

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU032619\
 Data File : VU030459.D
 Acq On : 26 Mar 2019 21:27
 Operator : JC/SP
 Sample : K2063-04DL 100X
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 DB6R6DL

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	35	rVB	16158	15550	0.37%	0.062%
2	1.399	88	96	106	rBV	294469	297094	7.03%	1.194%
3	1.672	173	181	195	rVV	206844	266443	6.30%	1.071%
4	1.749	198	205	224	rVB	62213	90052	2.13%	0.362%
5	1.891	242	249	256	rBV6	7687	14005	0.33%	0.056%
6	1.942	258	265	279	rVB	40469	60544	1.43%	0.243%
7	2.045	291	297	307	rVB	18075	24718	0.58%	0.099%
8	2.132	319	324	332	rBV3	10254	11653	0.28%	0.047%
9	2.183	332	340	354	rVB5	20040	35380	0.84%	0.142%
10	2.267	354	366	380	rBV	420441	650186	15.38%	2.613%
11	2.440	414	420	424	rBV	1889	2527	0.06%	0.010%
12	2.508	438	441	450	rVB2	1976	2466	0.06%	0.010%
13	2.678	485	494	505	rBV4	16034	37973	0.90%	0.153%
14	2.903	558	564	568	rBV	1765	2149	0.05%	0.009%
15	2.939	572	575	578	rVV2	1962	1538	0.04%	0.006%
16	2.994	580	592	607	rVB5	18739	39778	0.94%	0.160%
17	3.193	651	654	663	rVV3	4054	5322	0.13%	0.021%
18	3.260	671	675	676	rBV	2334	1657	0.04%	0.007%
19	3.299	679	687	698	rVV4	9729	19505	0.46%	0.078%
20	3.373	706	710	715	rVB	1589	1575	0.04%	0.006%
21	3.421	722	725	734	rVB5	3013	3811	0.09%	0.015%
22	3.874	860	866	867	rBV2	2736	2650	0.06%	0.011%
23	4.016	907	910	913	rBV	1850	1541	0.04%	0.006%
24	4.090	918	933	947	rBV3	35633	94638	2.24%	0.380%
25	4.174	948	959	987	rVV	202754	518346	12.26%	2.083%
26	4.347	1010	1013	1016	rVB	2442	1606	0.04%	0.006%
27	4.383	1021	1024	1031	rVB	2895	2990	0.07%	0.012%
28	4.418	1031	1035	1040	rBV	1740	1921	0.05%	0.008%
29	4.569	1077	1082	1088	rBV	1410	1803	0.04%	0.007%
30	4.643	1089	1105	1124	rBV	300483	709905	16.79%	2.852%
31	4.881	1176	1179	1183	rVB	1942	1638	0.04%	0.007%
32	4.990	1199	1213	1228	rVV5	30050	73879	1.75%	0.297%
33	5.215	1273	1283	1286	rBV4	4905	6371	0.15%	0.026%
34	5.337	1298	1321	1329	rBV2	630629	1849363	43.75%	7.431%

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

35	5.710	1430	1437	1441	rBV	2172	2533	0.06%	0.010%
36	5.881	1477	1490	1527	rBV	546567	1121878	26.54%	4.508%
37	6.077	1549	1551	1558	rVB	2367	2567	0.06%	0.010%
38	6.112	1558	1562	1566	rBV2	2575	2471	0.06%	0.010%
39	6.209	1587	1592	1600	rBV3	2451	3451	0.08%	0.014%
40	6.328	1614	1629	1646	rBV	423711	864815	20.46%	3.475%
41	6.624	1715	1721	1723	rBV2	2976	3071	0.07%	0.012%
42	6.826	1775	1784	1786	rBV2	2569	3193	0.08%	0.013%
43	6.910	1803	1810	1813	rBV	1715	2257	0.05%	0.009%
44	7.061	1851	1857	1858	rBV	3705	3416	0.08%	0.014%
45	7.132	1876	1879	1882	rBV	1976	1723	0.04%	0.007%
46	7.225	1895	1908	1927	rBV	231824	422522	10.00%	1.698%
47	7.562	2000	2013	2024	rBV	794959	1382073	32.69%	5.553%
48	7.633	2024	2035	2072	rVB	2424483	4227293	100.00%	16.986%
49	7.845	2092	2101	2119	rBV	159359	267150	6.32%	1.073%
50	8.038	2157	2161	2163	rBV2	2035	1694	0.04%	0.007%
51	8.090	2171	2177	2183	rVB2	3338	4342	0.10%	0.017%
52	8.119	2183	2186	2195	rBV	1529	2254	0.05%	0.009%
53	8.308	2234	2245	2287	rBV	663295	1203338	28.47%	4.835%
54	8.582	2322	2330	2333	rBV3	7528	9159	0.22%	0.037%
55	8.643	2345	2349	2354	rBV2	1978	2300	0.05%	0.009%
56	8.935	2437	2440	2446	rVB	2031	2160	0.05%	0.009%
57	9.035	2460	2471	2474	rBV3	2883	4854	0.11%	0.020%
58	9.087	2475	2487	2509	rBV	808833	1369777	32.40%	5.504%
59	9.247	2526	2537	2562	rBV	365860	603226	14.27%	2.424%
60	9.369	2562	2575	2600	rBV	1262548	2089373	49.43%	8.395%
61	9.672	2664	2669	2677	rBV	2190	3321	0.08%	0.013%
62	9.775	2689	2701	2720	rBV	588851	978199	23.14%	3.931%
63	10.167	2814	2823	2833	rBV2	18643	30188	0.71%	0.121%
64	10.273	2852	2856	2859	rBV	1806	1584	0.04%	0.006%
65	10.427	2892	2904	2921	rBV	577107	929680	21.99%	3.736%
66	10.511	2928	2930	2937	rVB	2042	1766	0.04%	0.007%
67	10.585	2944	2953	2963	rBV2	40182	65534	1.55%	0.263%
68	10.678	2971	2982	3001	rBV	200010	433152	10.25%	1.740%
69	10.768	3001	3010	3024	rVB	93643	152114	3.60%	0.611%
70	10.971	3062	3073	3097	rVB	85503	143035	3.38%	0.575%
71	11.086	3105	3109	3111	rBV	1986	1644	0.04%	0.007%

