

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120221\
 Data File : VU046046.D
 Acq On : 02 Dec 2021 12:50
 Operator : SY/MD
 Sample : M4868-07
 Misc : 6.94g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BGKP2

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\SFAMULM112921WMA.M
 Title : VOC Analysis

Signal : TIC: VU046046.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.485	41	45	51	rVB	2610	2638	0.32%	0.029%
2	1.597	72	80	101	rVB	85595	118099	14.19%	1.296%
3	1.880	157	168	170	rBV2	1624	2656	0.32%	0.029%
4	1.916	170	179	197	rVB	52367	91024	10.94%	0.999%
5	2.546	363	375	387	rBV	169420	276622	33.23%	3.035%
6	3.035	521	527	539	rBV4	1606	2941	0.35%	0.032%
7	3.182	558	573	587	rBV2	5479	13129	1.58%	0.144%
8	4.658	1014	1032	1075	rBV2	42045	192710	23.15%	2.114%
9	5.067	1141	1159	1183	rVV	131959	323971	38.92%	3.554%
10	5.723	1342	1363	1385	rBV2	326028	832364	100.00%	9.132%
11	5.980	1433	1443	1461	rBV5	3919	9943	1.19%	0.109%
12	6.250	1511	1527	1548	rBV	246847	501544	60.26%	5.503%
13	6.530	1603	1614	1627	rBV6	4631	11605	1.39%	0.127%
14	6.604	1633	1637	1650	rVB5	1952	3102	0.37%	0.034%
15	6.694	1650	1665	1691	rVB	180772	373495	44.87%	4.098%
16	7.186	1811	1818	1825	rBV3	1155	1733	0.21%	0.019%
17	7.568	1921	1937	1959	rBV	95152	188181	22.61%	2.065%
18	7.655	1959	1964	1970	rVV2	1218	1661	0.20%	0.018%
19	7.851	2019	2025	2026	rVV3	1104	1113	0.13%	0.012%
20	7.899	2026	2040	2058	rVV	359362	683351	82.10%	7.497%
21	7.970	2059	2062	2077	rVB4	5059	8266	0.99%	0.091%
22	8.182	2116	2128	2154	rBV2	59627	128158	15.40%	1.406%
23	8.536	2230	2238	2247	rVB4	1609	2097	0.25%	0.023%
24	8.681	2259	2283	2314	rBV	136058	426379	51.23%	4.678%
25	8.864	2334	2340	2349	rVB3	4440	6679	0.80%	0.073%
26	8.925	2349	2359	2367	rBV4	8197	15141	1.82%	0.166%
27	9.066	2392	2403	2411	rBV6	1615	3630	0.44%	0.040%
28	9.121	2411	2420	2426	rBV3	6520	10240	1.23%	0.112%
29	9.169	2426	2435	2442	rVV6	8036	15251	1.83%	0.167%
30	9.211	2442	2448	2461	rVB4	8012	14421	1.73%	0.158%
31	9.337	2475	2487	2500	rBV4	9900	17783	2.14%	0.195%
32	9.427	2500	2515	2535	rBV	290870	633225	76.08%	6.947%
33	9.510	2538	2541	2547	rVB4	927	1123	0.13%	0.012%
34	9.568	2547	2559	2570	rBV7	7421	14502	1.74%	0.159%
35	9.668	2570	2590	2596	rBV8	8521	20968	2.52%	0.230%
36	9.719	2597	2606	2613	rVV3	16763	26717	3.21%	0.293%

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Integrator: RTE

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Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
 Title : VOC Analysis

37	9.761	2613	2619	2640	rVB6	12507	28124	3.38%	0.309%
38	9.883	2646	2657	2668	rVB8	3274	6718	0.81%	0.074%
39	9.957	2668	2680	2689	rBV5	3809	7984	0.96%	0.088%
40	10.025	2689	2701	2714	rBV6	29674	73808	8.87%	0.810%
41	10.121	2723	2731	2739	rBV2	19995	29990	3.60%	0.329%
42	10.173	2741	2747	2755	rVB4	13264	20823	2.50%	0.228%
43	10.253	2755	2772	2785	rBV5	63358	136695	16.42%	1.500%
44	10.317	2786	2792	2794	rBV3	7622	8142	0.98%	0.089%
45	10.356	2795	2804	2811	rVV4	48410	90088	10.82%	0.988%
46	10.398	2813	2817	2828	rVB	35915	50880	6.11%	0.558%
47	10.468	2829	2839	2851	rVB7	14246	29768	3.58%	0.327%
48	10.558	2851	2867	2879	rBV5	42455	103412	12.42%	1.135%
49	10.626	2880	2888	2895	rVV2	24816	34962	4.20%	0.384%
50	10.674	2896	2903	2915	rVB4	9168	16225	1.95%	0.178%
51	10.771	2922	2933	2943	rBV	252728	430279	51.69%	4.721%
52	10.957	2985	2991	3001	rVB2	12704	19688	2.37%	0.216%
53	11.015	3002	3009	3017	rBV7	15629	27759	3.33%	0.305%
54	11.066	3017	3025	3033	rVB2	35553	54016	6.49%	0.593%
55	11.124	3034	3043	3056	rVB7	34272	75793	9.11%	0.832%
56	11.201	3058	3067	3074	rVB9	10331	19653	2.36%	0.216%
57	11.246	3075	3081	3087	rBV6	8211	11825	1.42%	0.130%
58	11.314	3089	3102	3113	rBV8	37683	73982	8.89%	0.812%
59	11.378	3115	3122	3127	rBV3	29894	41770	5.02%	0.458%
60	11.417	3127	3134	3142	rVB	88897	121086	14.55%	1.328%
61	11.465	3142	3149	3155	rBV5	23127	41491	4.98%	0.455%
62	11.574	3174	3183	3190	rBV2	40832	70615	8.48%	0.775%
63	11.642	3199	3204	3214	rVB4	28777	41949	5.04%	0.460%
64	11.706	3214	3224	3233	rBV6	49612	96430	11.59%	1.058%
65	11.819	3249	3259	3271	rVV	406686	661436	79.46%	7.257%
66	11.877	3271	3277	3284	rVB5	21014	30669	3.68%	0.336%
67	11.944	3292	3298	3302	rBV	18174	25567	3.07%	0.281%
68	12.086	3336	3342	3343	rBV5	7266	6998	0.84%	0.077%
69	12.198	3361	3377	3390	rBV	412540	722013	86.74%	7.922%
70	12.327	3411	3417	3429	rVV5	25595	45117	5.42%	0.495%
71	12.404	3429	3441	3450	rVB7	29605	73462	8.83%	0.806%
72	12.494	3458	3469	3474	rBV9	23322	50824	6.11%	0.558%
73	12.607	3497	3504	3514	rBV5	16109	28455	3.42%	0.312%
74	12.668	3515	3523	3531	rVB2	34985	53995	6.49%	0.592%
75	12.841	3571	3577	3590	rVB6	41096	72745	8.74%	0.798%
76	12.935	3592	3606	3624	rBV6	31797	109094	13.11%	1.197%

