

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051222\
 Data File : VU048539.D
 Acq On : 12 May 2022 14:16
 Operator : SY/MD
 Sample : N2801-06
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 EXLB9

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM050622WMA.M
 Title : VOC Analysis

Signal : TIC: VU048539.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.601	74	81	94	rVB	112838	126494	11.57%	1.357%
2	1.913	170	178	192	rBV	88709	113490	10.38%	1.217%
3	2.517	364	366	369	rBV2	329	148	0.01%	0.002%
4	2.568	370	382	399	rBV	221019	371868	34.01%	3.988%
5	2.710	420	426	429	rBV2	458	583	0.05%	0.006%
6	2.765	436	443	447	rBV3	1772	2367	0.22%	0.025%
7	2.958	499	503	505	rBV	269	233	0.02%	0.002%
8	3.044	514	530	550	rBV4	15135	34843	3.19%	0.374%
9	3.134	550	558	560	rBV4	1228	1147	0.10%	0.012%
10	3.359	614	628	643	rBV3	16133	34973	3.20%	0.375%
11	3.446	652	655	658	rBV	226	154	0.01%	0.002%
12	3.482	664	666	669	rBV	267	181	0.02%	0.002%
13	3.568	690	693	694	rBV2	337	198	0.02%	0.002%
14	3.652	716	719	721	rBV	291	234	0.02%	0.003%
15	3.758	749	752	753	rBV	256	163	0.01%	0.002%
16	3.800	763	765	767	rBV	311	219	0.02%	0.002%
17	3.880	786	790	792	rBV	221	221	0.02%	0.002%
18	4.089	849	855	856	rBV2	992	912	0.08%	0.010%
19	4.157	873	876	879	rBV	302	303	0.03%	0.003%
20	4.218	892	895	897	rBV	212	169	0.02%	0.002%
21	4.340	930	933	936	rBV2	454	384	0.04%	0.004%
22	4.440	953	964	970	rBV5	1548	2675	0.24%	0.029%
23	4.533	978	993	1008	rBV3	28550	68246	6.24%	0.732%
24	4.620	1008	1020	1054	rBV	113186	282204	25.81%	3.027%
25	4.803	1074	1077	1085	rVB3	697	693	0.06%	0.007%
26	5.064	1143	1158	1176	rBV	205210	447935	40.97%	4.804%
27	5.301	1228	1232	1235	rVV2	856	722	0.07%	0.008%
28	5.391	1243	1260	1272	rBV3	18964	41342	3.78%	0.443%
29	5.588	1317	1321	1327	rBV4	870	886	0.08%	0.010%
30	5.639	1327	1337	1344	rBV7	2016	4125	0.38%	0.044%
31	5.726	1344	1364	1392	rVV2	402987	1093322	100.00%	11.726%
32	6.035	1456	1460	1466	rBV3	498	588	0.05%	0.006%
33	6.131	1487	1490	1493	rBV	400	261	0.02%	0.003%
34	6.170	1498	1502	1503	rBV2	404	256	0.02%	0.003%
35	6.250	1511	1527	1557	rBV	325353	628865	57.52%	6.745%
36	6.398	1566	1573	1582	rBV6	2230	3374	0.31%	0.036%

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Integration Parameters: LSCINT.P

Integrator: RTE

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Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM050622WMA.M
 Title : VOC Analysis

37	6.530	1604	1614	1619	rBV6	3726	6777	0.62%	0.073%
38	6.623	1640	1643	1646	rBV2	305	267	0.02%	0.003%
39	6.690	1650	1664	1678	rBV	246802	488327	44.66%	5.237%
40	6.854	1712	1715	1716	rBV	625	315	0.03%	0.003%
41	6.883	1720	1724	1728	rVB2	545	469	0.04%	0.005%
42	6.903	1728	1730	1734	rBV2	341	231	0.02%	0.002%
43	6.983	1748	1755	1762	rVB5	1011	1509	0.14%	0.016%
44	7.047	1771	1775	1779	rVB	675	598	0.05%	0.006%
45	7.102	1789	1792	1799	rVB	369	356	0.03%	0.004%
46	7.170	1800	1813	1814	rBV6	1703	2516	0.23%	0.027%
47	7.250	1834	1838	1839	rBV3	1074	865	0.08%	0.009%
48	7.391	1867	1882	1894	rVB9	9285	20780	1.90%	0.223%
49	7.504	1915	1917	1921	rBV2	274	182	0.02%	0.002%
50	7.565	1923	1936	1953	rBV2	142456	257710	23.57%	2.764%
51	7.761	1988	1997	2006	rVV6	3714	6248	0.57%	0.067%
52	7.899	2026	2040	2055	rBV	507243	867478	79.34%	9.304%
53	8.179	2114	2127	2153	rBV	109048	186596	17.07%	2.001%
54	8.475	2208	2219	2233	rVB2	5197	9529	0.87%	0.102%
55	8.552	2234	2243	2253	rVB4	9448	15402	1.41%	0.165%
56	8.633	2256	2268	2310	rBV	349766	649325	59.39%	6.964%
57	8.841	2331	2333	2336	rBV	301	185	0.02%	0.002%
58	8.887	2344	2347	2350	rVB2	501	353	0.03%	0.004%
59	9.054	2395	2399	2403	rVB2	404	334	0.03%	0.004%
60	9.076	2403	2406	2408	rBV	375	246	0.02%	0.003%
61	9.163	2429	2433	2435	rVV2	226	216	0.02%	0.002%
62	9.301	2473	2476	2479	rBV2	265	226	0.02%	0.002%
63	9.417	2499	2512	2534	rBV	469144	779487	71.30%	8.360%
64	9.703	2594	2601	2605	rBV4	1155	1316	0.12%	0.014%
65	9.735	2609	2611	2614	rBV3	441	222	0.02%	0.002%
66	9.874	2652	2654	2658	rVB2	333	210	0.02%	0.002%
67	10.095	2721	2723	2726	rBV	358	213	0.02%	0.002%
68	10.250	2768	2771	2773	rBV2	525	388	0.04%	0.004%
69	10.314	2788	2791	2795	rVB	271	165	0.02%	0.002%
70	10.349	2800	2802	2805	rBV	324	211	0.02%	0.002%
71	10.395	2813	2816	2819	rVB	381	261	0.02%	0.003%
72	10.440	2826	2830	2834	rBV2	307	308	0.03%	0.003%
73	10.485	2836	2844	2855	rVB2	4906	7868	0.72%	0.084%
74	10.542	2860	2862	2864	rBV2	353	200	0.02%	0.002%
75	10.655	2895	2897	2898	rBV	494	174	0.02%	0.002%
76	10.758	2916	2929	2949	rVB	367572	601250	54.99%	6.449%