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Integration Parameters: LSCINT.P

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72	11.144	3116	3127	3144	rBV	346225	548787	12.98%	2.205%
73	11.318	3178	3181	3187	rVB2	2808	2319	0.05%	0.009%
74	11.347	3187	3190	3195	rVB	2694	2548	0.06%	0.010%
75	11.424	3206	3214	3220	rBV5	5949	8729	0.21%	0.035%
76	11.482	3220	3232	3250	rBV	743905	1226713	29.02%	4.929%
77	11.569	3251	3259	3275	rVB2	97857	161578	3.82%	0.649%
78	11.765	3309	3320	3329	rBV	86441	153551	3.63%	0.617%
79	11.858	3338	3349	3368	rVB	724756	1238402	29.30%	4.976%
80	11.977	3380	3386	3398	rVB5	13019	21635	0.51%	0.087%
81	12.054	3406	3410	3419	rVB3	5764	7492	0.18%	0.030%
82	12.144	3431	3438	3443	rBV2	13144	19782	0.47%	0.079%
83	12.250	3463	3471	3478	rBV2	19960	31017	0.73%	0.125%
84	12.353	3496	3503	3513	rVB4	15244	22862	0.54%	0.092%
85	12.553	3558	3565	3577	rVB4	6053	11579	0.27%	0.047%
86	12.649	3588	3595	3601	rVV	11621	15406	0.36%	0.062%
87	12.701	3605	3611	3622	rVB2	16524	26038	0.62%	0.105%
88	12.868	3657	3663	3666	rBV2	1583	1508	0.04%	0.006%
89	12.906	3671	3675	3678	rBV	1680	1642	0.04%	0.007%
90	12.967	3685	3694	3704	rBV3	15011	24840	0.59%	0.100%
91	13.122	3733	3742	3756	rVB5	31454	51900	1.23%	0.209%
92	13.360	3813	3816	3820	rVB	2021	1568	0.04%	0.006%
93	13.405	3825	3830	3833	rBV	1711	1914	0.05%	0.008%
94	13.511	3857	3863	3867	rBV5	4073	5272	0.12%	0.021%
95	13.617	3889	3896	3898	rBV4	2357	2302	0.05%	0.009%
96	13.745	3927	3936	3959	rBV	41850	96038	2.27%	0.386%
97	14.045	4026	4029	4033	rBV	1857	1820	0.04%	0.007%
98	15.417	4452	4456	4460	rBV2	2199	2251	0.05%	0.009%
99	15.897	4602	4605	4608	rBV2	2189	1770	0.04%	0.007%
100	16.035	4645	4648	4653	rVB2	2859	2170	0.05%	0.009%

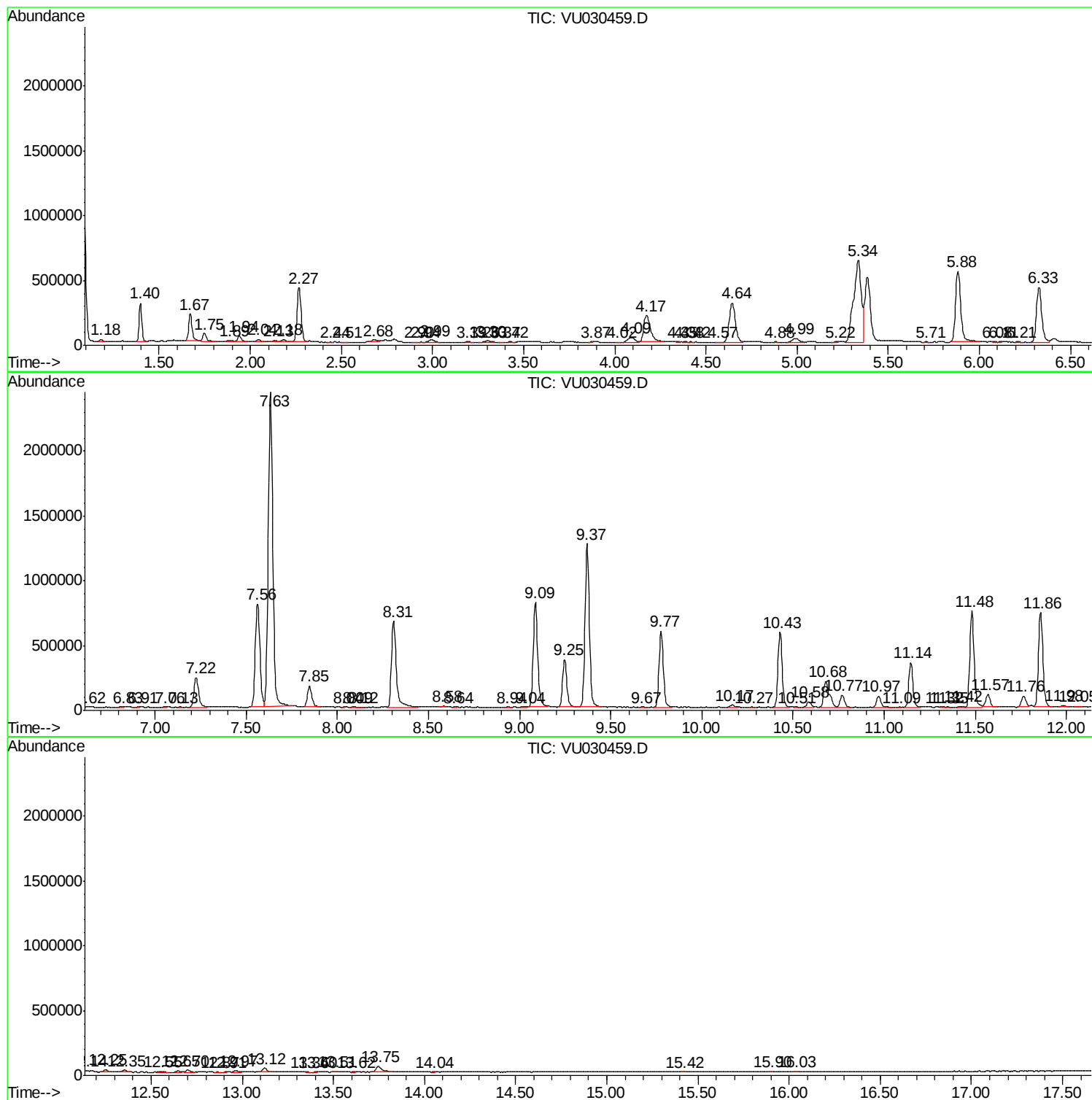
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ALS Vial : 33 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
DB6R6DL