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Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
 Title : VOC Analysis

77	13.057	3637	3644	3651	rBV5	18422	26841	3.22%	0.294%
78	13.195	3681	3687	3702	rVB5	9421	19456	2.34%	0.213%
79	13.356	3732	3737	3749	rVB5	4871	6770	0.81%	0.074%
80	13.587	3801	3809	3817	rBV	45511	76384	9.18%	0.838%
81	13.729	3848	3853	3868	rBV	7673	14029	1.69%	0.154%
82	14.050	3946	3953	3959	rBV3	12178	16878	2.03%	0.185%
83	14.089	3959	3965	3971	rVV	19166	27268	3.28%	0.299%
84	14.124	3972	3976	3985	rVB8	13759	19931	2.39%	0.219%
85	14.192	3986	3997	4011	rBV3	33584	64254	7.72%	0.705%
86	14.452	4074	4078	4083	rBV7	4857	5314	0.64%	0.058%
87	14.574	4112	4116	4122	rVB6	7034	7303	0.88%	0.080%
88	14.941	4225	4230	4235	rBV4	6695	8096	0.97%	0.089%
89	15.021	4249	4255	4264	rBV4	6337	11110	1.33%	0.122%
90	15.076	4267	4272	4278	rVB4	8700	11329	1.36%	0.124%
91	15.211	4304	4314	4327	rBV6	24169	50223	6.03%	0.551%
92	15.272	4329	4333	4347	rVB6	8411	13539	1.63%	0.149%
93	15.349	4350	4357	4360	rBV8	5310	7549	0.91%	0.083%
94	15.404	4371	4374	4382	rVB2	7516	8329	1.00%	0.091%
95	15.510	4402	4407	4416	rVB10	10490	17007	2.04%	0.187%
96	15.979	4549	4553	4564	rVB5	10533	13862	1.67%	0.152%
97	16.192	4614	4619	4629	rVB5	11620	16609	2.00%	0.182%
98	16.410	4680	4687	4694	rBV2	10362	15384	1.85%	0.169%
99	16.545	4725	4729	4737	rVB6	12899	16195	1.95%	0.178%
100	17.134	4905	4912	4921	rVB7	11443	20081	2.41%	0.220%

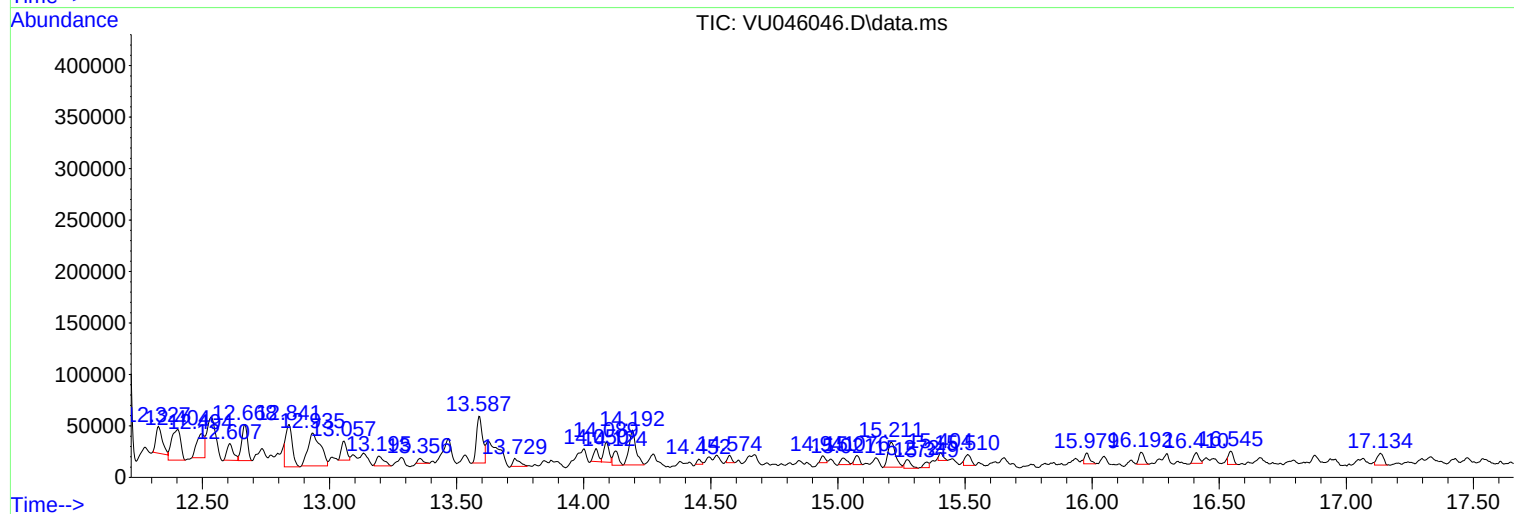
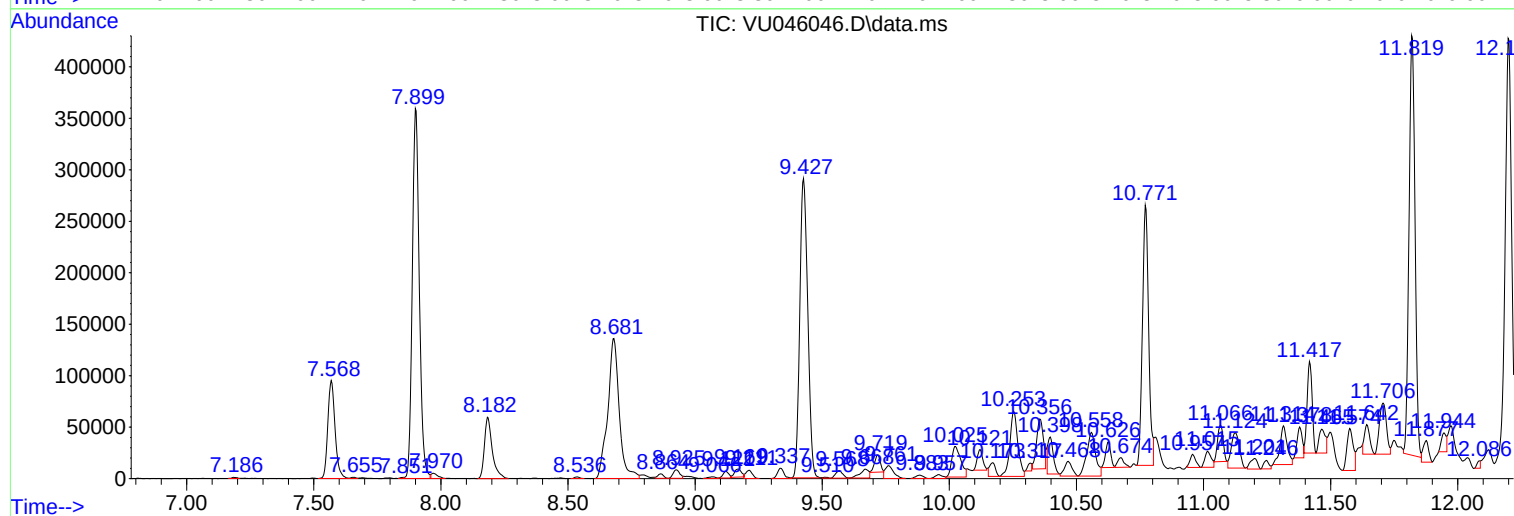
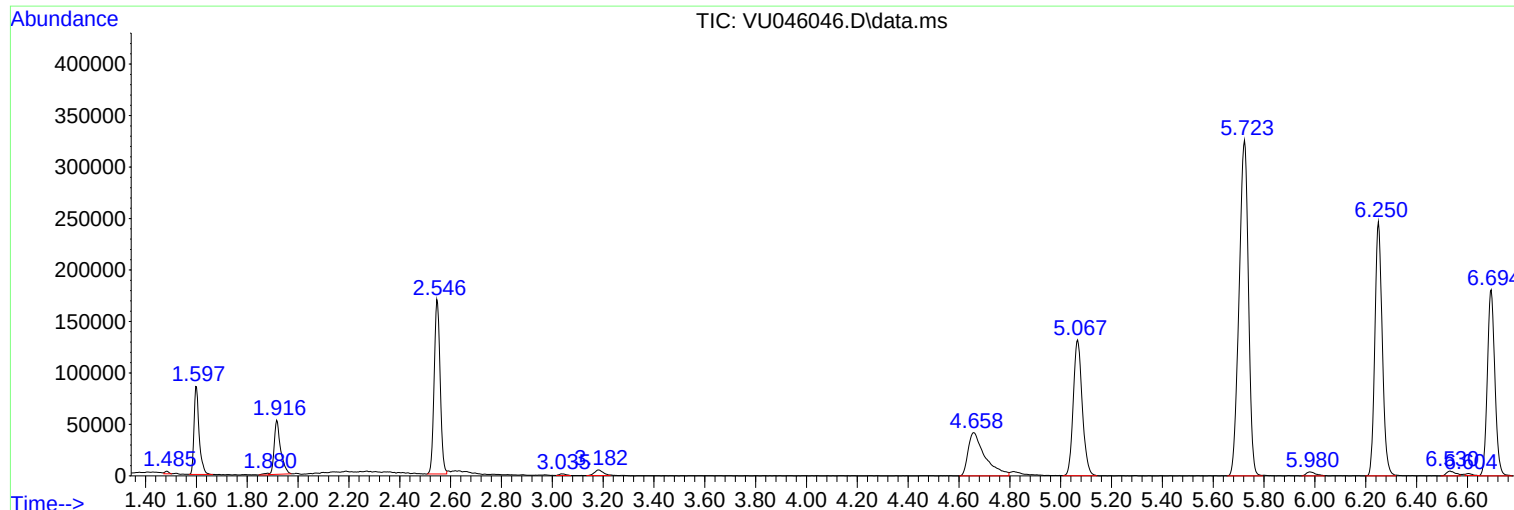
Sum of corrected areas: 9114533

