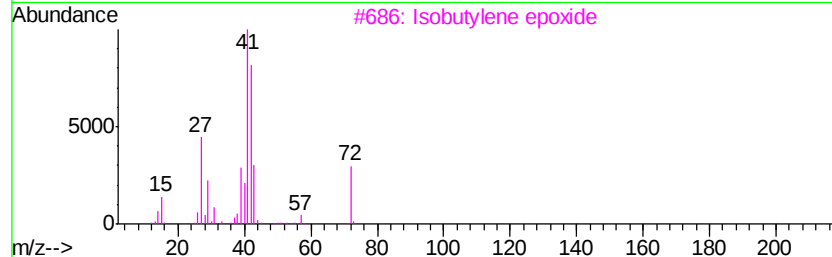
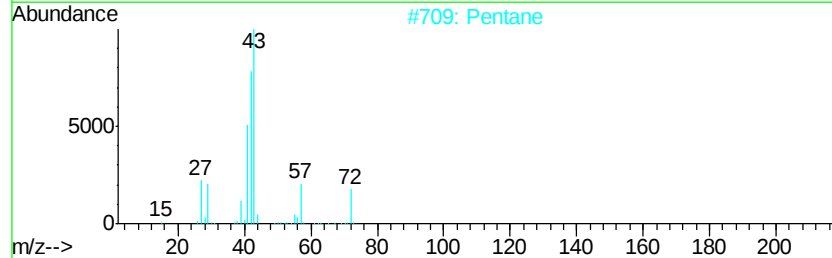
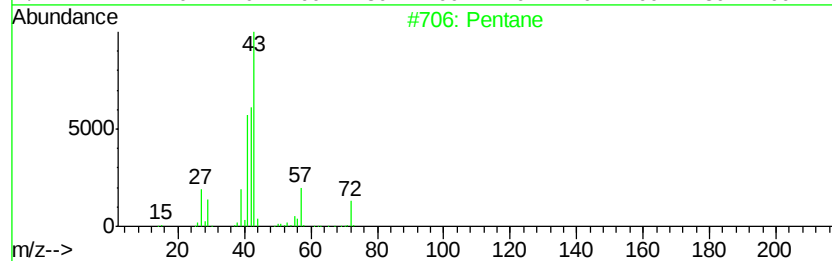
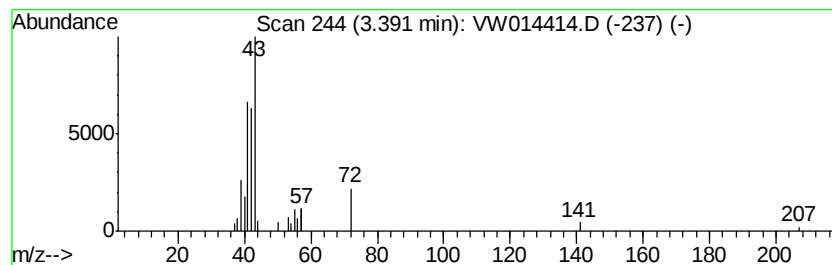
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Quant Title : VOC Analysis

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.39	1.62 ug/L	14585	1,4-Difluorobenzene	8.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane	72	C5H12	000109-66-0	72
2		Pentane	72	C5H12	000109-66-0	64
3		Isobutylene epoxide	72	C4H8O	000558-30-5	49
4		Diaziridine,3,3-dimethyl-	72	C3H8N2	004901-76-2	47
5		Diaziridine,3,3-dimethyl-	72	C3H8N2	004901-76-2	38