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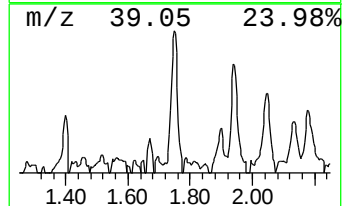
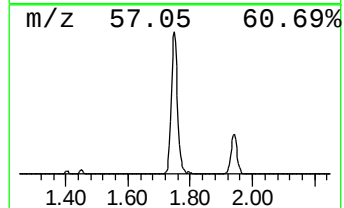
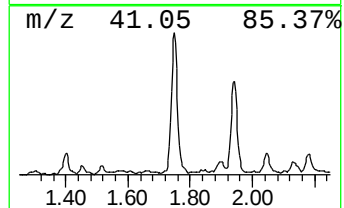
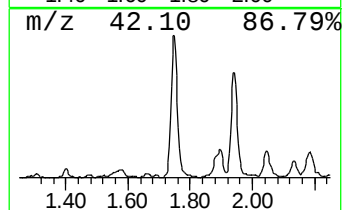
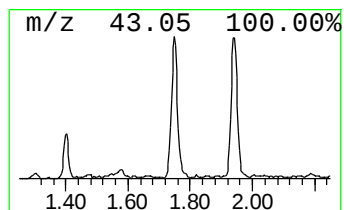
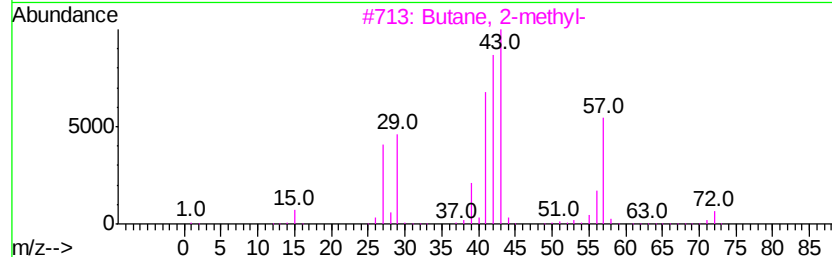
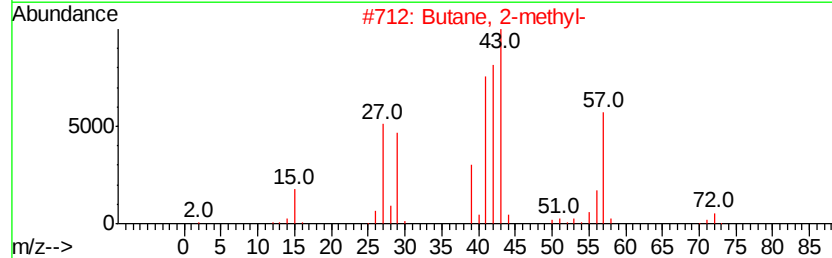
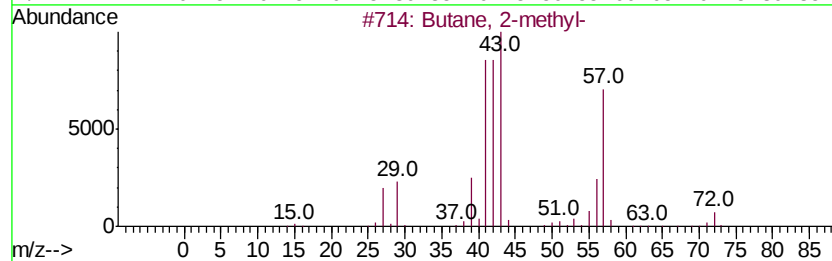
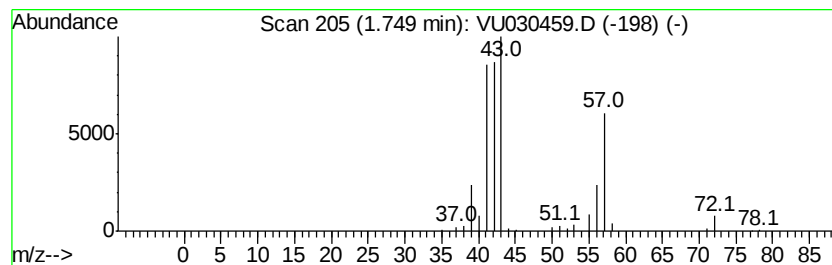
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM032319WMA.M
Quant Title : VOC Analysis

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.75	4.01 ug/L	90052	1,4-Difluorobenzene	5.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methyl-	72	C5H12	000078-78-4	91
2		Butane, 2-methyl-	72	C5H12	000078-78-4	86
3		Butane, 2-methyl-	72	C5H12	000078-78-4	86
4		3-Buten-1-ol	72	C4H8O	000627-27-0	47
5		Pentane	72	C5H12	000109-66-0	39



