

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU032719\
 Data File : VU030494.D
 Acq On : 27 Mar 2019 19:31
 Operator : JC/SP
 Sample : K2063-14DL 4X
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 DB6S1DL

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_U\METHOD\SOMULM032319WMA.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	1.183	24	29	36	rVB	16141	14832	0.79%	0.078%
2	1.399	89	96	109	rBV	255829	261011	13.85%	1.373%
3	1.675	175	182	197	rVB	191427	255207	13.54%	1.342%
4	1.752	199	206	220	rVB	80987	114180	6.06%	0.601%
5	1.942	260	265	275	rVB	23097	30398	1.61%	0.160%
6	2.048	294	298	307	rVB3	11259	13750	0.73%	0.072%
7	2.270	356	367	380	rBV	389793	603165	32.00%	3.172%
8	2.604	467	471	474	rBV	1797	1467	0.08%	0.008%
9	2.640	479	482	486	rVV2	1833	1522	0.08%	0.008%
10	2.701	492	501	507	rBV3	19565	37641	2.00%	0.198%
11	3.000	578	594	611	rBV3	24576	57109	3.03%	0.300%
12	3.106	624	627	633	rBV	1678	1899	0.10%	0.010%
13	3.264	672	676	679	rBV	1860	1620	0.09%	0.009%
14	3.299	679	687	699	rVV5	8006	17363	0.92%	0.091%
15	3.498	741	749	758	rBV7	2896	6224	0.33%	0.033%
16	3.572	760	772	793	rVB4	14289	41103	2.18%	0.216%
17	3.704	810	813	820	rVB	2543	2874	0.15%	0.015%
18	3.743	820	825	827	rBV	1902	2025	0.11%	0.011%
19	3.913	873	878	881	rBV	1693	1386	0.07%	0.007%
20	3.981	888	899	900	rBV3	4852	6904	0.37%	0.036%
21	4.093	918	934	948	rBV3	71795	193059	10.24%	1.015%
22	4.173	948	959	985	rVV	181957	465082	24.67%	2.446%
23	4.360	1015	1017	1020	rVB	2498	1410	0.07%	0.007%
24	4.379	1020	1023	1028	rBV	1998	1999	0.11%	0.011%
25	4.646	1090	1106	1129	rBV2	306922	800885	42.49%	4.212%
26	4.919	1187	1191	1195	rVB	1960	1387	0.07%	0.007%
27	4.993	1199	1214	1226	rBV3	27210	66744	3.54%	0.351%
28	5.077	1229	1240	1254	rVB5	15336	37621	2.00%	0.198%
29	5.215	1272	1283	1292	rBV5	8363	20533	1.09%	0.108%
30	5.337	1298	1321	1330	rBV2	611657	1805195	95.77%	9.494%
31	5.620	1403	1409	1421	rVB8	9213	17898	0.95%	0.094%
32	5.771	1448	1456	1459	rBV2	3679	4732	0.25%	0.025%
33	5.884	1478	1491	1519	rBV	529461	1065958	56.55%	5.606%
34	6.148	1570	1573	1576	rBV	1671	1366	0.07%	0.007%

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Integration Parameters: LSCINT.P

Integrator: RTE
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Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM032319WMA.M
 Title : VOC Analysis

35	6.328	1613	1629	1644	rBV	412197	840350	44.58%	4.420%
36	6.759	1752	1763	1774	rVB3	8058	16913	0.90%	0.089%
37	6.874	1795	1799	1802	rBV	1791	1625	0.09%	0.009%
38	6.923	1803	1814	1824	rVV4	16968	32561	1.73%	0.171%
39	7.045	1838	1852	1867	rBV3	30346	67377	3.57%	0.354%
40	7.186	1892	1896	1897	rBV	2744	1762	0.09%	0.009%
41	7.225	1897	1908	1924	rVV	219997	397915	21.11%	2.093%
42	7.369	1949	1953	1963	rVB2	8371	12356	0.66%	0.065%
43	7.562	1993	2013	2024	rBV	768026	1341013	71.14%	7.053%
44	7.633	2025	2035	2056	rVB	1075503	1884909	100.00%	9.913%
45	7.849	2092	2102	2119	rBV	150875	261793	13.89%	1.377%
46	8.006	2148	2151	2155	rBV	2042	1830	0.10%	0.010%
47	8.086	2172	2176	2181	rBV	2231	2266	0.12%	0.012%
48	8.138	2187	2192	2195	rBV	1712	1475	0.08%	0.008%
49	8.308	2234	2245	2277	rBV	580410	1085072	57.57%	5.707%
50	8.649	2347	2351	2353	rBV	1782	1429	0.08%	0.008%
51	8.710	2366	2370	2372	rBV2	2196	1438	0.08%	0.008%
52	8.858	2411	2416	2423	rVB2	1829	2603	0.14%	0.014%
53	8.913	2431	2433	2439	rVB	2360	2095	0.11%	0.011%
54	8.948	2439	2444	2449	rBV2	1895	1931	0.10%	0.010%
55	9.029	2465	2469	2474	rBV4	2612	3145	0.17%	0.017%
56	9.086	2475	2487	2509	rBV	782287	1306643	69.32%	6.872%
57	9.247	2527	2537	2559	rBV	218092	357349	18.96%	1.879%
58	9.369	2564	2575	2600	rVV	524394	900860	47.79%	4.738%
59	9.540	2625	2628	2633	rBV2	1397	1309	0.07%	0.007%
60	9.710	2678	2681	2685	rVB	1729	1321	0.07%	0.007%
61	9.778	2685	2702	2716	rBV	265835	445600	23.64%	2.344%
62	9.935	2746	2751	2753	rBV2	1911	2071	0.11%	0.011%
63	10.032	2773	2781	2783	rBV	2246	2918	0.15%	0.015%
64	10.109	2800	2805	2807	rBV	1479	1380	0.07%	0.007%
65	10.164	2812	2822	2835	rVV2	13810	24476	1.30%	0.129%
66	10.427	2890	2904	2923	rBV	553923	901861	47.85%	4.743%
67	10.537	2935	2938	2942	rBV	2103	2193	0.12%	0.012%
68	10.588	2946	2954	2965	rVB2	30004	48784	2.59%	0.257%
69	10.678	2972	2982	3002	rBV	97132	221613	11.76%	1.166%
70	10.768	3002	3010	3023	rVB2	55856	87792	4.66%	0.462%
71	10.910	3051	3054	3059	rVB2	1785	1543	0.08%	0.008%