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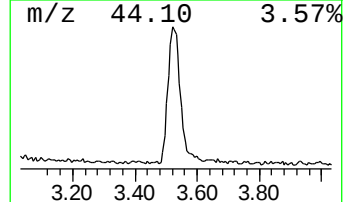
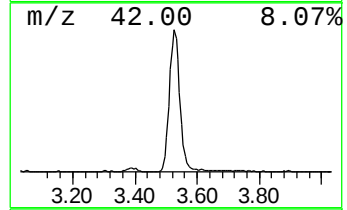
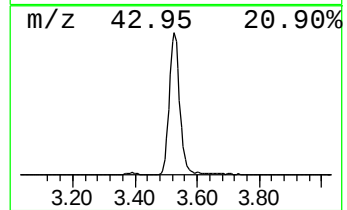
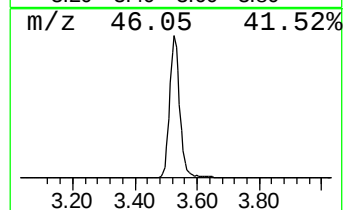
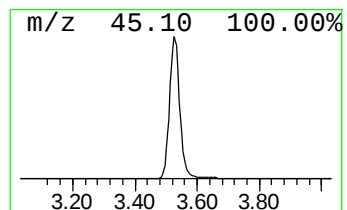
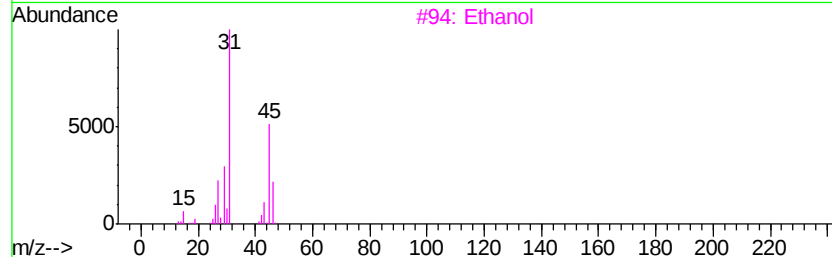
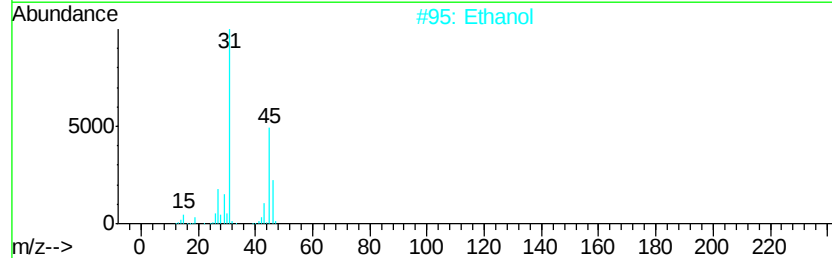
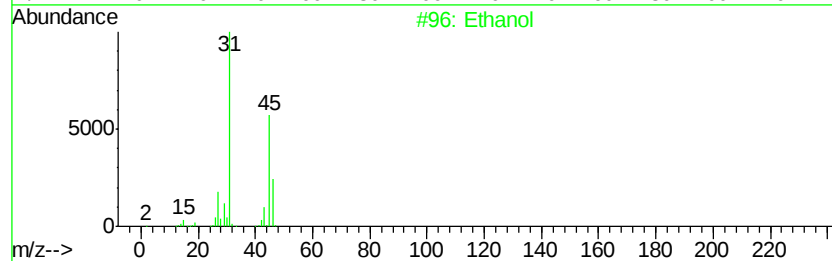
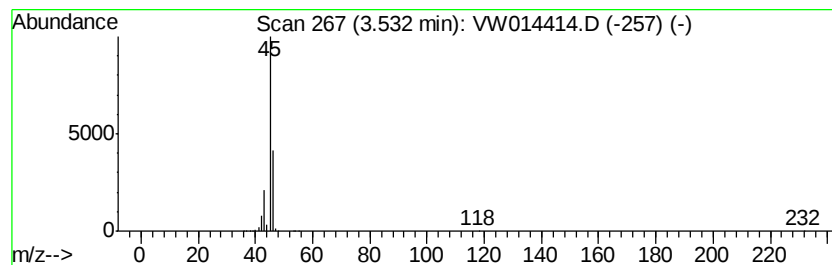
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Quant Title : VOC Analysis

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

Peak Number 4 Ethanol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.53	257.95 ug/L	2320280	1,4-Difluorobenzene	8.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol	46	C2H6O	000064-17-5	64
2		Ethanol	46	C2H6O	000064-17-5	64
3		Ethanol	46	C2H6O	000064-17-5	56
4		Dimethyl ether	46	C2H6O	000115-10-6	9
5		Dimethyl ether	46	C2H6O	000115-10-6	9



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW121719\
Data File : VW014414.D
Acq On : 18 Dec 2019 01:20
Operator : SY/VA
Sample : K6279-10
Misc : 3.99G/10ML/MSVOA W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BFEY4