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77	10.864	2958	2962	2964	rBV2	460	345	0.03%	0.004%
78	10.893	2964	2971	2982	rVB3	6270	10190	0.93%	0.109%
79	10.980	2995	2998	3000	rBV	322	186	0.02%	0.002%
80	11.060	3020	3023	3029	rBV2	558	484	0.04%	0.005%
81	11.292	3086	3095	3106	rVB2	21955	33121	3.03%	0.355%
82	11.459	3141	3147	3149	rBV4	1948	2103	0.19%	0.023%
83	11.581	3182	3185	3187	rBV2	304	221	0.02%	0.002%
84	11.613	3187	3195	3199	rBV2	1492	2056	0.19%	0.022%
85	11.742	3229	3235	3244	rVB4	1432	2078	0.19%	0.022%
86	11.812	3244	3257	3281	rBV	507871	813224	74.38%	8.722%
87	12.095	3329	3345	3360	rBV3	60725	104440	9.55%	1.120%
88	12.192	3362	3375	3393	rVB	545015	890290	81.43%	9.549%
89	12.304	3398	3410	3417	rBV7	6000	9321	0.85%	0.100%
90	12.375	3424	3432	3444	rVB	11667	17857	1.63%	0.192%
91	12.465	3451	3460	3472	rBV2	17884	24860	2.27%	0.267%
92	12.571	3483	3493	3501	rBV	33265	49906	4.56%	0.535%
93	12.681	3516	3527	3543	rBV2	40312	68455	6.26%	0.734%
94	12.973	3609	3618	3628	rBV	28826	42578	3.89%	0.457%
95	13.253	3701	3705	3708	rBV4	890	669	0.06%	0.007%
96	13.295	3709	3718	3728	rVV2	13314	19251	1.76%	0.206%
97	13.410	3750	3754	3759	rBV4	1169	1286	0.12%	0.014%
98	13.455	3759	3768	3779	rVB2	32670	47366	4.33%	0.508%
99	13.626	3813	3821	3830	rBV8	6055	8939	0.82%	0.096%
100	14.246	4012	4014	4017	rBV	539	403	0.04%	0.004%

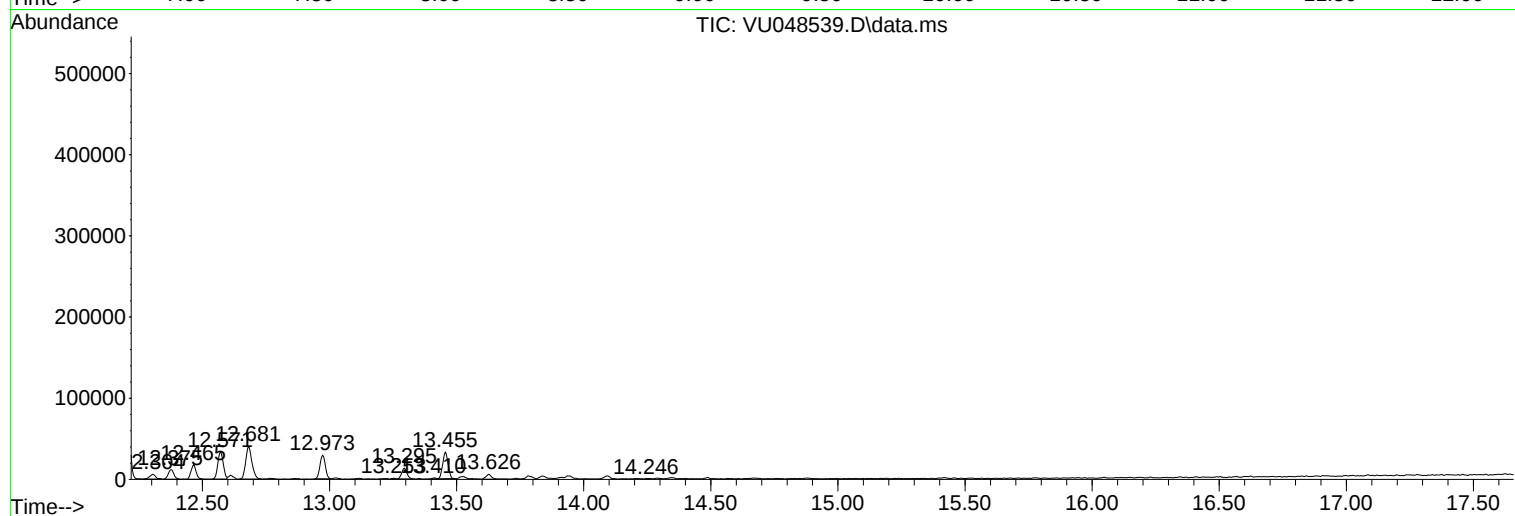
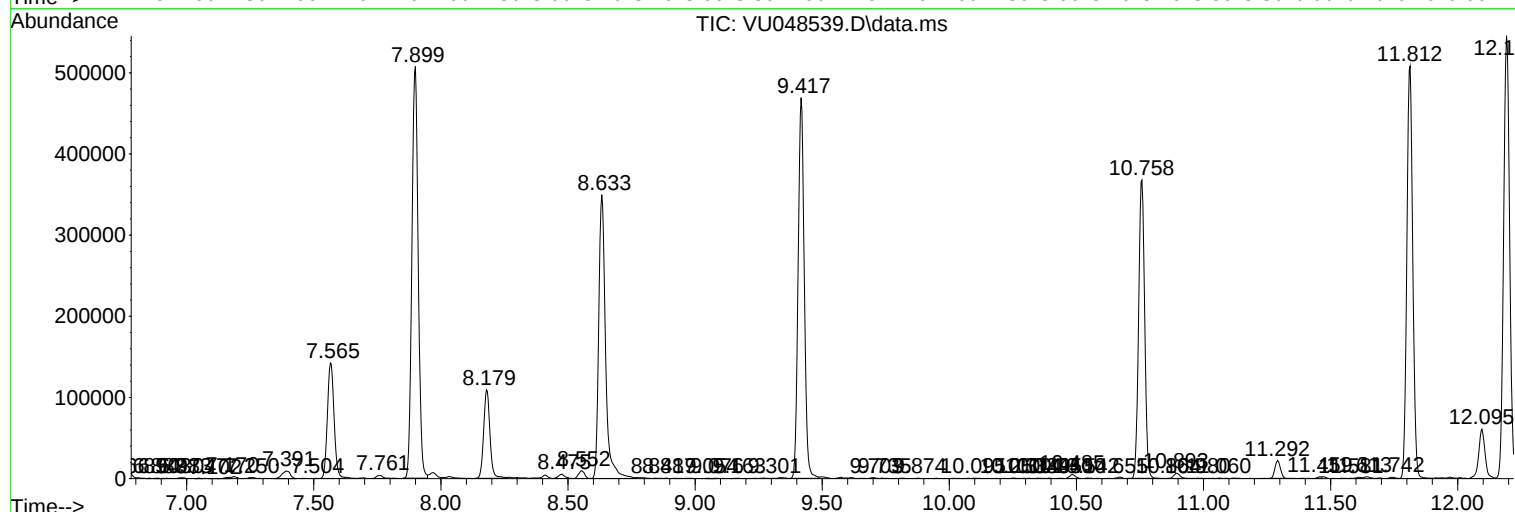
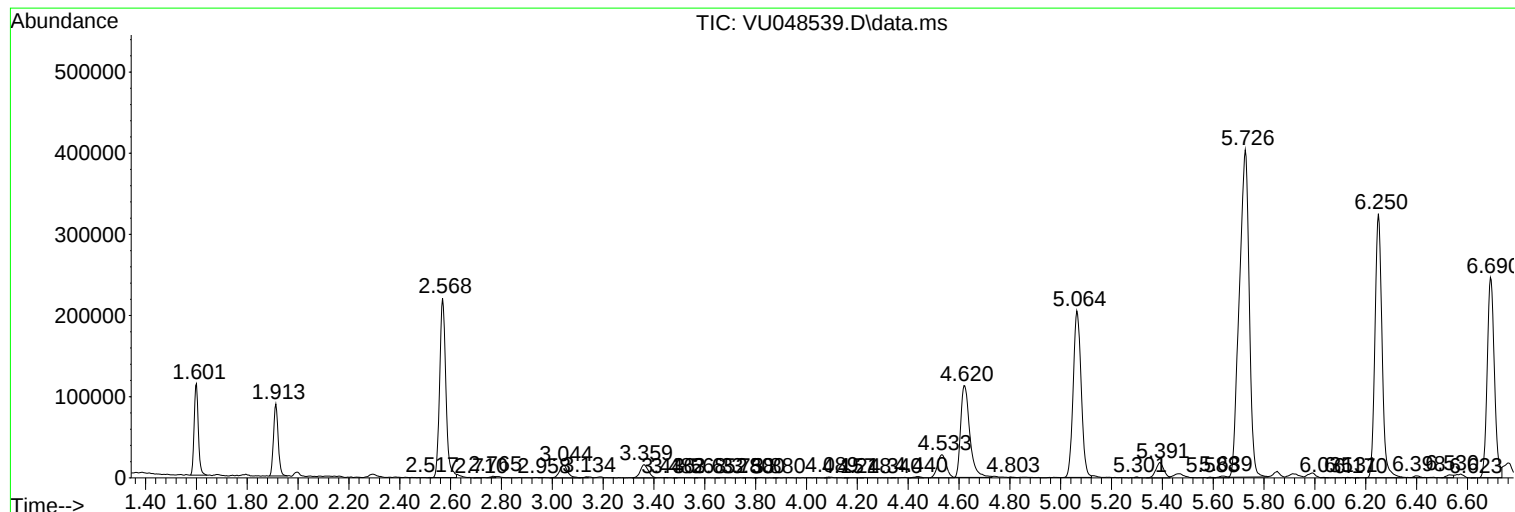
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Instrument :
MSVOA_U
ClientSampleId :
EXLB9