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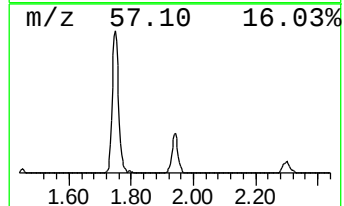
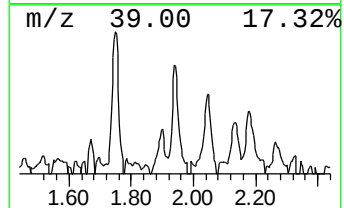
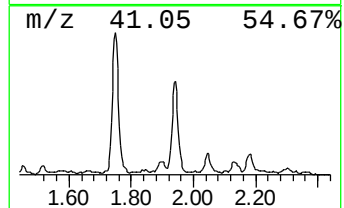
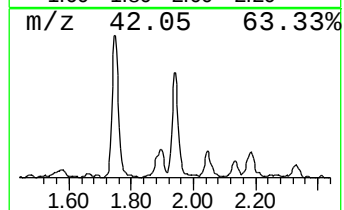
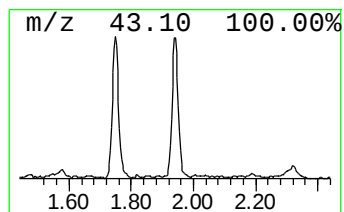
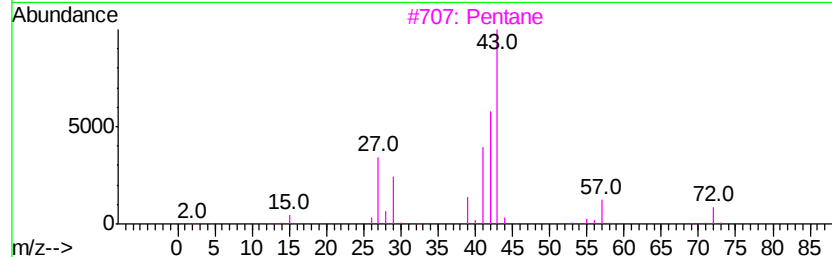
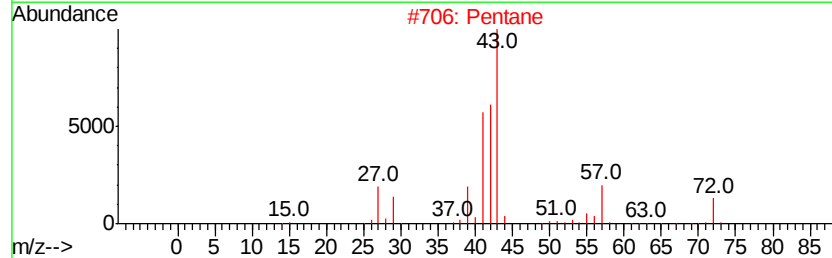
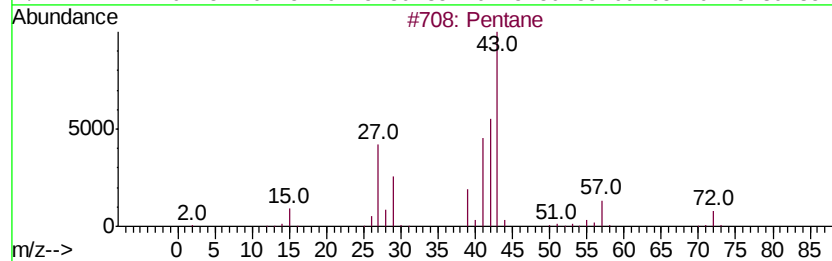
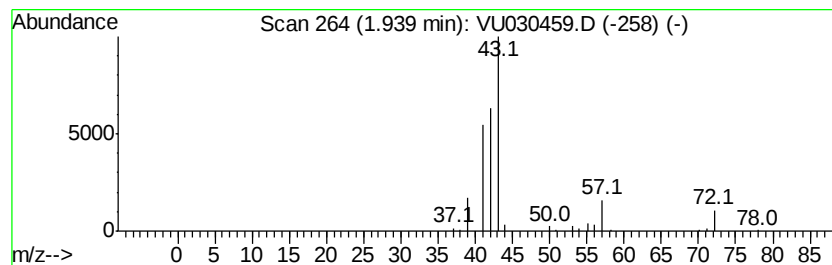
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Quant Title : VOC Analysis

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.94	2.70 ug/L	60544	1,4-Difluorobenzene	5.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane	72	C5H12	000109-66-0	86
2		Pentane	72	C5H12	000109-66-0	78
3		Pentane	72	C5H12	000109-66-0	72
4		Pentane	72	C5H12	000109-66-0	59
5		Pentane, 2-methyl-	86	C6H14	000107-83-5	39