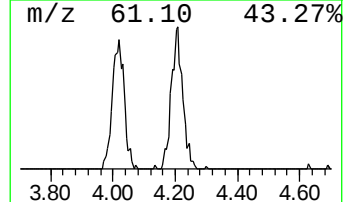
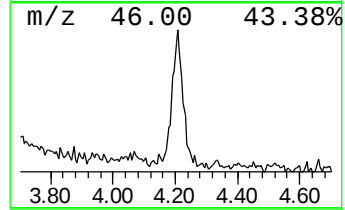
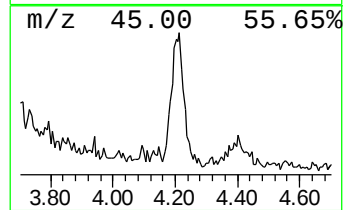
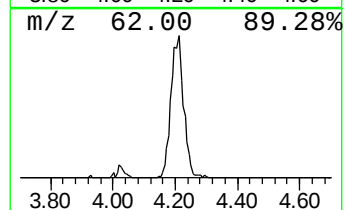
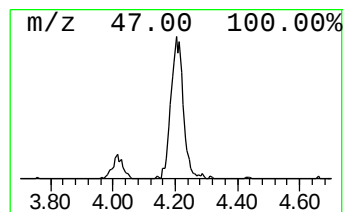
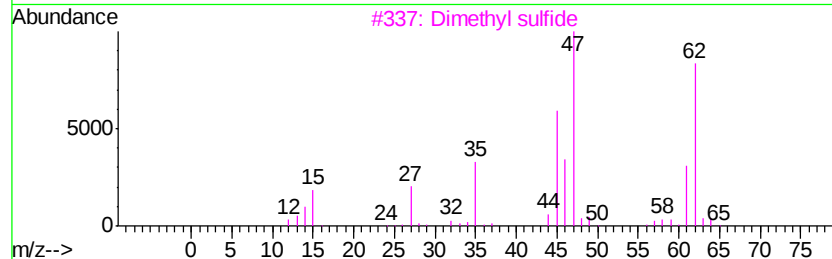
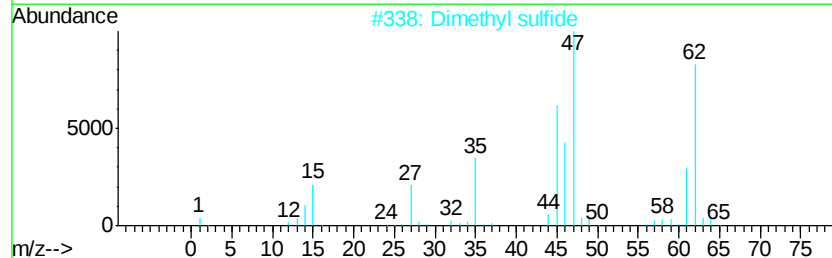
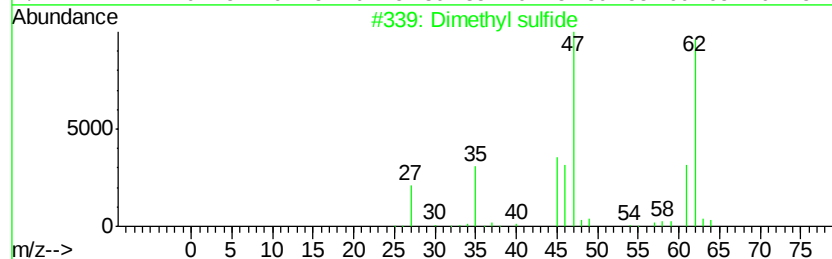
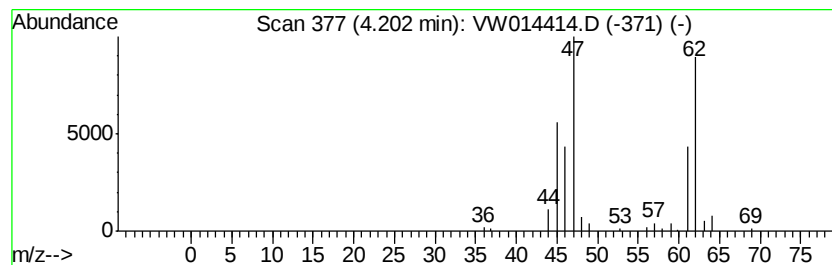
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Quant Title : VOC Analysis