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM050622WMA.M
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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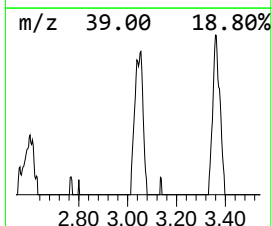
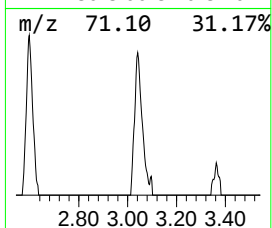
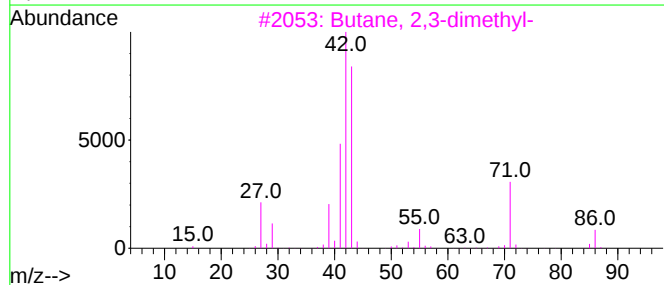
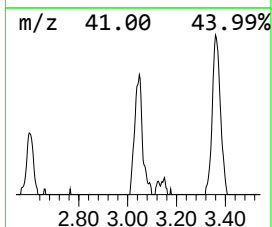
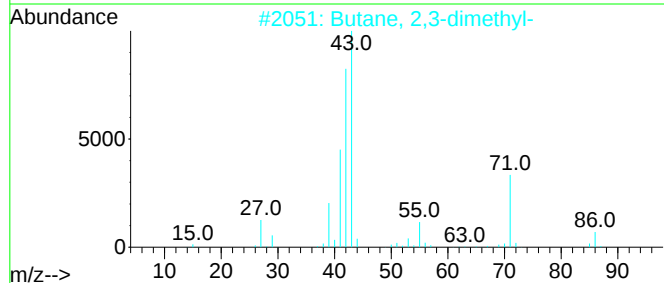
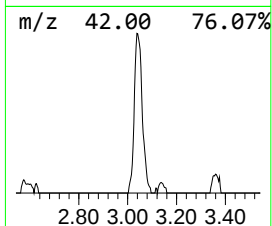
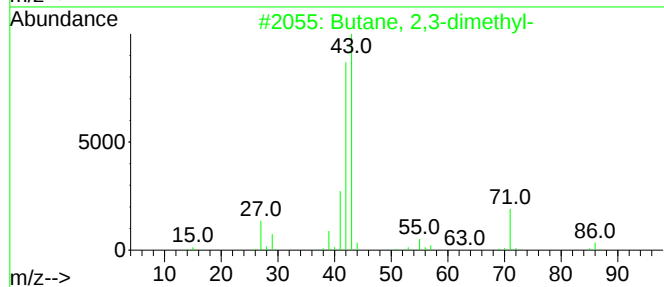
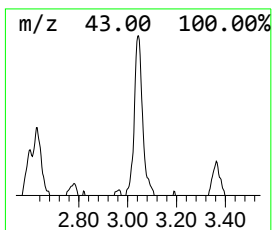
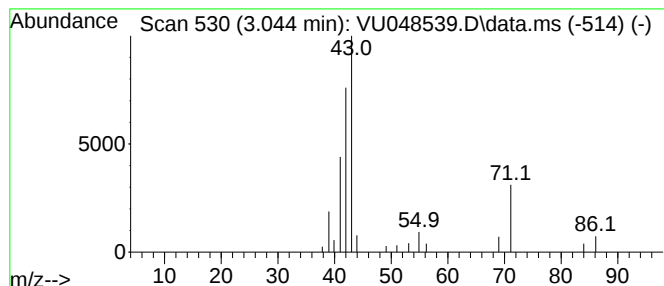
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM050622WMA.M
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.044	2.77 ug/L	34843	1,4-Difluorobenzene	6.250