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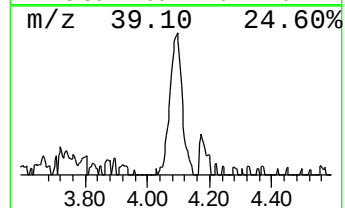
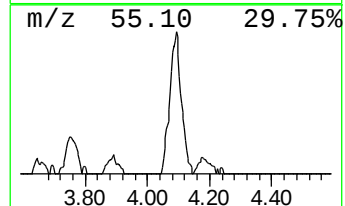
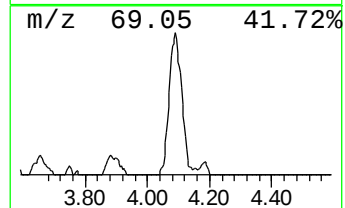
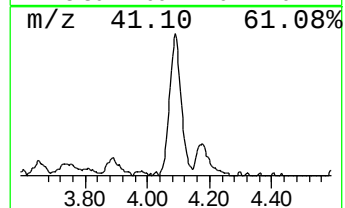
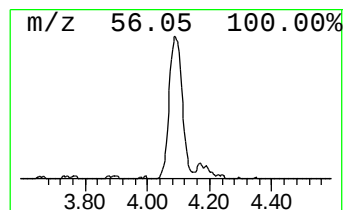
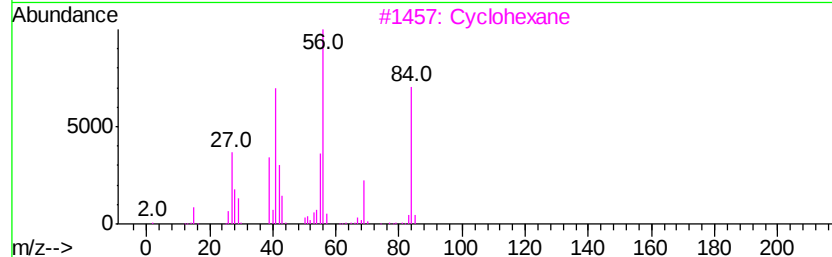
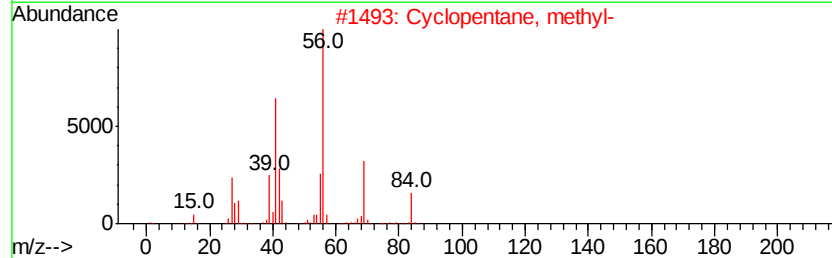
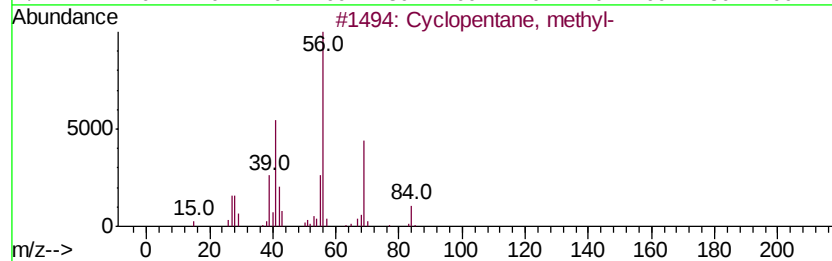
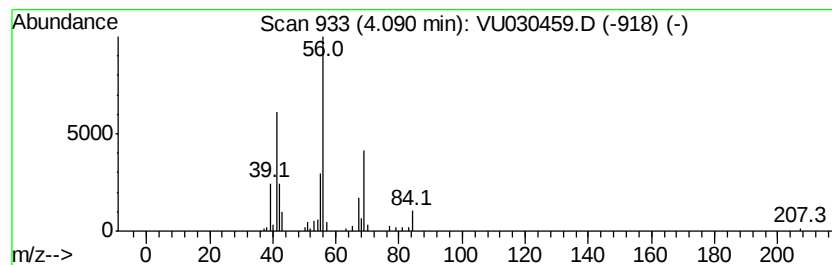
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Quant Title : VOC Analysis

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

Peak Number 3 (DEL) Alkane: Cyclic4.09 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.09	4.22 ug/L	94638	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	87
3		Cyclohexane	84	C6H12	000110-82-7	80
4		Cyclohexane	84	C6H12	000110-82-7	78
5		Cyclohexane	84	C6H12	000110-82-7	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU032619\
Data File : VU030459.D
Acq On : 26 Mar 2019 21:27
Operator : JC/SP
Sample : K2063-04DL 100X
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 33 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
DB6R6DL

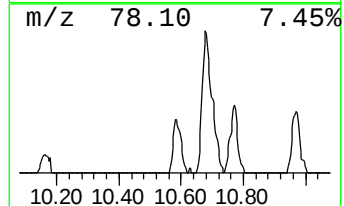
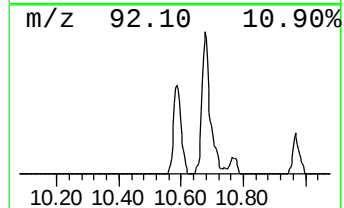
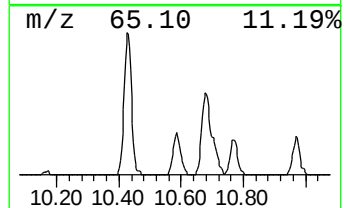
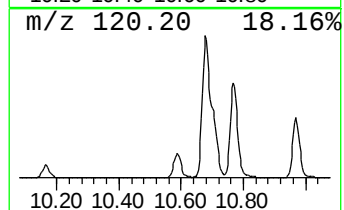
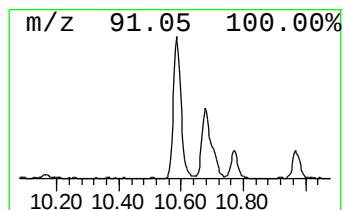
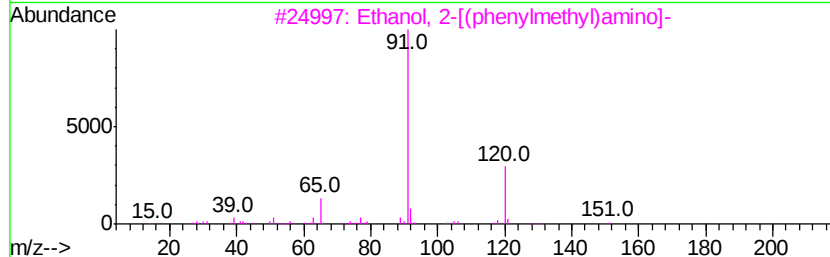
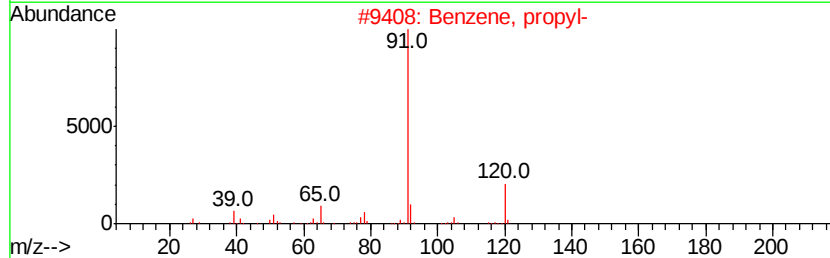
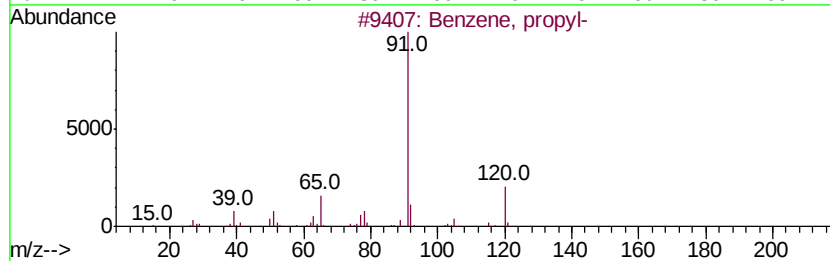
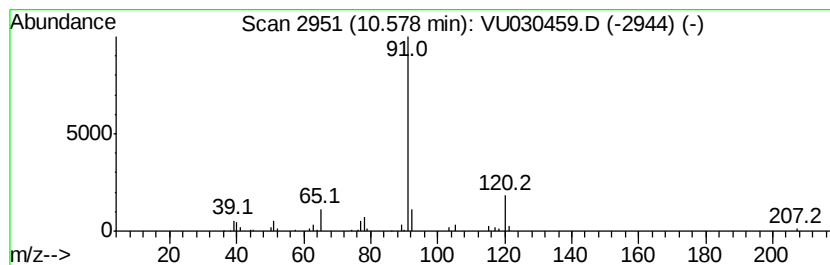
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Quant Title : VOC Analysis