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

Peak Number 5 Dimethyl sulfide Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.20	6.61 ug/L	59426	1,4-Difluorobenzene	8.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dimethyl sulfide	62	C2H6S	000075-18-3	91
2		Dimethyl sulfide	62	C2H6S	000075-18-3	91
3		Dimethyl sulfide	62	C2H6S	000075-18-3	91
4		Dimethyl sulfide	62	C2H6S	000075-18-3	91
5		Borane-methyl sulfide complex	76	C2H9BS	013292-87-0	83



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW121719\
Data File : VW014414.D
Acq On : 18 Dec 2019 01:20
Operator : SY/VA
Sample : K6279-10
Misc : 3.99G/10ML/MSVOA W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BFEY4

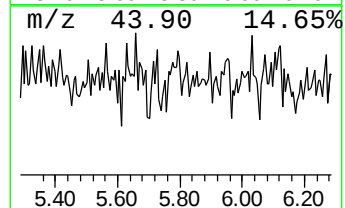
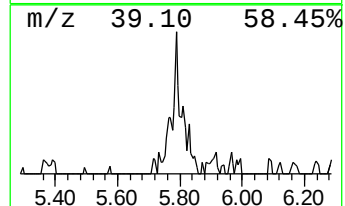
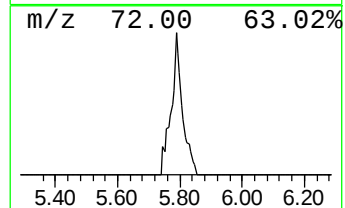
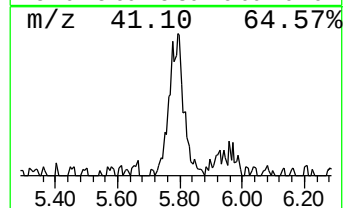
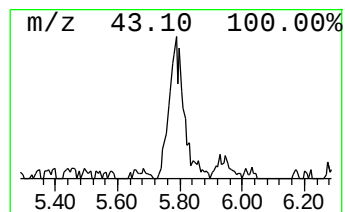
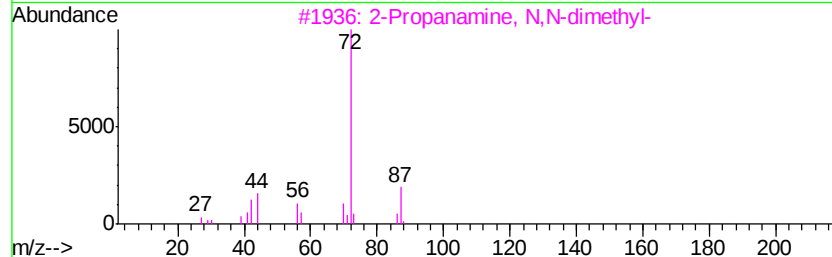
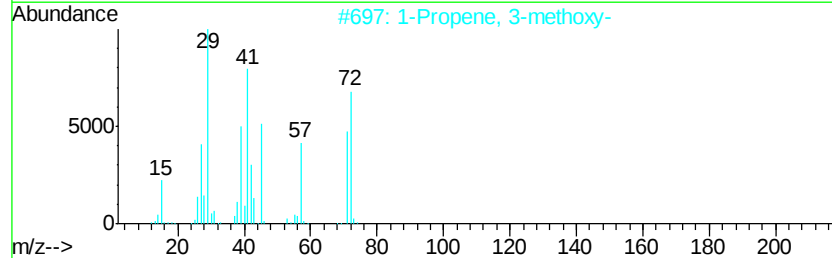
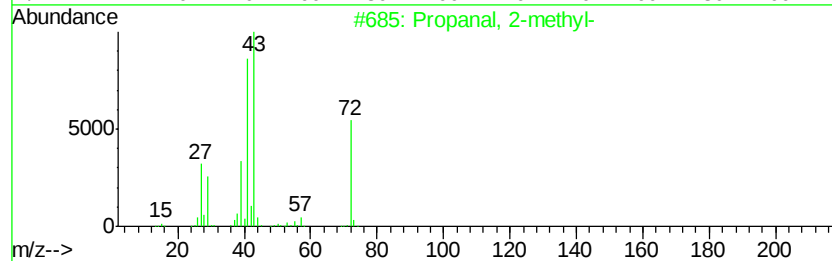
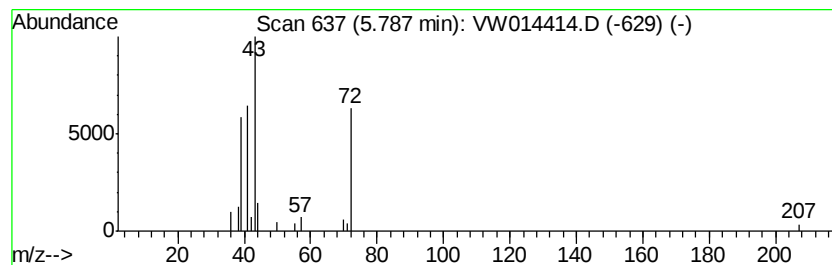
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Quant Title : VOC Analysis