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	80
2		Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	80
3		Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	64
4		Pentane, 2-methyl-	86	C6H14	000107-83-5	53
5		Pentane, 2-methyl-	86	C6H14	000107-83-5	53



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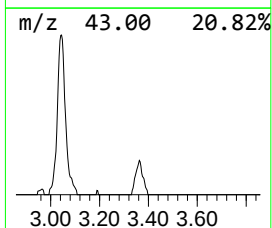
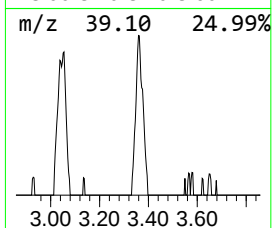
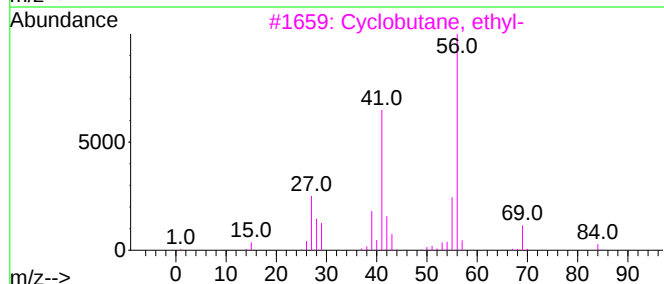
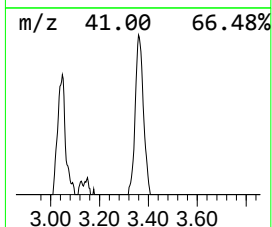
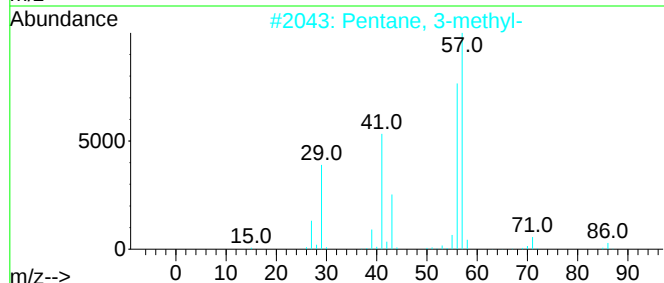
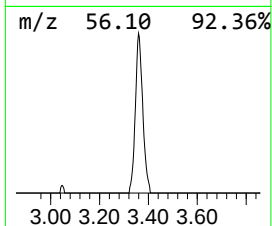
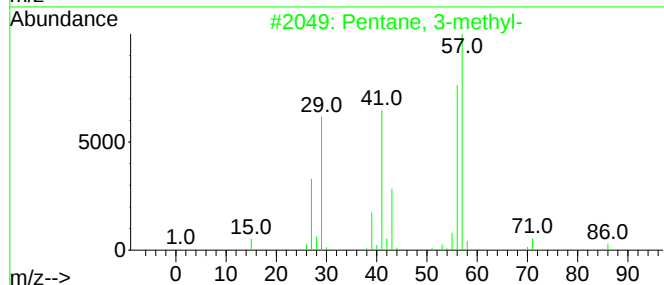
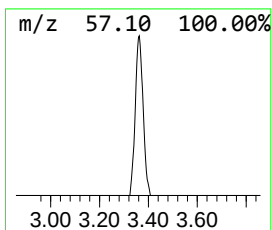
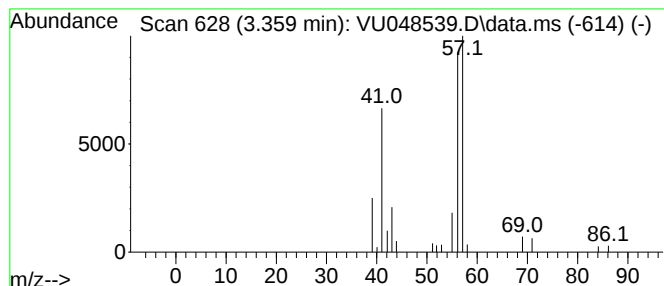
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM050622WMA.M
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.359	2.78 ug/L	34973	1,4-Difluorobenzene	6.250

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentane, 3-methyl-	86	C6H14	000096-14-0	64
2		Pentane, 3-methyl-	86	C6H14	000096-14-0	64
3		Cyclobutane, ethyl-	84	C6H12	004806-61-5	53
4		Cyclopropane, propyl-	84	C6H12	002415-72-7	47
5		Pentane, 3-methyl-	86	C6H14	000096-14-0	39