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72	10.967	3059	3072	3091	rBV2	52791	93078	4.94%	0.490%
73	11.147	3117	3128	3143	rBV	112524	182397	9.68%	0.959%
74	11.250	3158	3160	3163	rBV	1823	1332	0.07%	0.007%
75	11.286	3167	3171	3173	rBV3	1943	1612	0.09%	0.008%
76	11.421	3204	3213	3220	rVV6	13027	19050	1.01%	0.100%
77	11.479	3220	3231	3251	rVV	674021	1125831	59.73%	5.921%
78	11.569	3251	3259	3269	rVB	40360	64400	3.42%	0.339%
79	11.765	3311	3320	3329	rBV5	19252	37847	2.01%	0.199%
80	11.855	3337	3348	3371	rVB	666492	1145032	60.75%	6.022%
81	11.977	3383	3386	3389	rBV4	2278	1732	0.09%	0.009%
82	12.054	3406	3410	3420	rVB5	6592	9232	0.49%	0.049%
83	12.144	3434	3438	3443	rBV4	3797	4530	0.24%	0.024%
84	12.254	3463	3472	3478	rBV3	11621	17083	0.91%	0.090%
85	12.479	3536	3542	3544	rBV2	2043	1839	0.10%	0.010%
86	12.649	3588	3595	3600	rVV5	3534	5512	0.29%	0.029%
87	12.704	3605	3612	3619	rVB3	6725	11825	0.63%	0.062%
88	12.893	3666	3671	3675	rBV	1823	1874	0.10%	0.010%
89	13.173	3753	3758	3762	rBV	1282	1758	0.09%	0.009%
90	13.508	3852	3862	3876	rBV4	19135	37647	2.00%	0.198%
91	13.749	3932	3937	3952	rBV3	4165	10100	0.54%	0.053%
92	13.993	4010	4013	4020	rVB3	2472	2152	0.11%	0.011%
93	14.228	4083	4086	4088	rBV	1951	1346	0.07%	0.007%
94	14.315	4108	4113	4115	rBV	1259	1402	0.07%	0.007%
95	15.430	4457	4460	4462	rBV2	2007	1459	0.08%	0.008%
96	15.466	4468	4471	4474	rBV	1625	1275	0.07%	0.007%
97	15.511	4478	4485	4486	rBV	2368	2468	0.13%	0.013%
98	15.832	4582	4585	4588	rBV	2350	1332	0.07%	0.007%
99	15.896	4600	4605	4607	rBV2	2231	1799	0.10%	0.009%
100	16.028	4643	4646	4649	rBV2	2588	1885	0.10%	0.010%