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ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120221\
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ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGKP2

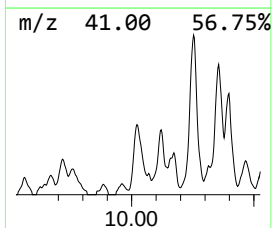
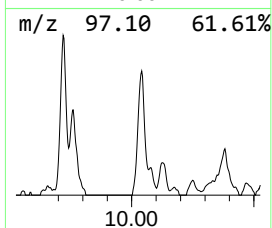
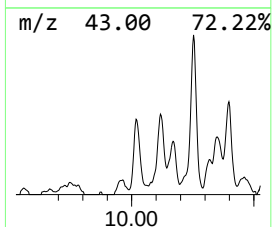
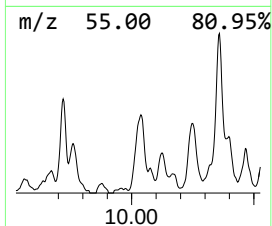
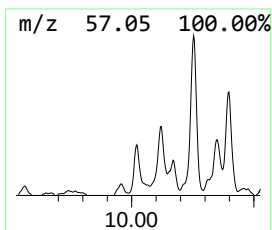
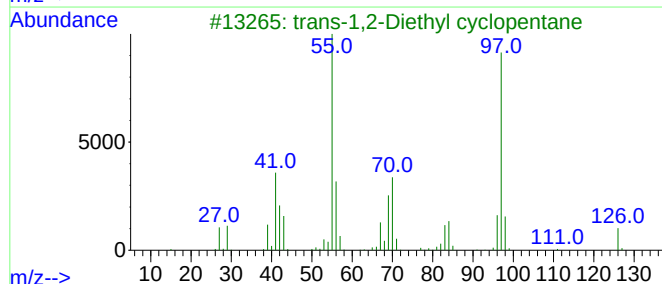
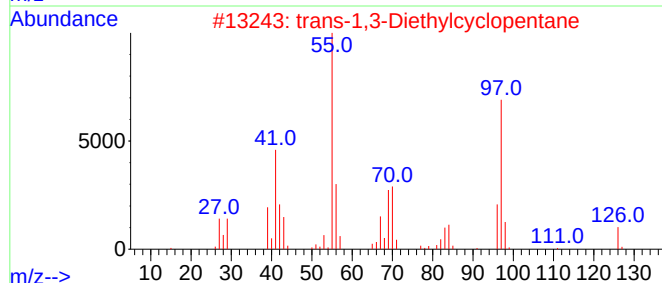
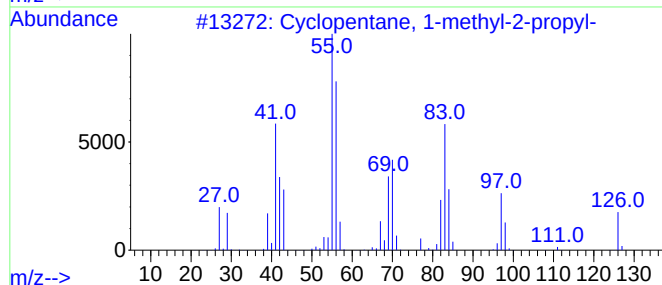
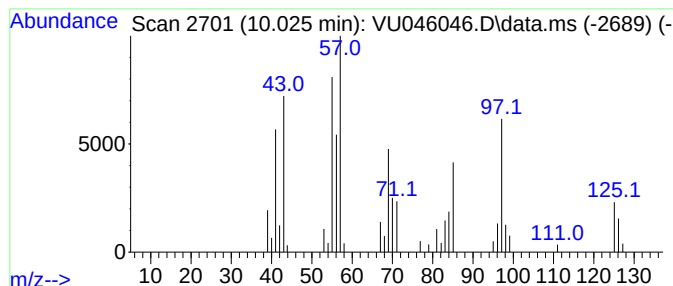
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