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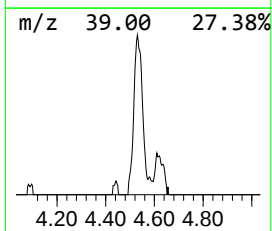
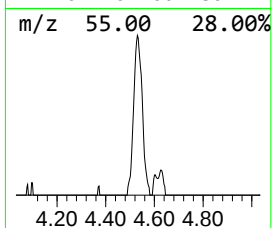
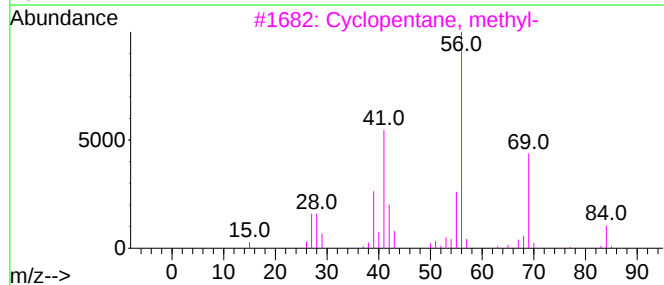
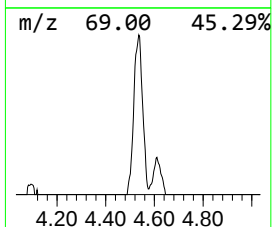
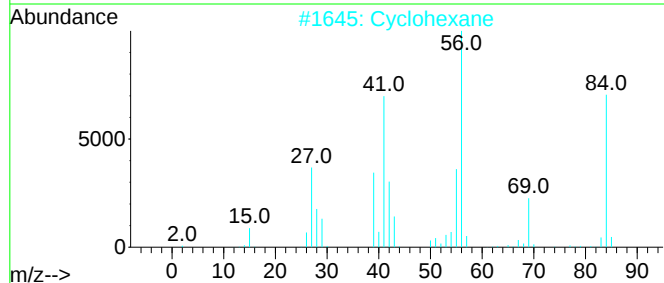
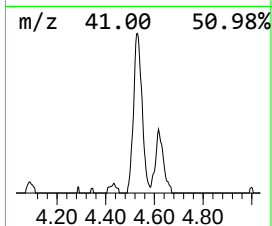
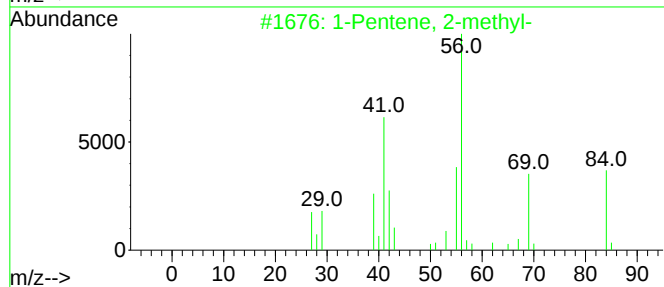
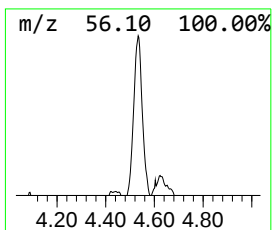
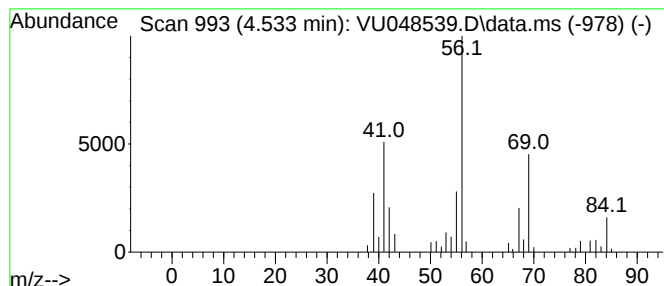
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM050622WMA.M
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.533	5.43 ug/L	68246	1,4-Difluorobenzene	6.250

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	74
2		Cyclohexane	84	C6H12	000110-82-7	72
3		Cyclopentane, methyl-	84	C6H12	000096-37-7	64
4		Cyclopentane, methyl-	84	C6H12	000096-37-7	64
5		Cyclohexane	84	C6H12	000110-82-7	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051222\
Data File : VU048539.D
Acq On : 12 May 2022 14:16
Operator : SY/MD
Sample : N2801-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXLB9

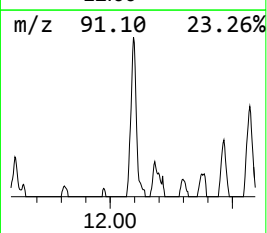
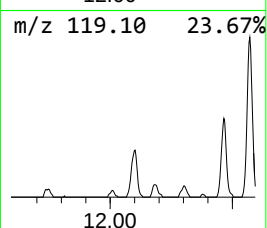
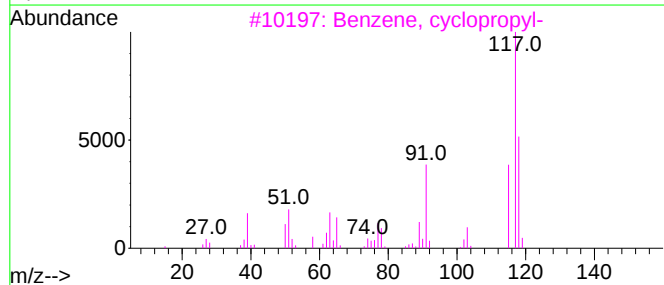
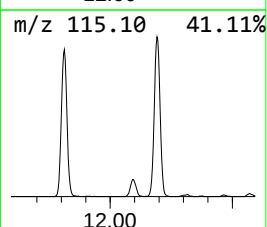
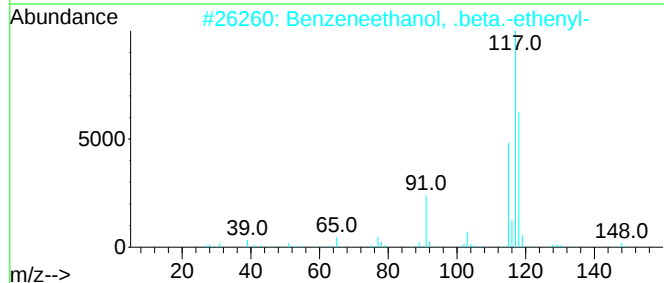
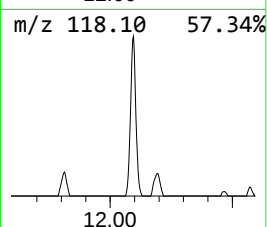
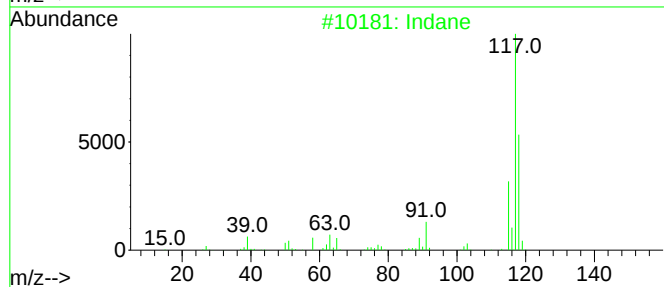
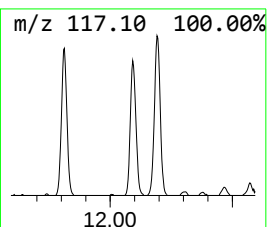
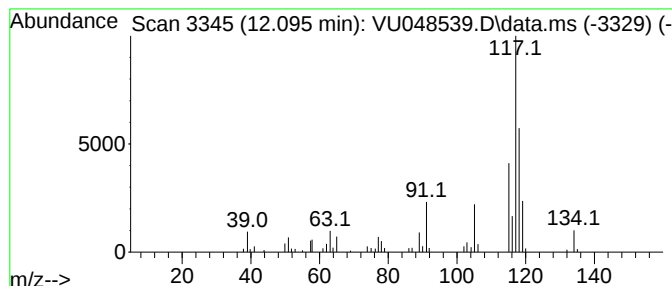
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.095	6.42 ug/L	104440	1,4-Dichlorobenzene-d4	11.813