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

Peak Number 5 Benzene, propyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.58	2.67 ug/L	65534	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, propyl-	120	C9H12	000103-65-1	90
2		Benzene, propyl-	120	C9H12	000103-65-1	87
3		Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	72
4		N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	72
5		Benzene, propyl-	120	C9H12	000103-65-1	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU032619\
Data File : VU030459.D
Acq On : 26 Mar 2019 21:27
Operator : JC/SP
Sample : K2063-04DL 100X
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 33 Sample Multiplier: 1

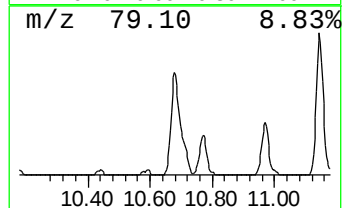
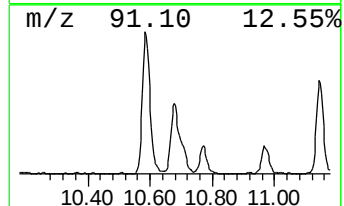
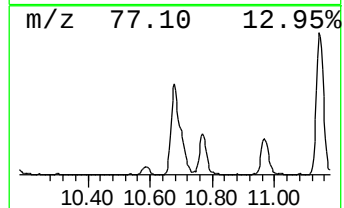
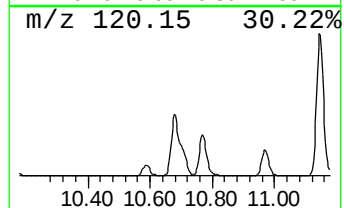
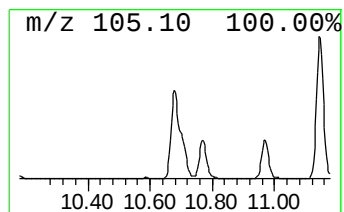
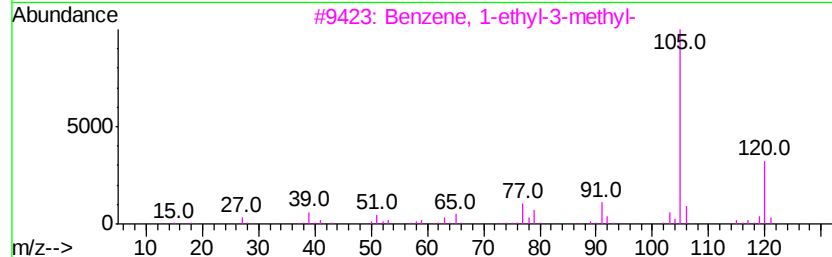
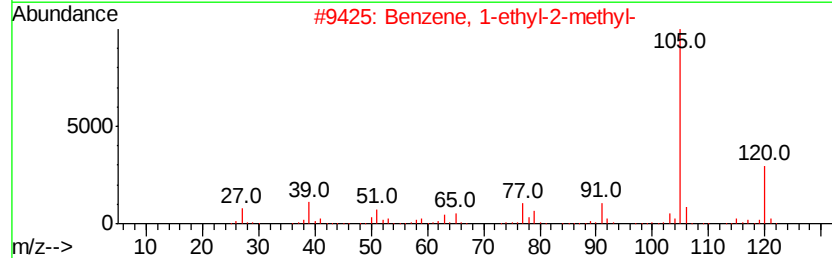
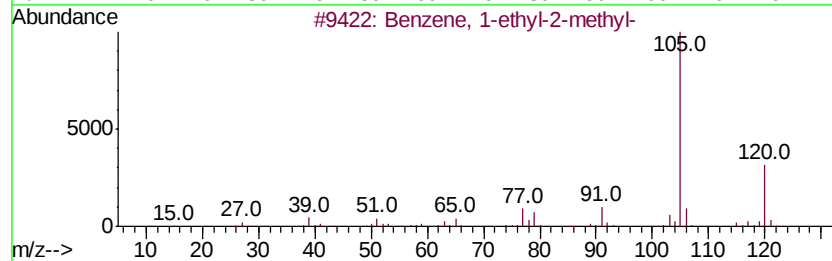
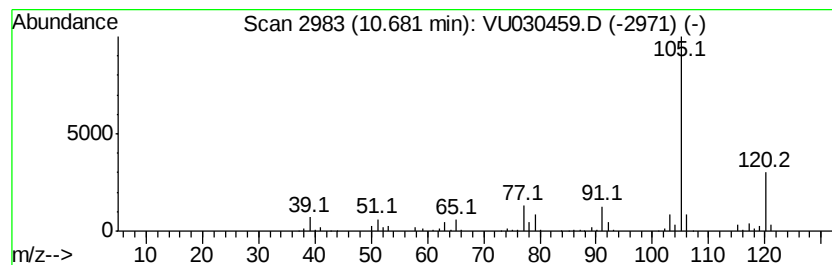
Instrument :
MSVOA_U
ClientSampleId :
DB6R6DL