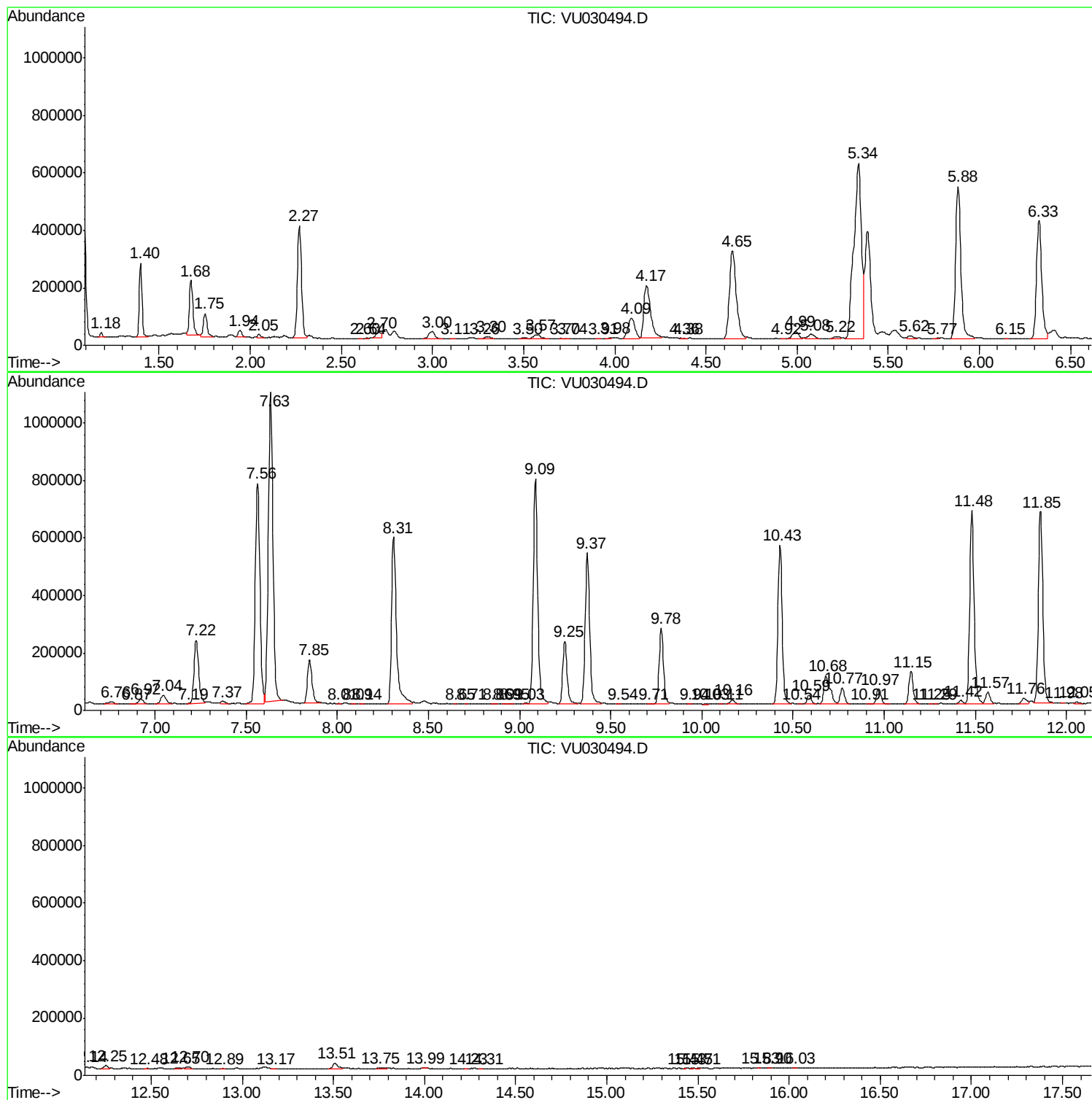
Sum of corrected areas: 19013949

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Instrument :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM032319WMA.M
Quant Title : VOC Analysis

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



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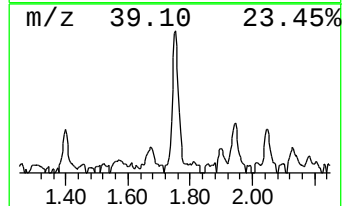
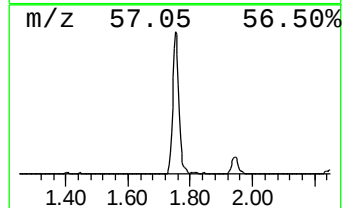
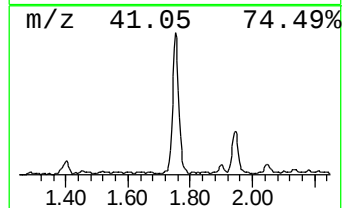
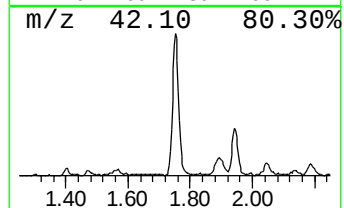
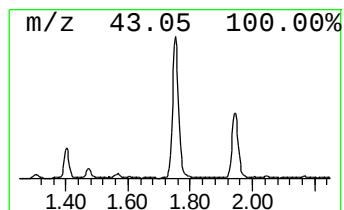
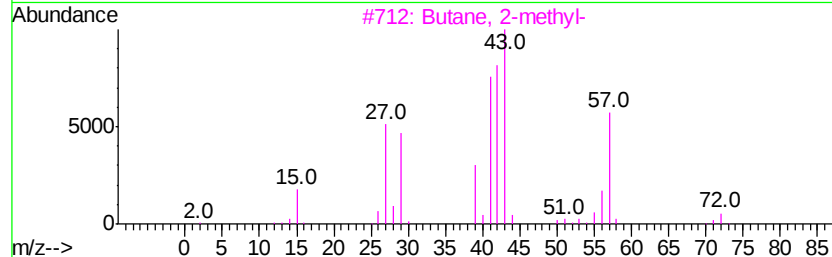
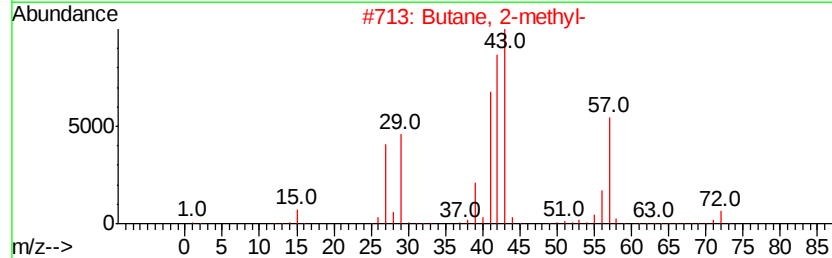
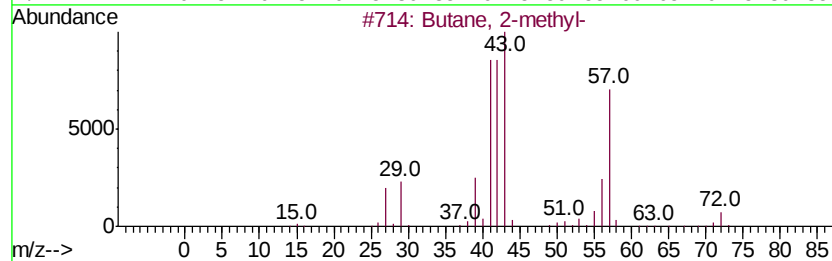
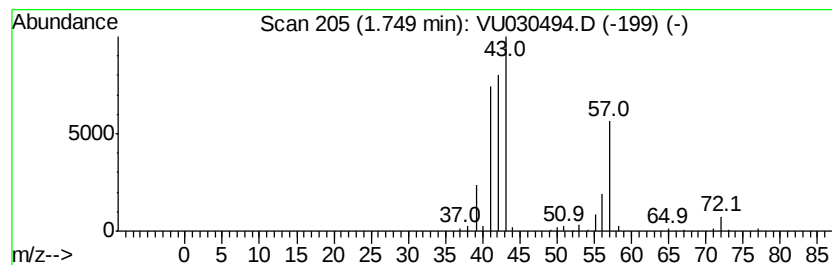
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM032319WMA.M
Quant Title : VOC Analysis

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.75	5.36 ug/L	114180	1,4-Difluorobenzene	5.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methyl-	72	C5H12	000078-78-4	94
2		Butane, 2-methyl-	72	C5H12	000078-78-4	91
3		Butane, 2-methyl-	72	C5H12	000078-78-4	90
4		Pentane	72	C5H12	000109-66-0	45
5		Isobutylene epoxide	72	C4H8O	000558-30-5	40