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

Peak Number 2 (DEL) Alkane: Cyclic10.025 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.025	5.83 ug/L	73808	Chlorobenzene-d5	9.427

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, 1-methyl-2-propyl-	126	C9H18	003728-57-2	38
2			trans-1,3-Diethylcyclopentane	126	C9H18	1000113-87-1	38
3			trans-1,2-Diethyl cyclopentane	126	C9H18	000932-40-1	35
4			Hexane, 1-(hexyloxy)-5-methyl-	200	C13H28O	074421-19-5	35
5			1-Pentanol, 2-ethyl-	116	C7H16O	027522-11-8	27



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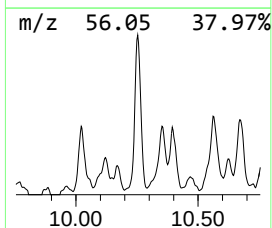
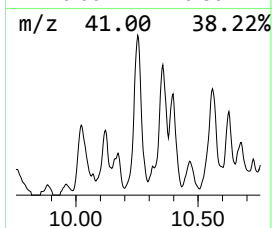
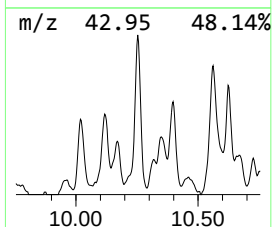
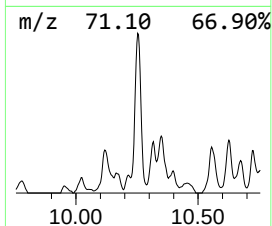
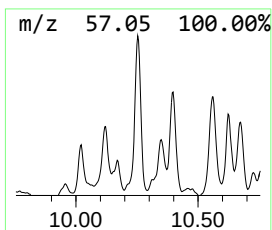
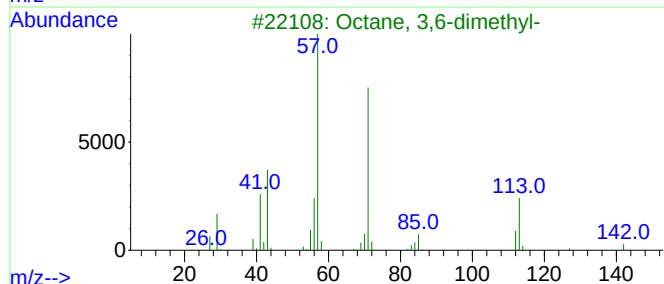
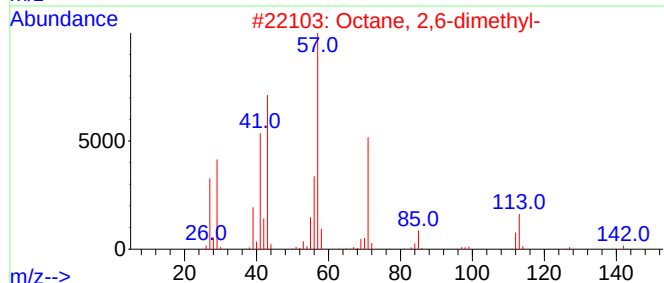
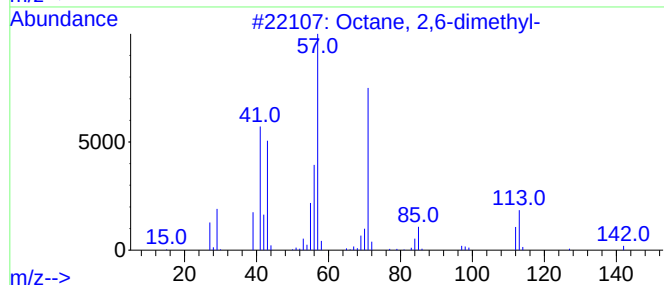
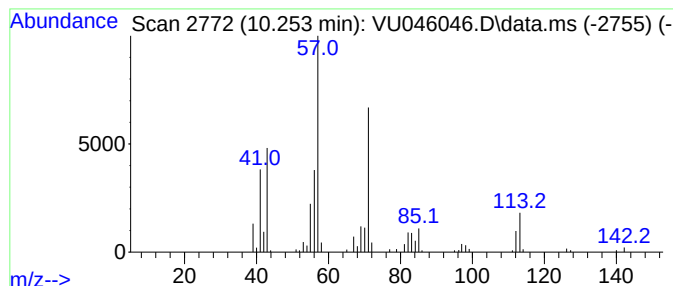
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.253	10.79 ug/L	136695	Chlorobenzene-d5	9.427

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	91
2			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	87
3			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	80
4			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	80
5			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	72