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

Peak Number 6 unknown-02 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.79	2.24 ug/L	20160	1,4-Difluorobenzene	8.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propanal, 2-methyl-	72	C4H8O	000078-84-2	38
2		1-Propene, 3-methoxy-	72	C4H8O	000627-40-7	28
3		2-Propanamine, N,N-dimethyl-	87	C5H13N	000996-35-0	12
4		Propanal, 2-methyl-	72	C4H8O	000078-84-2	9
5		Propanal, 2-methyl-	72	C4H8O	000078-84-2	9



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW121719\
Data File : VW014414.D
Acq On : 18 Dec 2019 01:20
Operator : SY/VA
Sample : K6279-10
Misc : 3.99G/10ML/MSVOA W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BFEY4

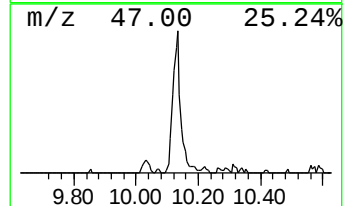
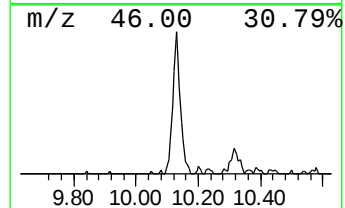
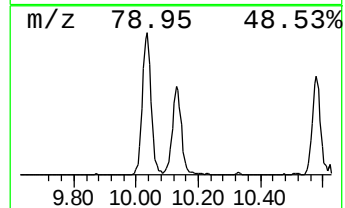
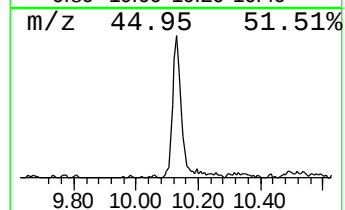
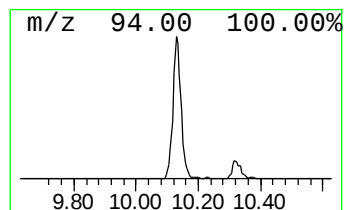
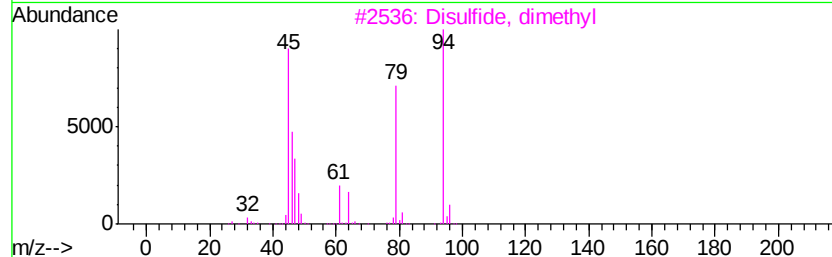
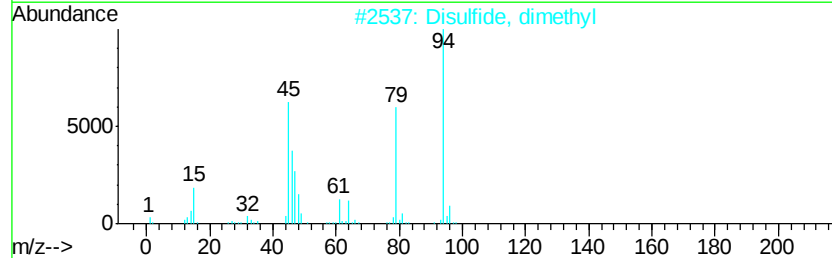
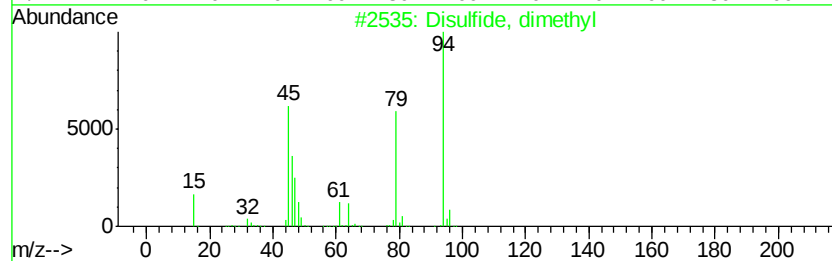
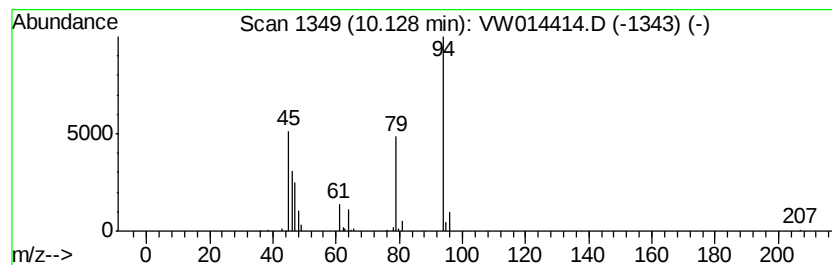
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM120519S.M
Quant Title : VOC Analysis