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-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.68	17.66 ug/L	433152	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 94
3		Benzene, 1-ethyl-3-methyl-	120 C9H12	000620-14-4 94
4		Benzene, 1-ethyl-3-methyl-	120 C9H12	000620-14-4 91
5		Mesitylene	120 C9H12	000108-67-8 91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU032619\
Data File : VU030459.D
Acq On : 26 Mar 2019 21:27
Operator : JC/SP
Sample : K2063-04DL 100X
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 33 Sample Multiplier: 1

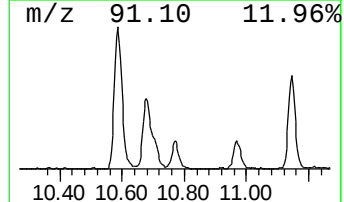
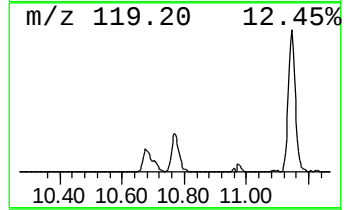
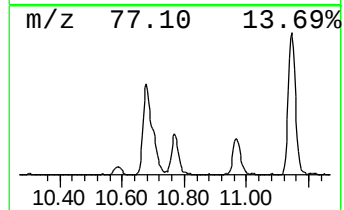
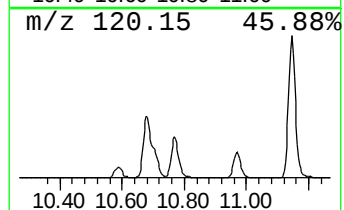
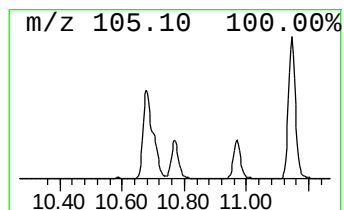
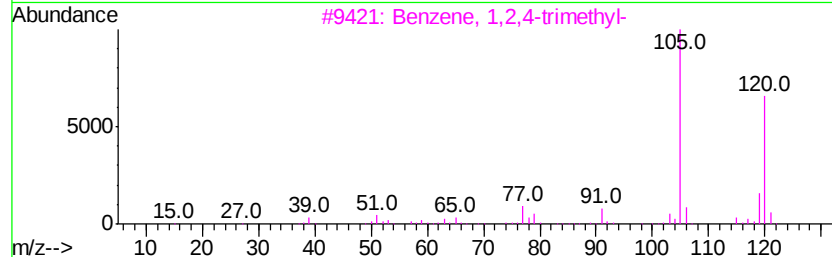
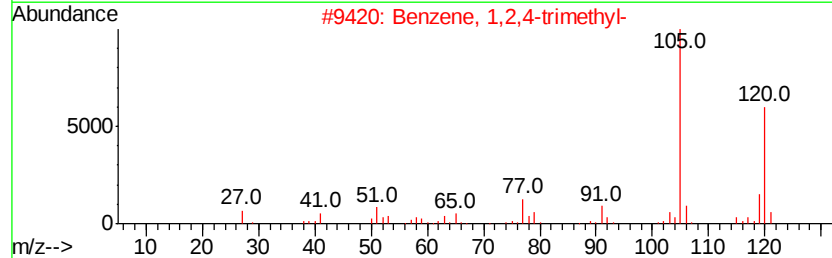
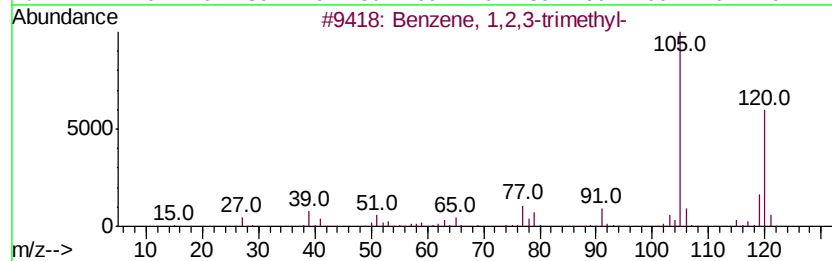
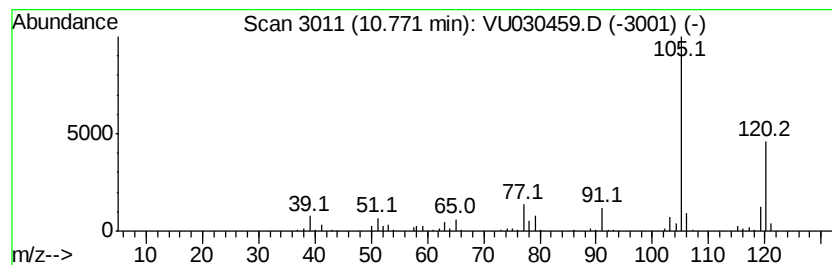
Instrument :
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ClientSampleId :
DB6R6DL