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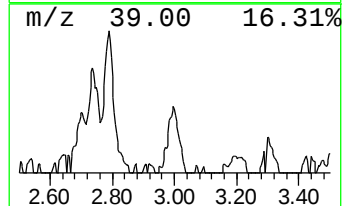
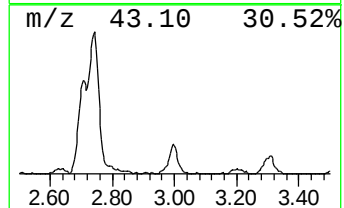
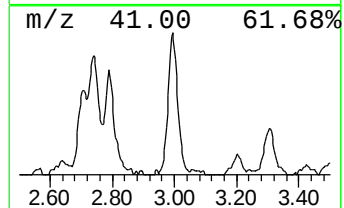
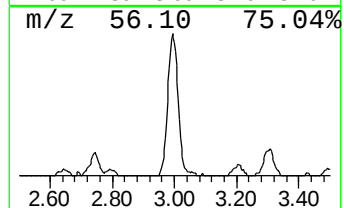
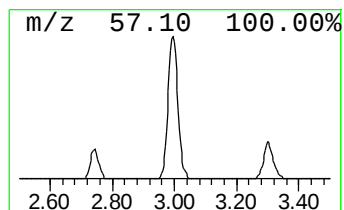
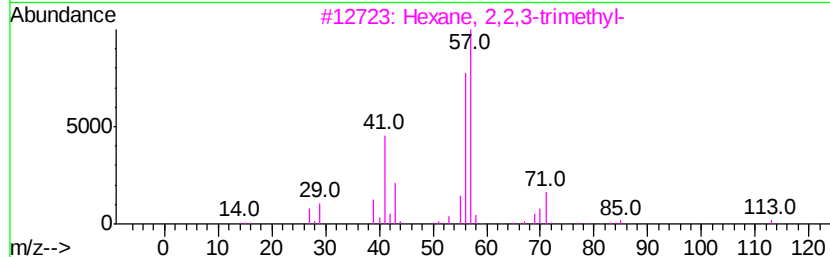
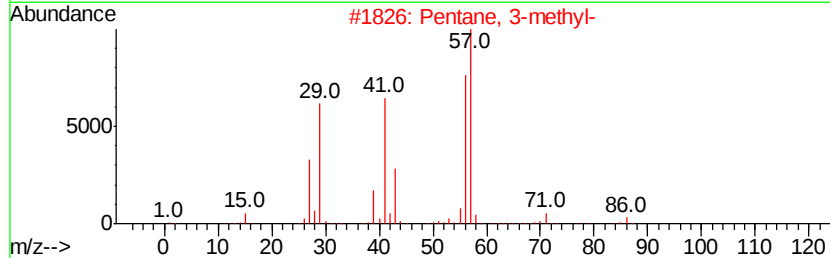
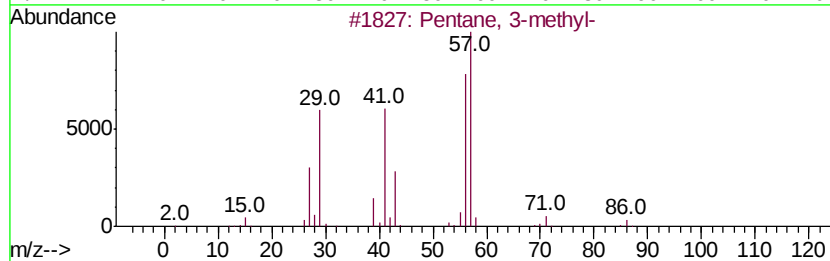
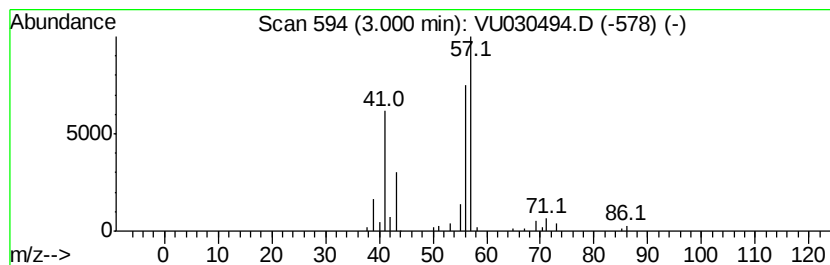
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Quant Title : VOC Analysis

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.00	2.68 ug/L	57109	1,4-Difluorobenzene	5.88

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 3-methyl-	86	C6H14	000096-14-0	86
2			Pentane, 3-methyl-	86	C6H14	000096-14-0	86
3			Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	78
4			Pentane, 3-methyl-	86	C6H14	000096-14-0	78
5			Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	64