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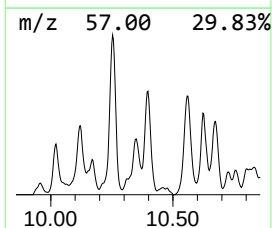
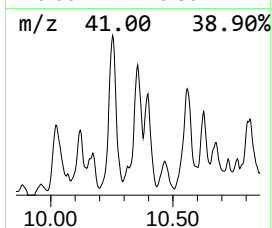
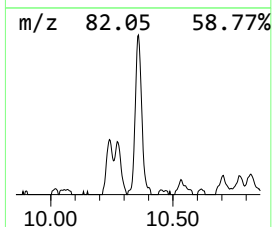
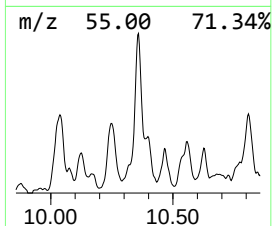
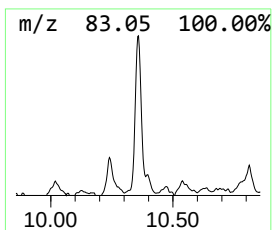
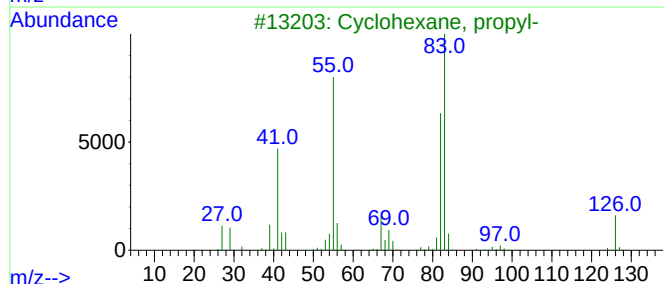
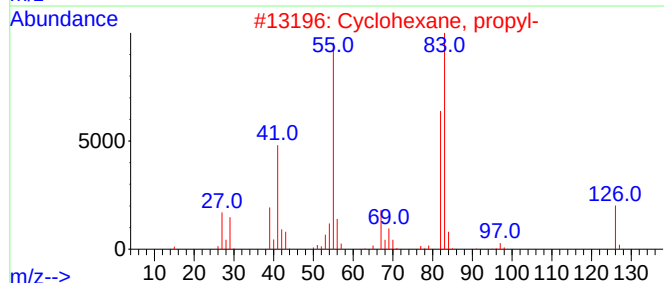
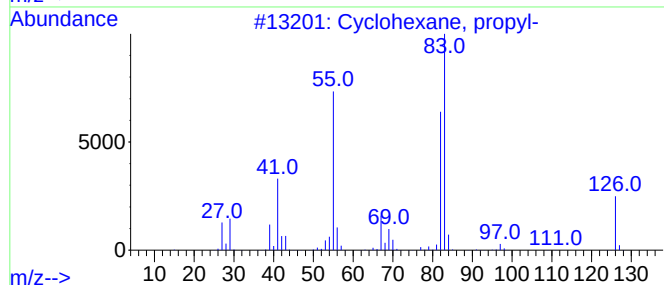
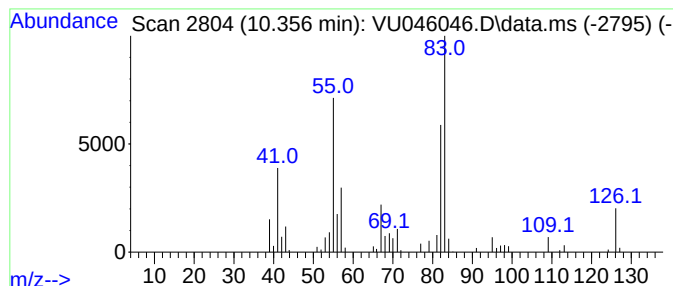
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Quant Title : VOC Analysis

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

Peak Number 4 (DEL) Alkane: Cyclic10.356 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.356	7.11 ug/L	90088	Chlorobenzene-d5	9.427

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, propyl-	126	C9H18	001678-92-8	93
2			Cyclohexane, propyl-	126	C9H18	001678-92-8	87
3			Cyclohexane, propyl-	126	C9H18	001678-92-8	83
4			Cyclohexane, propyl-	126	C9H18	001678-92-8	81
5			Cyclohexane, propyl-	126	C9H18	001678-92-8	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120221\
Data File : VU046046.D
Acq On : 02 Dec 2021 12:50
Operator : SY/MD
Sample : M4868-07
Misc : 6.94g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGKP2

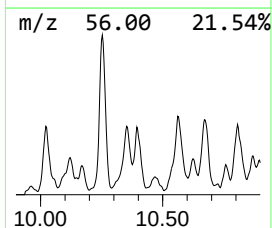
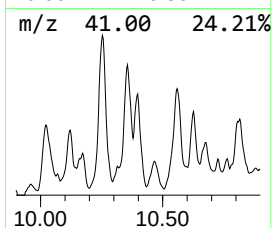
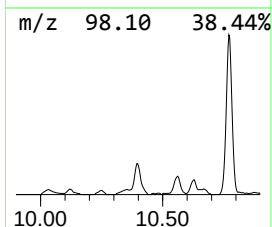
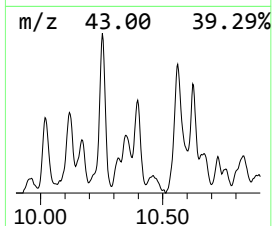
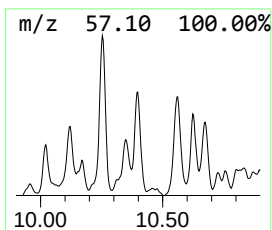
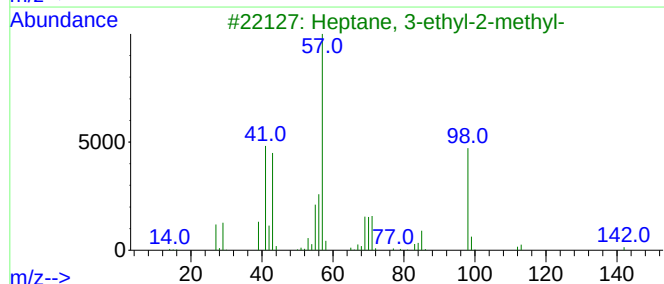
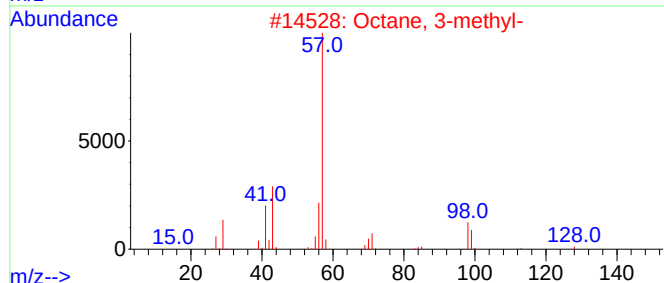
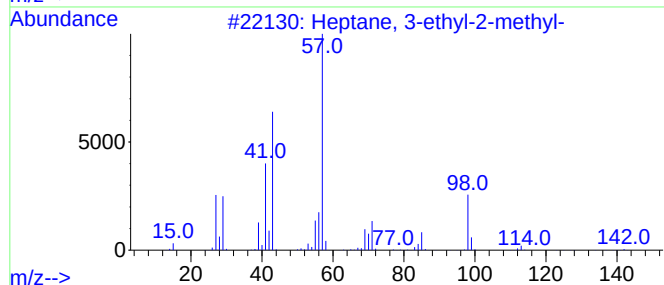
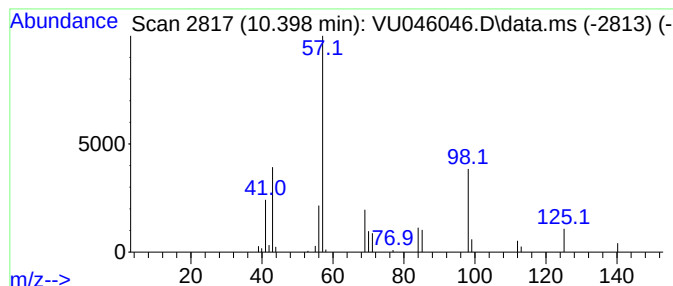
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Quant Title : VOC Analysis