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	60
2			Benzeneethanol, .beta.-ethenyl-	148	C10H12O	006052-63-7	59
3			Benzene, cyclopropyl-	118	C9H10	000873-49-4	55
4			Indane	118	C9H10	000496-11-7	53
5			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051222\
Data File : VU048539.D
Acq On : 12 May 2022 14:16
Operator : SY/MD
Sample : N2801-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXLB9

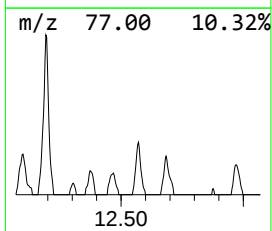
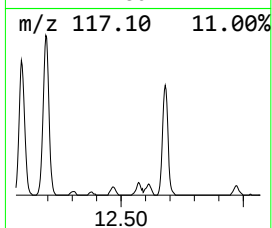
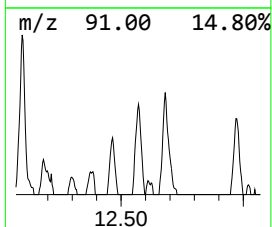
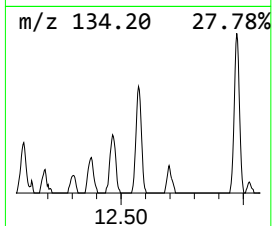
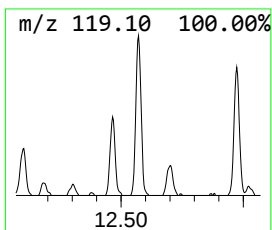
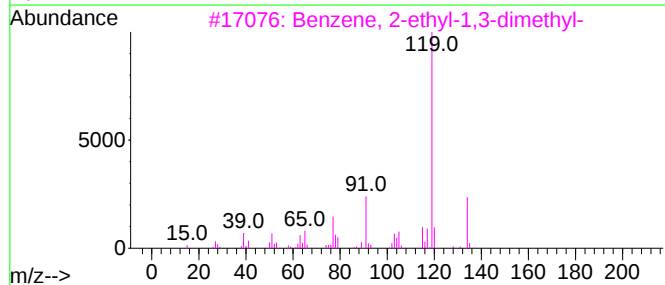
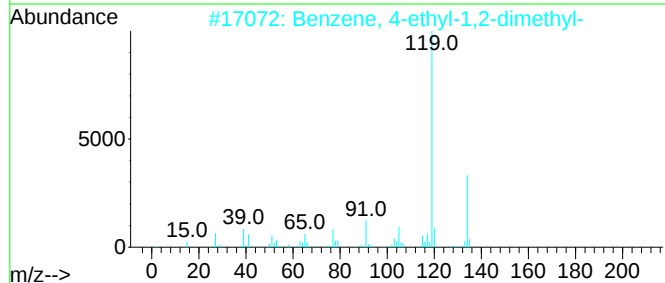
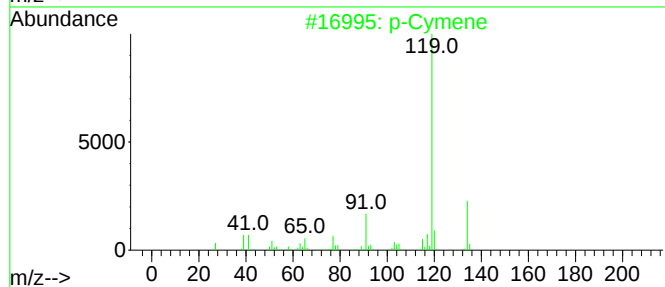
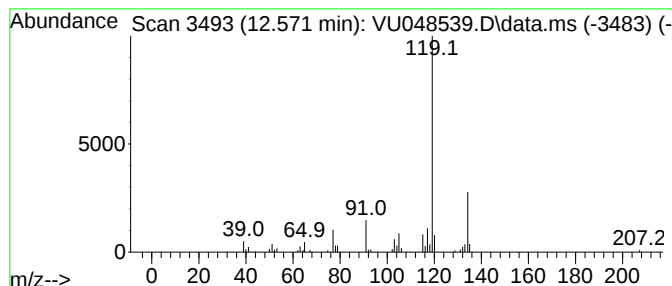
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 p-Cymene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.571	3.07 ug/L	49906	1,4-Dichlorobenzene-d4	11.813

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			p-Cymene	134	C10H14	000099-87-6	94
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
3			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	93
4			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	93
5			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051222\
Data File : VU048539.D
Acq On : 12 May 2022 14:16
Operator : SY/MD
Sample : N2801-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXLB9