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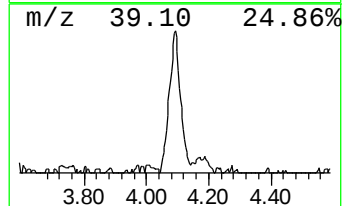
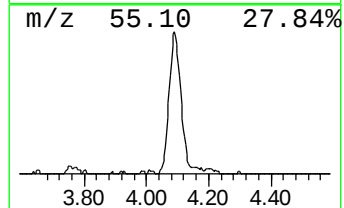
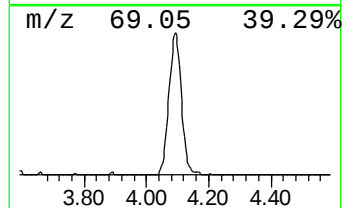
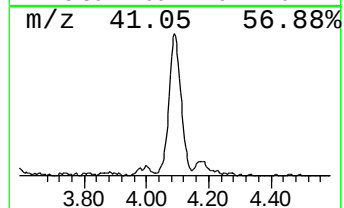
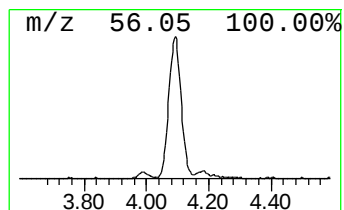
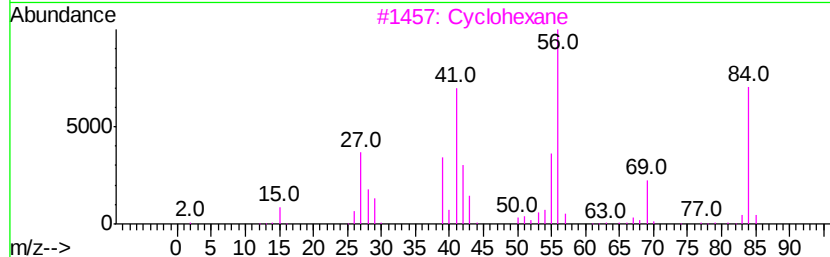
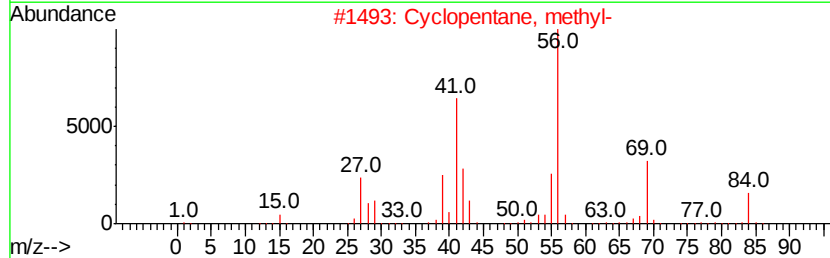
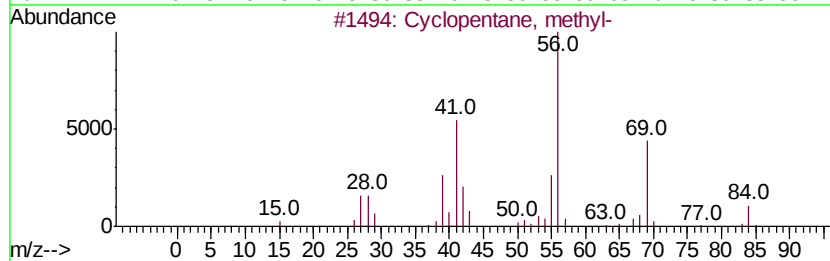
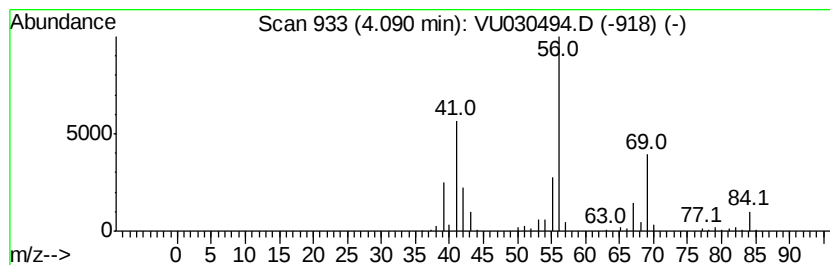
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Quant Title : VOC Analysis

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

Peak Number 3 (DEL) Alkane: Cyclic4.09 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.09	9.06 ug/L	193059	1,4-Difluorobenzene	5.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2		Cyclopentane, methyl-	84	C6H12	000096-37-7	91
3		Cyclohexane	84	C6H12	000110-82-7	87
4		Cyclohexane	84	C6H12	000110-82-7	86
5		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	80



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU032719\
Data File : VU030494.D
Acq On : 27 Mar 2019 19:31
Operator : JC/SP
Sample : K2063-14DL 4X
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
DB6S1DL