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.398	4.02 ug/L	50880	Chlorobenzene-d5	9.427

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	45
2			Octane, 3-methyl-	128	C9H20	002216-33-3	40
3			Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	36
4			Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	33
5			Hexadecane, 1-chloro-	260	C16H33Cl	004860-03-1	25



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120221\
Data File : VU046046.D
Acq On : 02 Dec 2021 12:50
Operator : SY/MD
Sample : M4868-07
Misc : 6.94g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGKP2

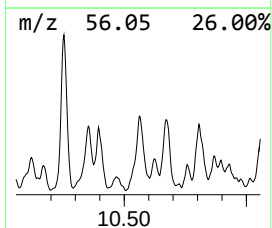
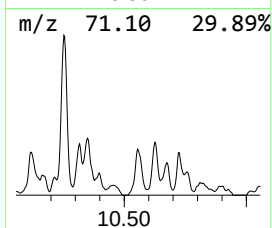
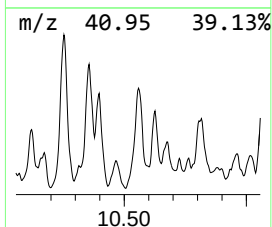
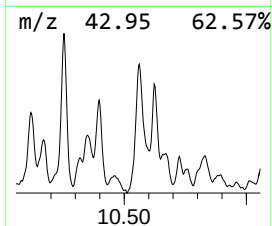
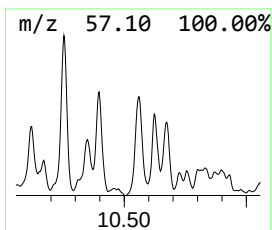
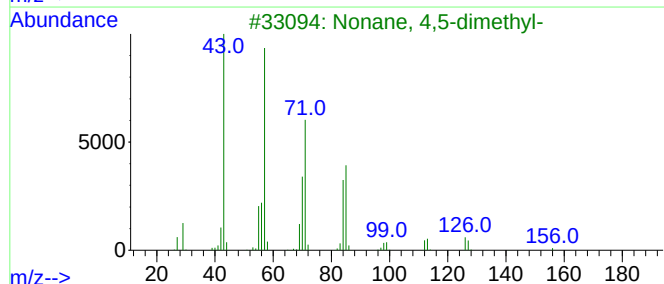
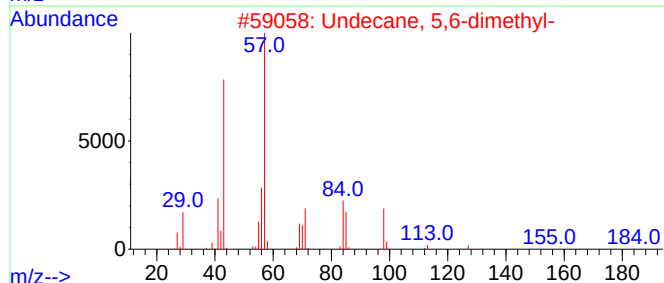
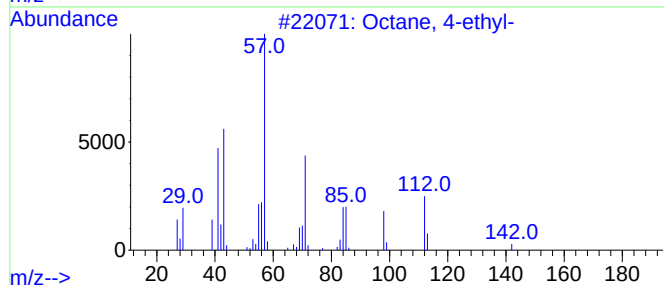
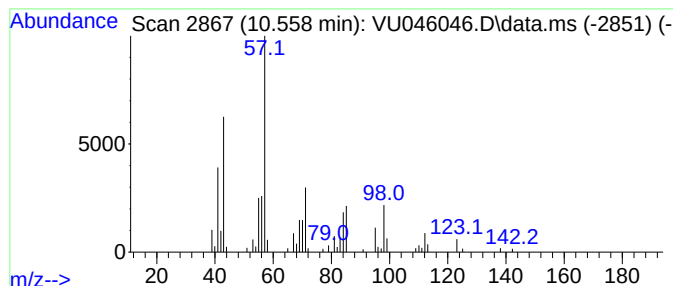
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Quant Title : VOC Analysis

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.558	8.17 ug/L	103412	Chlorobenzene-d5	9.427

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 4-ethyl-			142	C10H22	015869-86-0	47
2	Undecane, 5,6-dimethyl-			184	C13H28	017615-91-7	47
3	Nonane, 4,5-dimethyl-			156	C11H24	017302-23-7	43
4	Silane, trichlorooctadecyl-			386	C18H37Cl3Si	000112-04-9	43
5	Decane			142	C10H22	000124-18-5	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120221\
Data File : VU046046.D
Acq On : 02 Dec 2021 12:50
Operator : SY/MD
Sample : M4868-07
Misc : 6.94g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 6 Sample Multiplier: 1