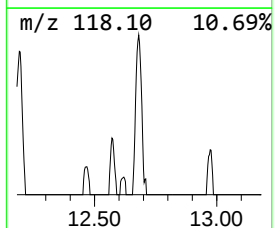
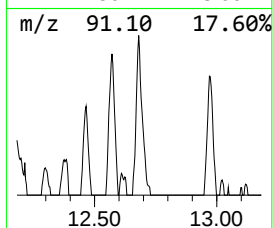
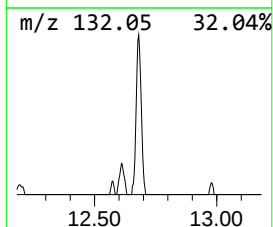
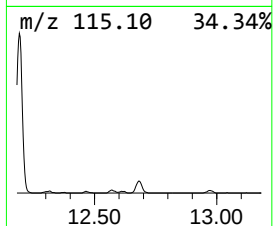
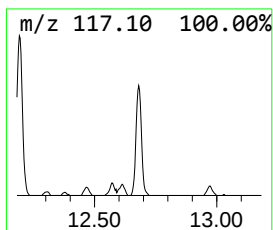
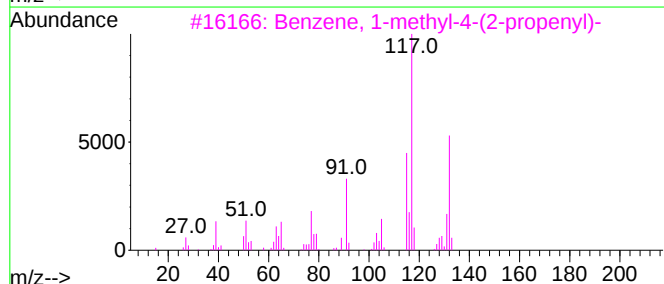
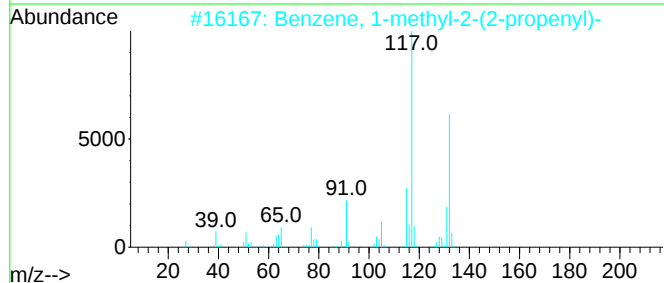
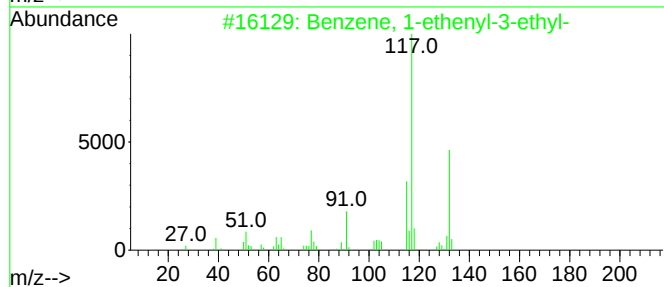
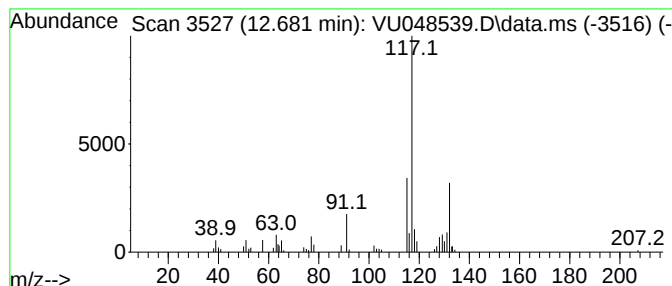
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.681	4.21 ug/L	68455	1,4-Dichlorobenzene-d4	11.813

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
2			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	87
3			Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	87
4			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
5			Indan, 1-methyl-	132	C10H12	000767-58-8	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051222\
Data File : VU048539.D
Acq On : 12 May 2022 14:16
Operator : SY/MD
Sample : N2801-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXLB9

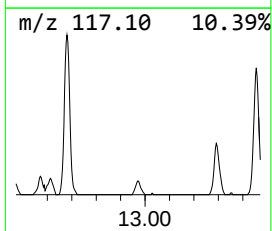
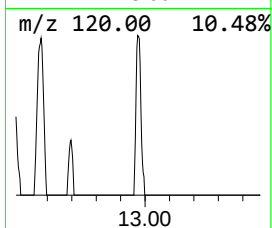
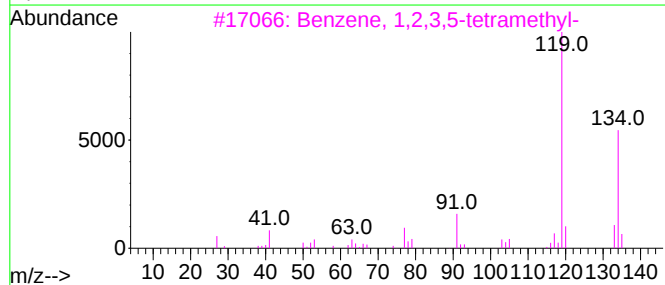
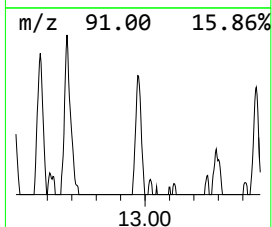
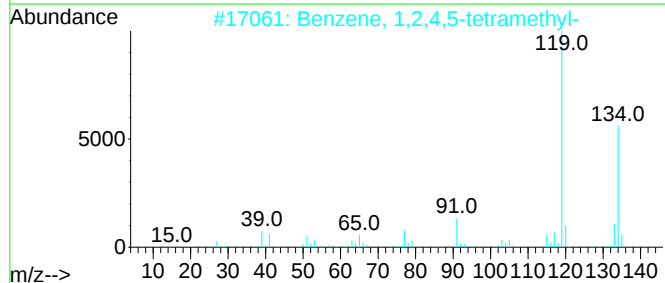
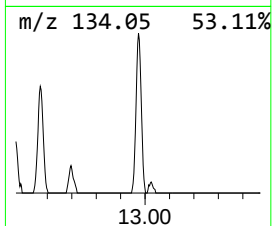
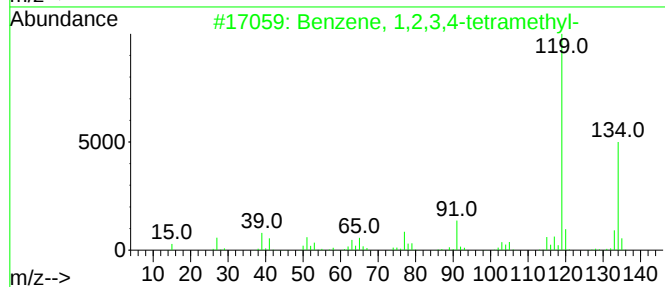
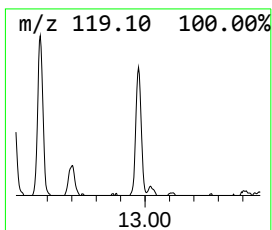
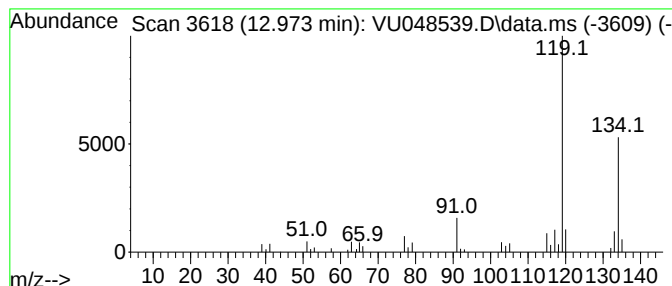
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.973	2.62 ug/L	42578	1,4-Dichlorobenzene-d4	11.813