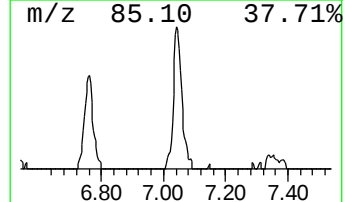
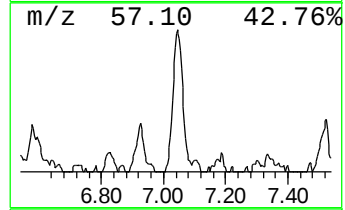
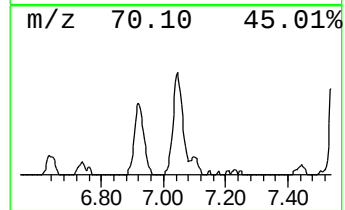
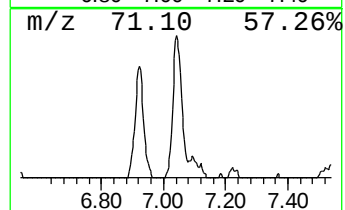
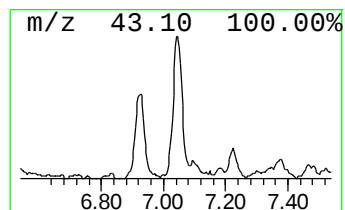
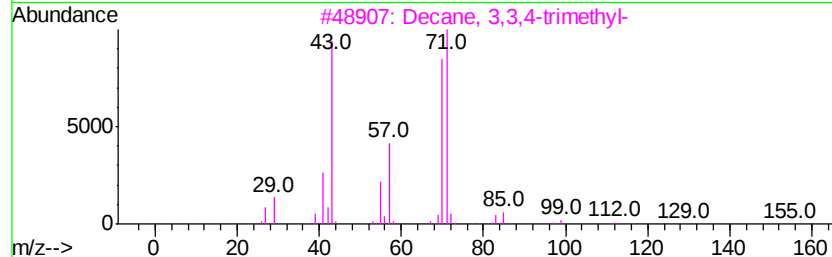
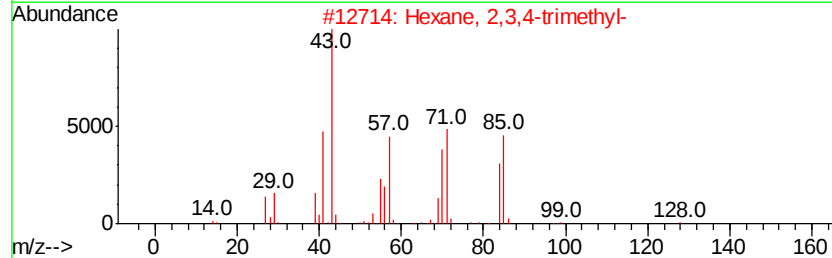
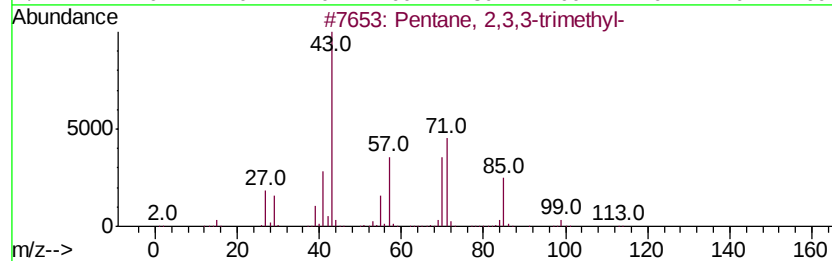
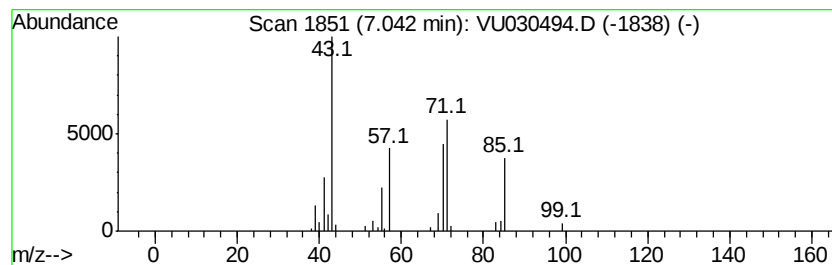
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Quant Title : VOC Analysis

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.04	3.16 ug/L	67377	1,4-Difluorobenzene	5.88

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	83
2			Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	72
3			Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	59
4			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	53
5			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU032719\
Data File : VU030494.D
Acq On : 27 Mar 2019 19:31
Operator : JC/SP
Sample : K2063-14DL 4X
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 24 Sample Multiplier: 1

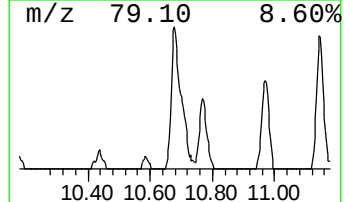
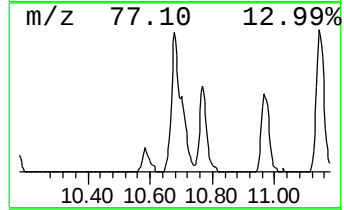
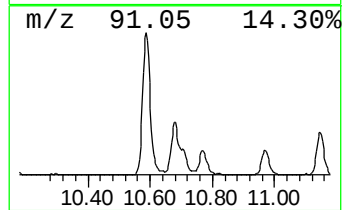
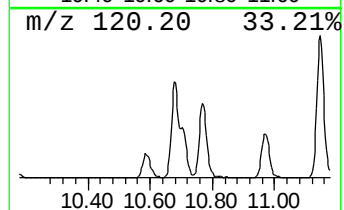
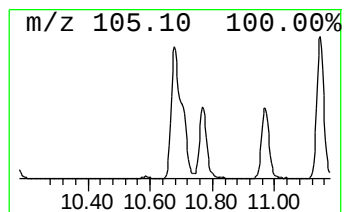
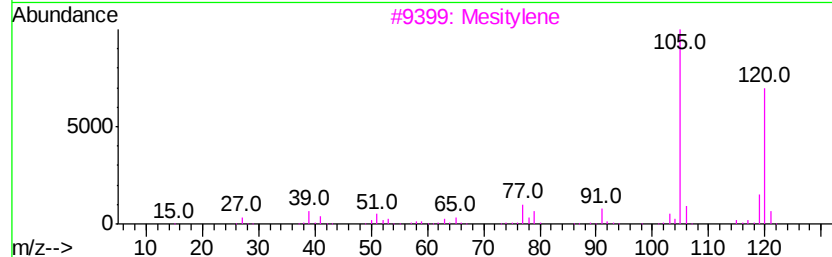
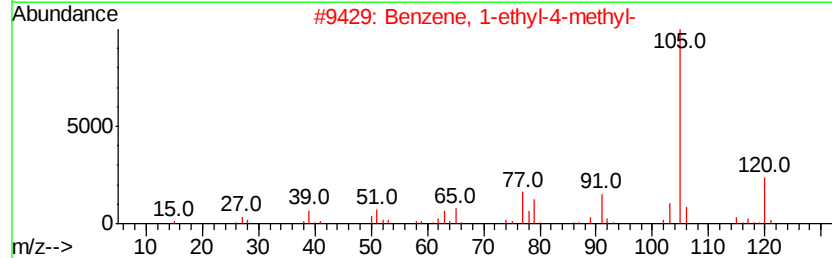
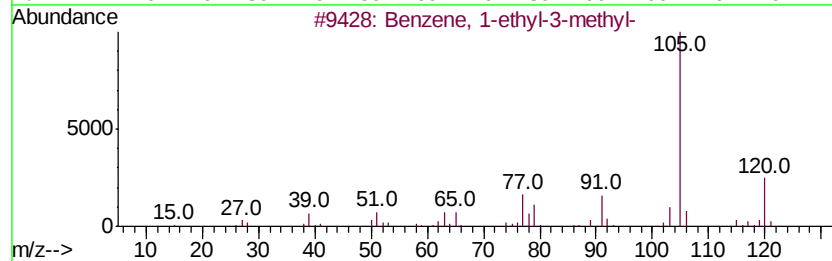
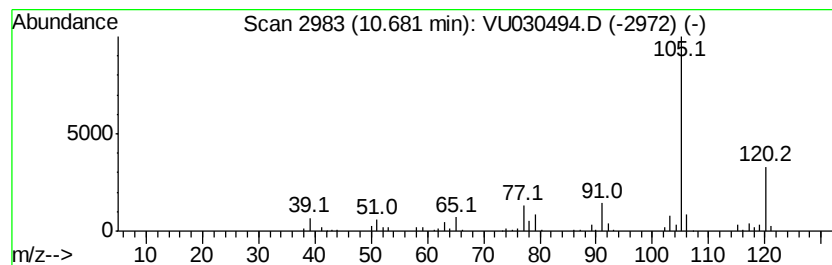
Instrument :
MSVOA_U
ClientSampleId :
DB6S1DL