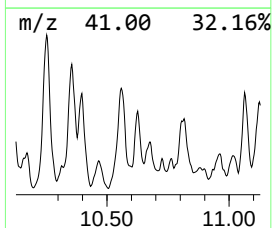
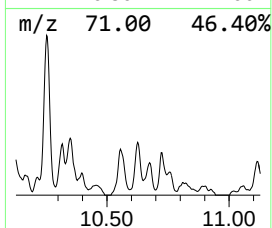
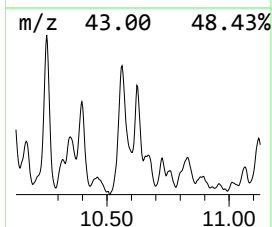
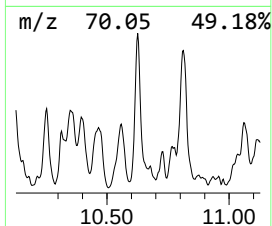
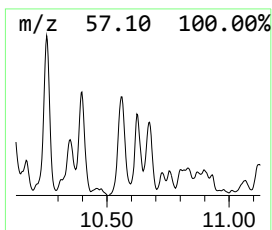
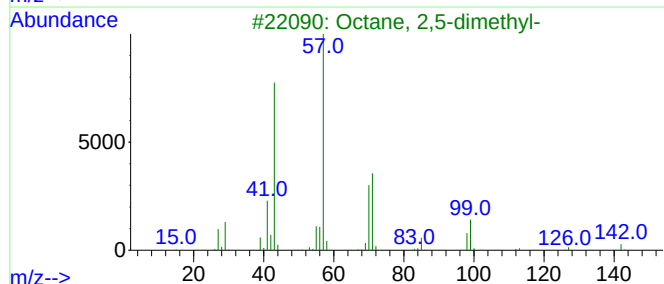
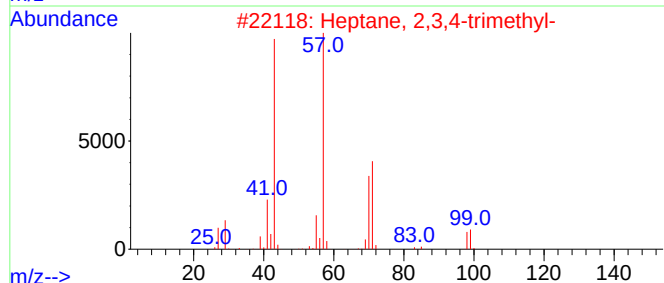
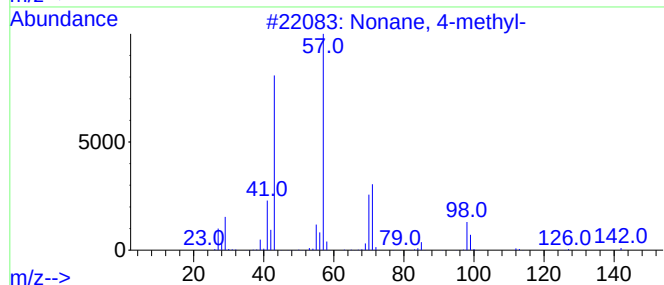
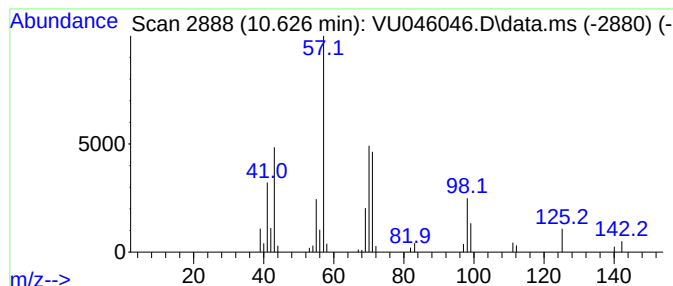
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ClientSampleId :
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Quant Title : VOC Analysis

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD			R.T.
10.626	2.64 ug/L	34962	1,4-Dichlorobenzene-d4			11.819
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonane, 4-methyl-		142	C10H22	017301-94-9	72
2	Heptane, 2,3,4-trimethyl-		142	C10H22	052896-95-4	59
3	Octane, 2,5-dimethyl-		142	C10H22	015869-89-3	50
4	Undecane, 4,5-dimethyl-		184	C13H28	017312-79-7	50
5	Nonane, 4-methyl-		142	C10H22	017301-94-9	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120221\
Data File : VU046046.D
Acq On : 02 Dec 2021 12:50
Operator : SY/MD
Sample : M4868-07
Misc : 6.94g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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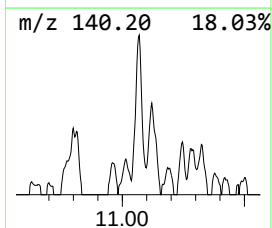
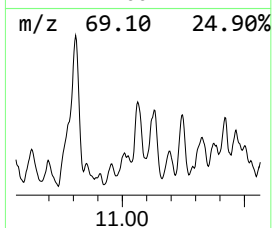
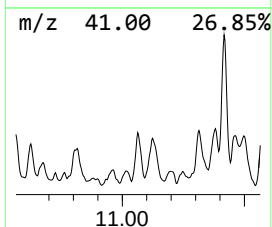
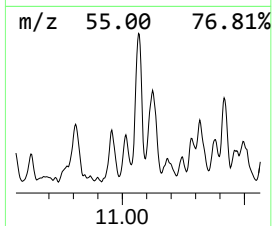
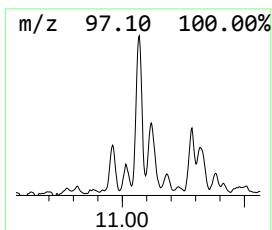
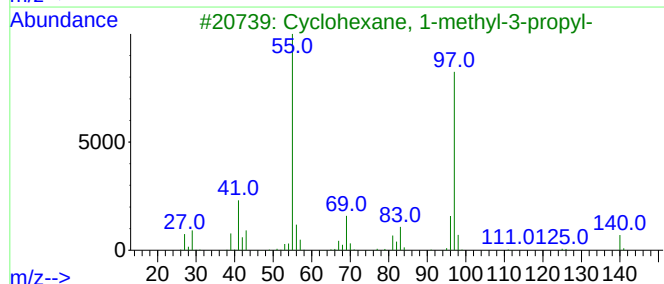
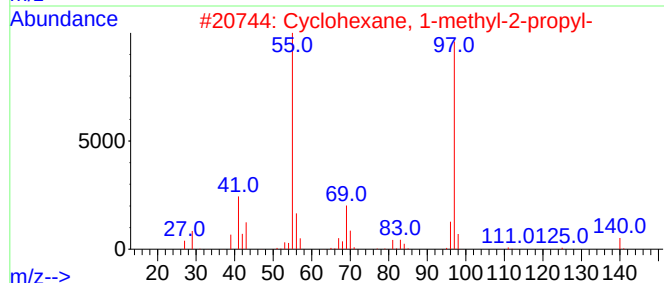
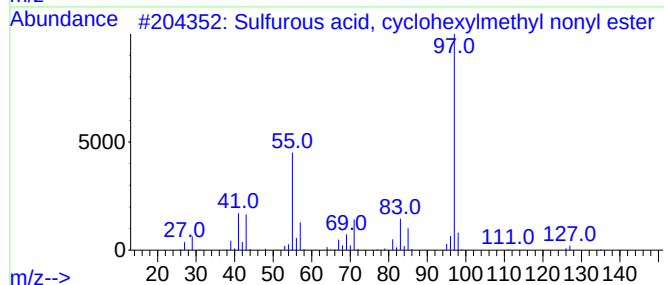
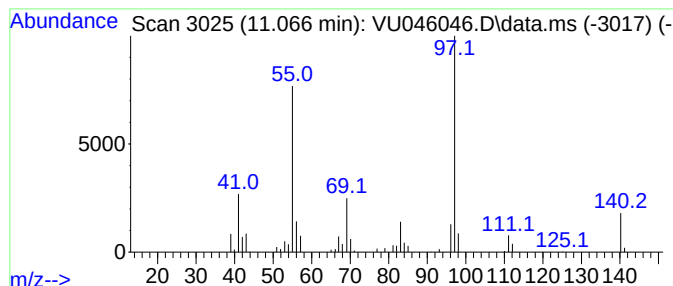
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Quant Title : VOC Analysis