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

Peak Number 8 Disulfide, dimethyl Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.13	8.13 ug/L	73140	1,4-Difluorobenzene	8.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Disulfide, dimethyl	94	C2H6S2	000624-92-0	97
2		Disulfide, dimethyl	94	C2H6S2	000624-92-0	95
3		Disulfide, dimethyl	94	C2H6S2	000624-92-0	95
4		Disulfide, dimethyl	94	C2H6S2	000624-92-0	94
5		Disulfide, dimethyl	94	C2H6S2	000624-92-0	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW121719\
Data File : VW014414.D
Acq On : 18 Dec 2019 01:20
Operator : SY/VA
Sample : K6279-10
Misc : 3.99G/10ML/MSVOA W/SOIL
ALS Vial : 1 Sample Multiplier: 1

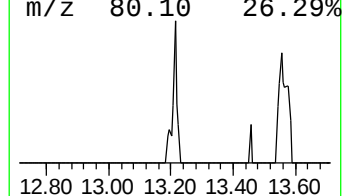
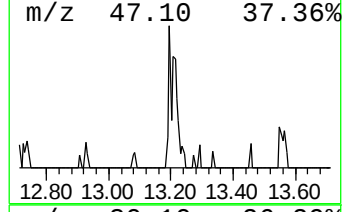
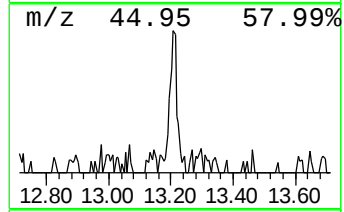
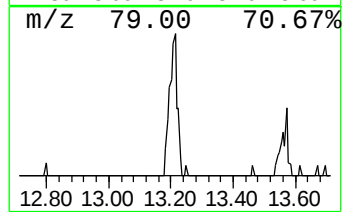
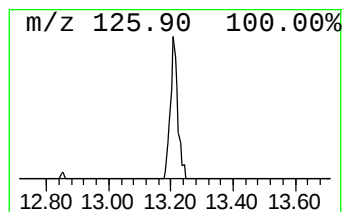
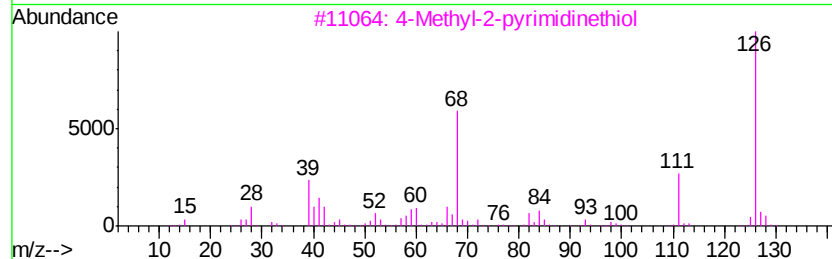
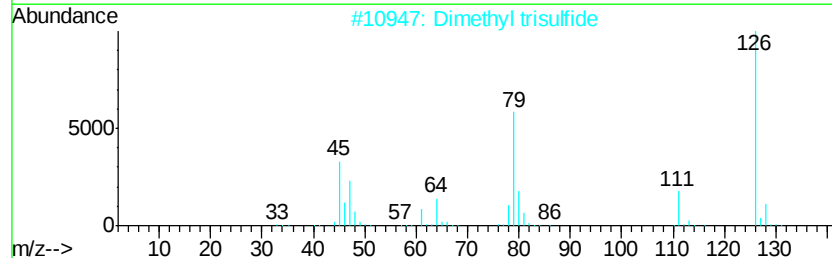
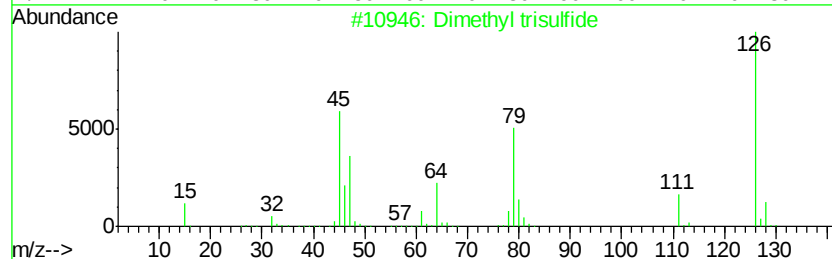
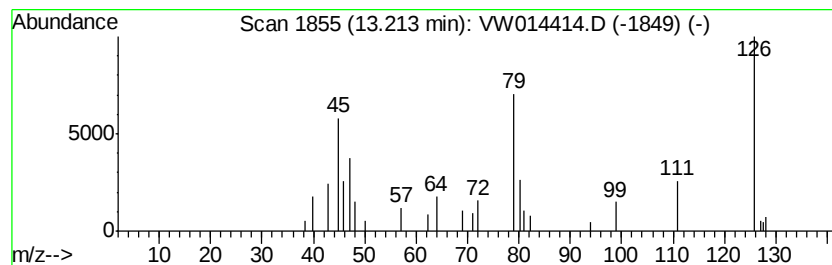
Instrument :
MSVOA_W
ClientSampleId :
BFEY4

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM120519S.M
Quant Title : VOC Analysis

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