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM032319WMA.M
Quant Title : VOC Analysis

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

Peak Number 6 Benzene, 1-ethyl-3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.68	9.84 ug/L	221613	1,4-Dichlorobenzene-d4	11.48	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	95
3	Mesitylene	120	C9H12	000108-67-8	91
4	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU032719\
Data File : VU030494.D
Acq On : 27 Mar 2019 19:31
Operator : JC/SP
Sample : K2063-14DL 4X
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
DB6S1DL

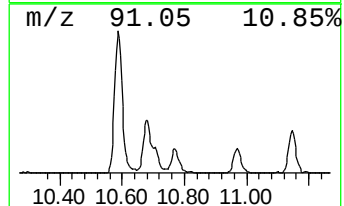
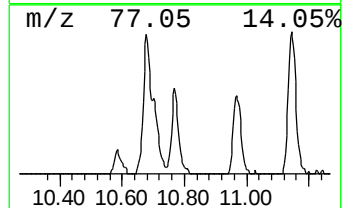
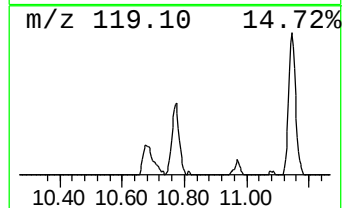
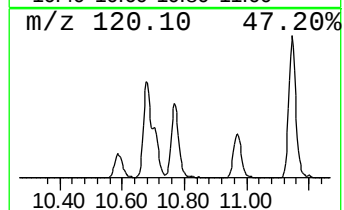
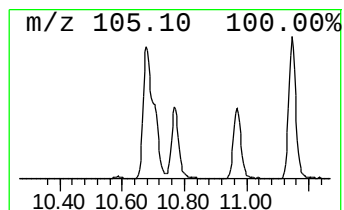
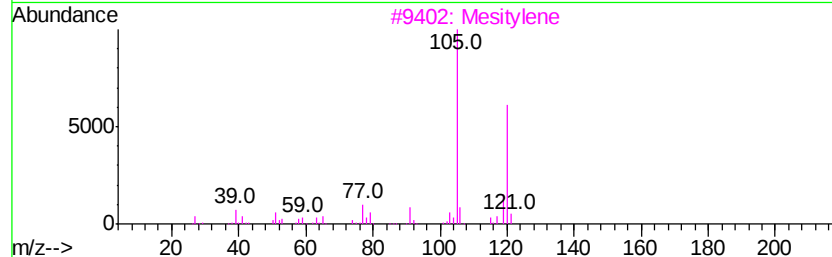
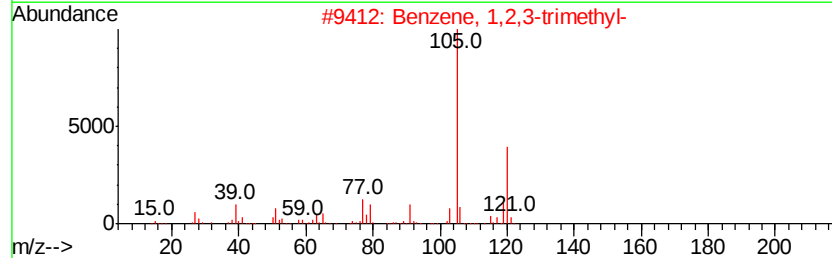
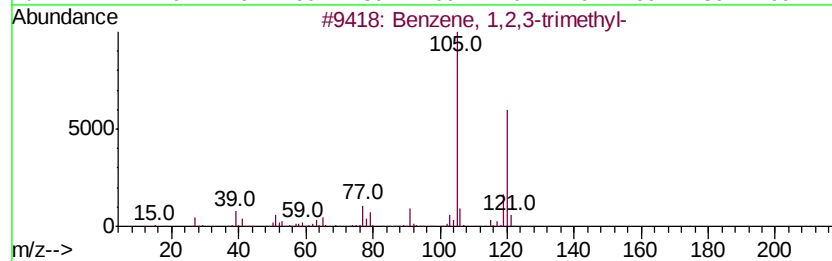
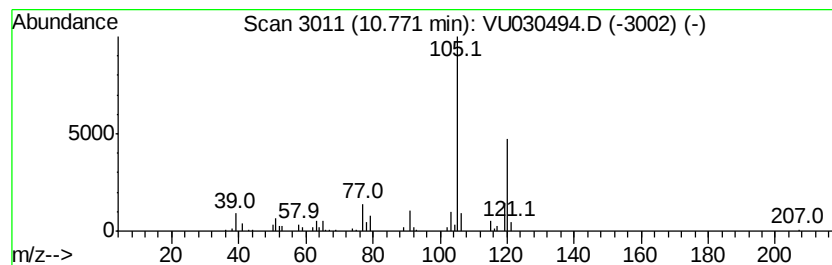
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Quant Title : VOC Analysis