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
4			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	96
5			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051222\
Data File : VU048539.D
Acq On : 12 May 2022 14:16
Operator : SY/MD
Sample : N2801-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXLB9

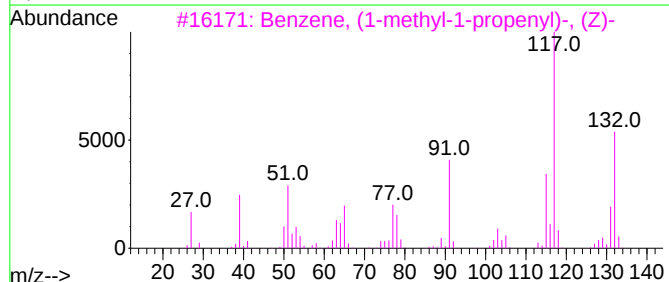
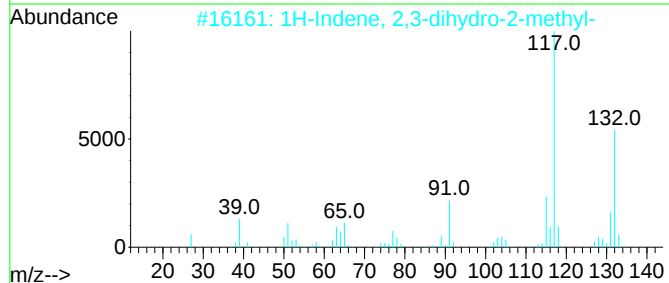
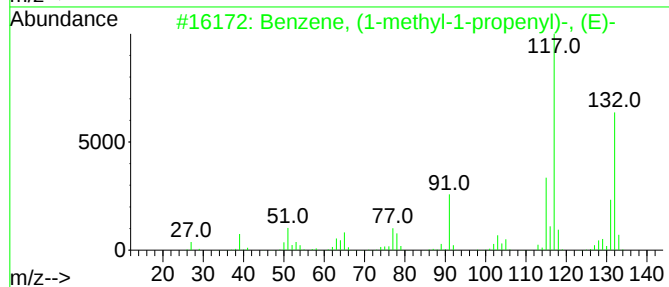
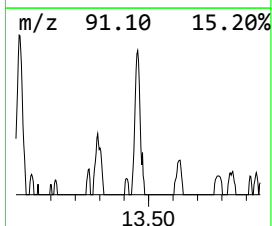
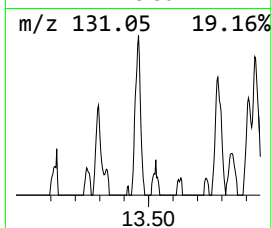
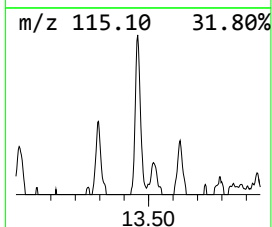
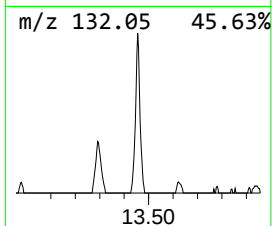
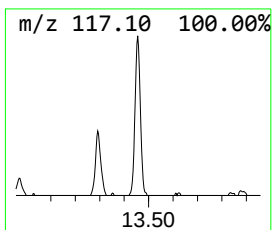
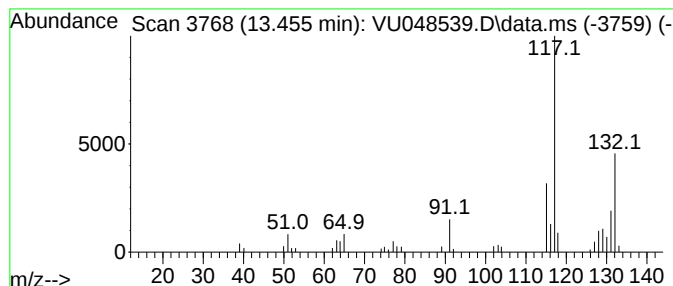
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Benzene, (1-methyl-1-propen... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.456	2.91 ug/L	47366	1,4-Dichlorobenzene-d4	11.813

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	86
2			1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	80
3			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	80
4			1-Phenyl-1-butene	132	C10H12	000824-90-8	80
5			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051222\
Data File : VU048539.D
Acq On : 12 May 2022 14:16
Operator : SY/MD
Sample : N2801-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXLB9

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM050622WMA.M
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	3.044	2.8	ug/L	34843	1	6.250	628865	50.0
(DEL) Alkane: S...	3.359	2.8	ug/L	34973	1	6.250	628865	50.0
(DEL) Alkane: S...	4.533	5.4	ug/L	68246	1	6.250	628865	50.0
Indane	12.095	6.4	ug/L	104440	3	11.813	813224	50.0
p-Cymene	12.571	3.1	ug/L	49906	3	11.813	813224	50.0
Benzene, 1-ethe...	12.681	4.2	ug/L	68455	3	11.813	813224	50.0
Benzene, 1,2,3,...	12.973	2.6	ug/L	42578	3	11.813	813224	50.0
Benzene, (1-met...	13.456	2.9	ug/L	47366	3	11.813	813224	50.0