Peak Number 13 Dimethyl trisulfide Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.21	3.52 ug/L	11429	1,4-Dichlorobenzene-d4	13.56	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dimethyl trisulfide	126	C2H6S3	003658-80-8	87
2	Dimethyl trisulfide	126	C2H6S3	003658-80-8	56
3	4-Methyl-2-pyrimidinethiol	126	C5H6N2S	035071-17-1	22
4	2-Cyclopenten-1-one, 3-ethyl-2-h...	126	C7H10O2	021835-01-8	17
5	2(1H)-Pyrimidinone, 4,6-diamino-	126	C4H6N4O	031458-45-4	12



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_W\DATA\VW121719\
Data File : VW014414.D
Acq On : 18 Dec 2019 01:20
Operator : SY/VA
Sample : K6279-10
Misc : 3.99G/10ML/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BFEY4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\SOM2WLM120519S.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
unknown-01	2.16	1.3	ug/L	11845	1	8.84	224880	25.0
(DEL) Alkane: Str...	3.39	1.6	ug/L	14585	1	8.84	224880	25.0
Ethanol	3.53	257.9	ug/L	2320280	1	8.84	224880	25.0
Dimethyl sulfide	4.20	6.6	ug/L	59426	1	8.84	224880	25.0
unknown-02	5.79	2.2	ug/L	20160	1	8.84	224880	25.0
Disulfide, dimethyl	10.13	8.1	ug/L	73140	1	8.84	224880	25.0
Dimethyl trisulfide	13.21	3.5	ug/L	11429	3	13.56	81243	25.0