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

Peak Number 7 Benzene, 1,2,3-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.77	3.90 ug/L	87792	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3		Mesitylene	120	C9H12	000108-67-8	95
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU032719\
Data File : VU030494.D
Acq On : 27 Mar 2019 19:31
Operator : JC/SP
Sample : K2063-14DL 4X
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 24 Sample Multiplier: 1

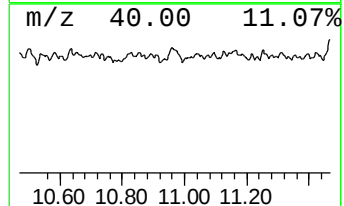
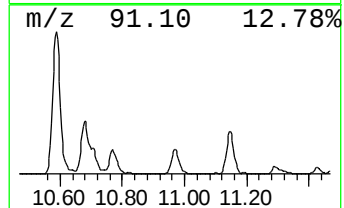
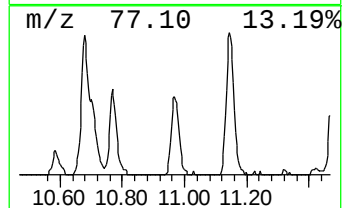
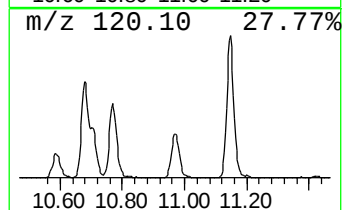
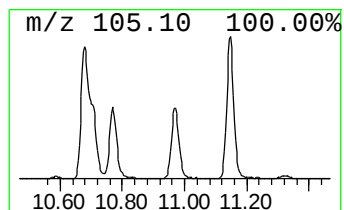
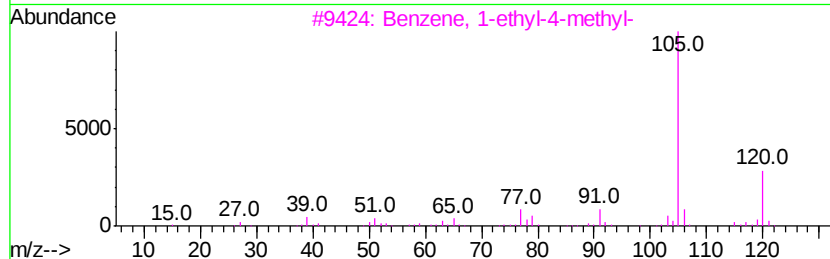
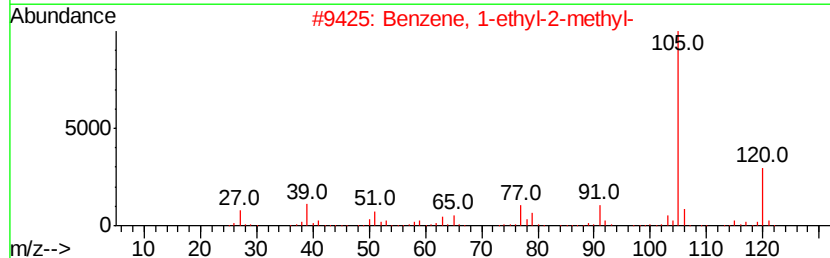
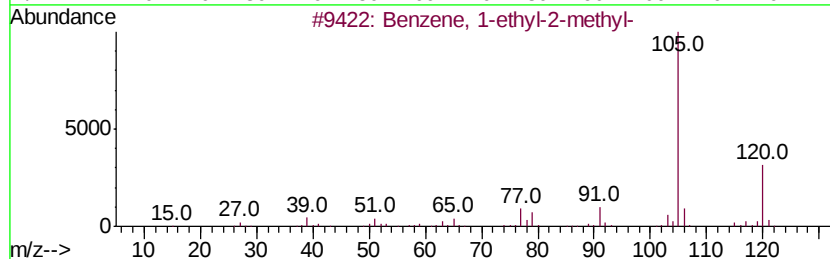
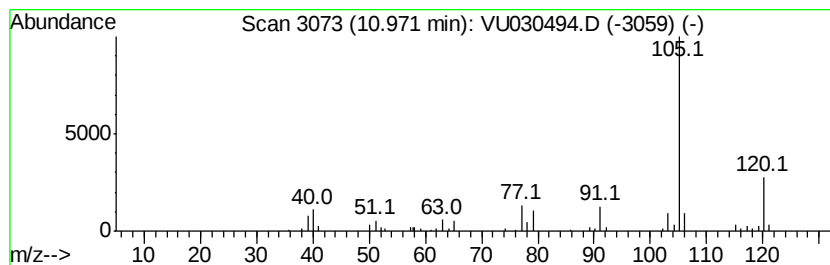
Instrument :
MSVOA_U
ClientSampleId :
DB6S1DL

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM032319WMA.M
Quant Title : VOC Analysis

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

Peak Number 8 Benzene, 1-ethyl-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.97	4.13 ug/L	93078	1,4-Dichlorobenzene-d4	11.48	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5	Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU032719\
Data File : VU030494.D
Acq On : 27 Mar 2019 19:31
Operator : JC/SP
Sample : K2063-14DL 4X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
DB6S1DL

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM032319WMA.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: Str...	1.75	5.4	ug/L	114180	1	5.88	1065960	50.0
(DEL) Alkane: Str...	3.00	2.7	ug/L	57109	1	5.88	1065960	50.0
(DEL) Alkane: Cyc...	4.09	9.1	ug/L	193059	1	5.88	1065960	50.0
(DEL) Alkane: Str...	7.04	3.2	ug/L	67377	1	5.88	1065960	50.0
Benzene, 1-ethyl-...	10.68	9.8	ug/L	221613	3	11.48	1125830	50.0
Benzene, 1,2,3-tr...	10.77	3.9	ug/L	87792	3	11.48	1125830	50.0
Benzene, 1-ethyl-...	10.97	4.1	ug/L	93078	3	11.48	1125830	50.0