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 7 Benzene, 1,2,3-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.77	6.20 ug/L	152114	1,4-Dichlorobenzene-d4	11.48	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4	Mesitylene	120	C9H12	000108-67-8	94
5	Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU032619\
Data File : VU030459.D
Acq On : 26 Mar 2019 21:27
Operator : JC/SP
Sample : K2063-04DL 100X
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 33 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
DB6R6DL

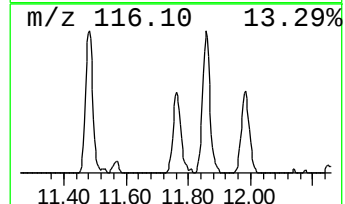
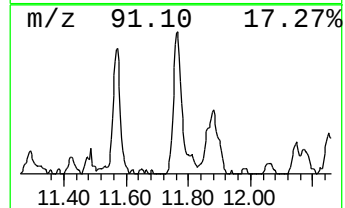
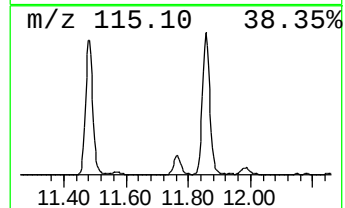
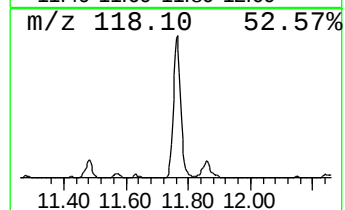
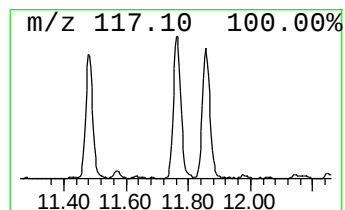
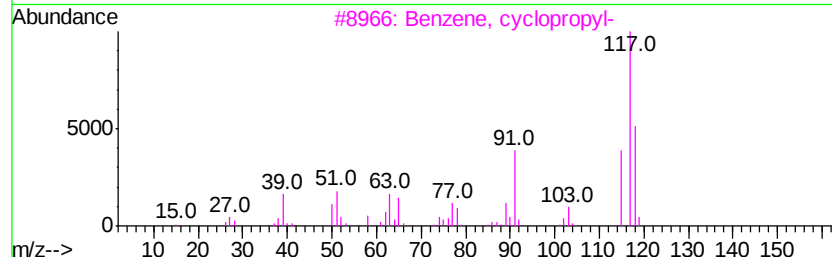
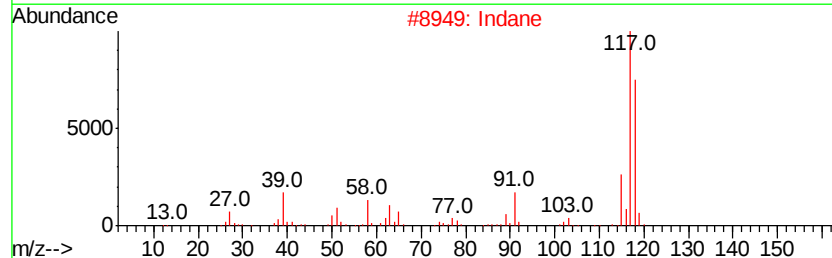
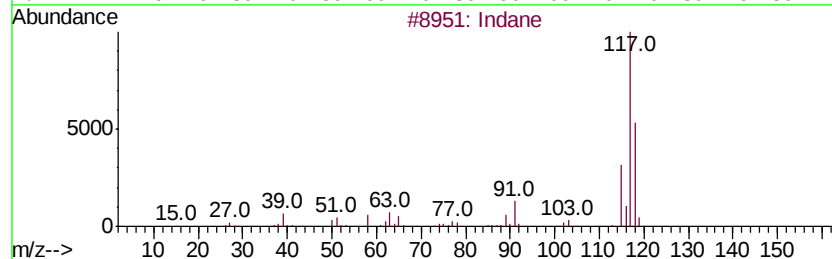
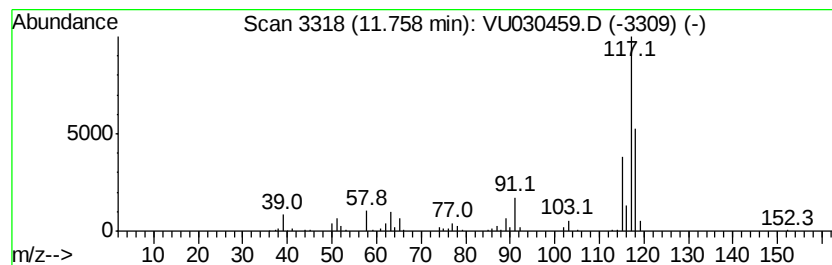
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Quant Title : VOC Analysis

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

Peak Number 11 Indane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.76	6.26 ug/L	153551	1,4-Dichlorobenzene-d4	11.48

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	95
2	Indane		118	C9H10	000496-11-7	95
3	Benzene, cvclopopyl-		118	C9H10	000873-49-4	94
4	Indane		118	C9H10	000496-11-7	94
5	Benzene, 1-ethenyl-2-methyl-		118	C9H10	000611-15-4	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU032619\
Data File : VU030459.D
Acq On : 26 Mar 2019 21:27
Operator : JC/SP
Sample : K2063-04DL 100X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 33 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
DB6R6DL

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: Str...	1.75	4.0	ug/L	90052	1	5.88	1121880	50.0
(DEL) Alkane: Str...	1.94	2.7	ug/L	60544	1	5.88	1121880	50.0
(DEL) Alkane: Cyc...	4.09	4.2	ug/L	94638	1	5.88	1121880	50.0
Benzene, propyl-	10.58	2.7	ug/L	65534	3	11.48	1226710	50.0
Benzene, 1-ethyl-...	10.68	17.7	ug/L	433152	3	11.48	1226710	50.0
Benzene, 1,2,3-tr...	10.77	6.2	ug/L	152114	3	11.48	1226710	50.0
Indane	11.76	6.3	ug/L	153551	3	11.48	1226710	50.0