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

Peak Number 8 Sulfurous acid, cyclohexylm... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.066	4.08 ug/L	54016	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, cyclohexylmethyl...	304	C16H32O3S	1000309-21-8	72
2			Cyclohexane, 1-methyl-2-propyl-	140	C10H20	004291-79-6	72
3			Cyclohexane, 1-methyl-3-propyl-	140	C10H20	004291-80-9	72
4			Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-77-7	72
5			Sulfurous acid, cyclohexylmethyl...	346	C19H38O3S	1000309-22-0	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120221\
Data File : VU046046.D
Acq On : 02 Dec 2021 12:50
Operator : SY/MD
Sample : M4868-07
Misc : 6.94g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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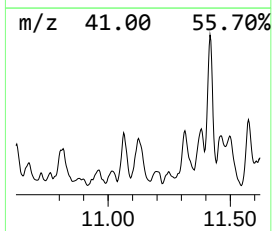
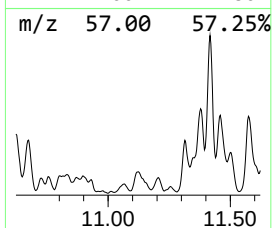
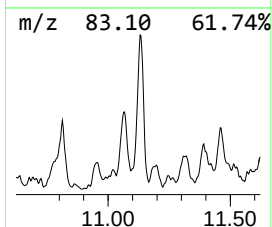
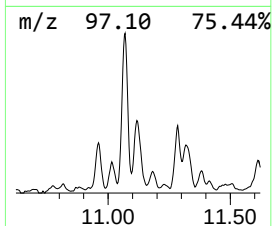
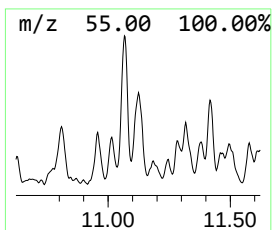
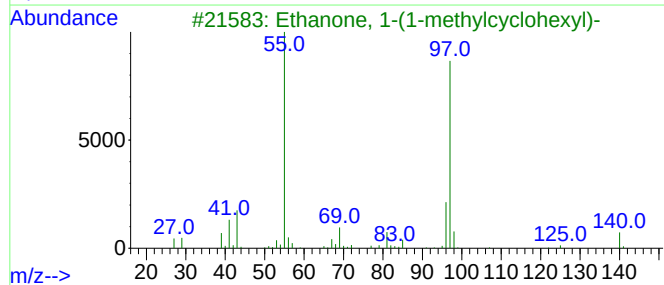
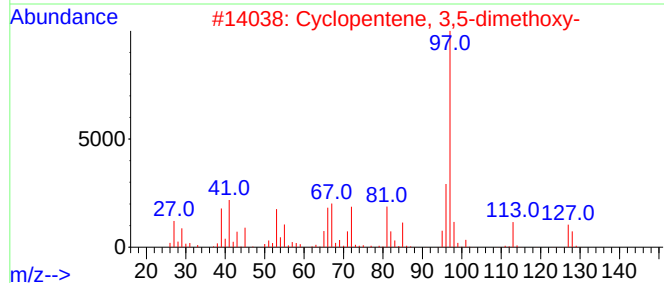
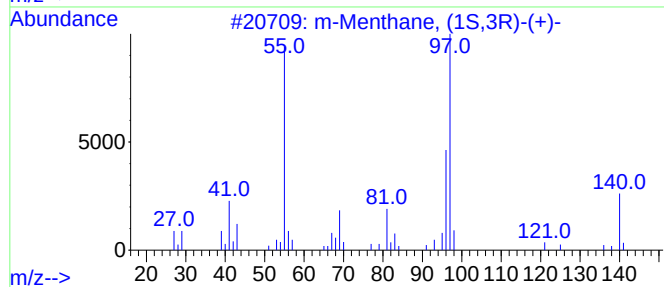
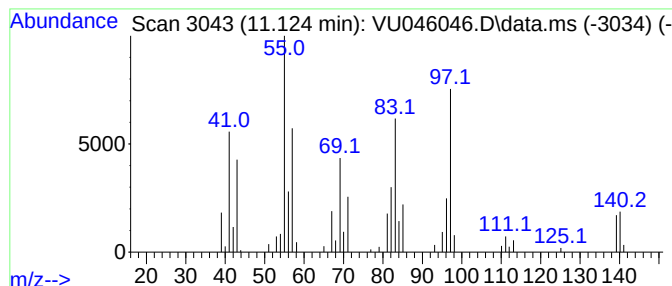
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Quant Title : VOC Analysis

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.124	5.73 ug/L	75793	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			m-Menthane, (1S,3R)-(+)-	140	C10H20	013837-66-6	46
2			Cyclopentene, 3,5-dimethoxy-	128	C7H12O2	089897-05-2	35
3			Ethanone, 1-(1-methylcyclohexyl)-	140	C9H16O	002890-62-2	35
4			Cyclohexane, 1-methyl-3-(1-methy...	140	C10H20	016580-24-8	35
5			Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	006069-98-3	30



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120221\
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Operator : SY/MD
Sample : M4868-07
Misc : 6.94g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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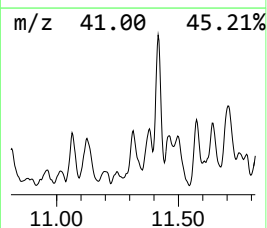
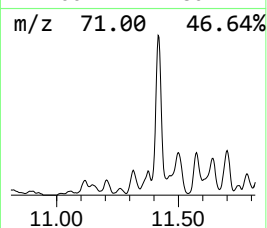
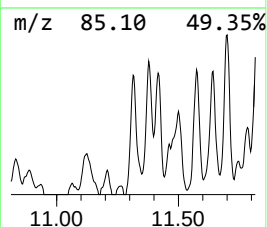
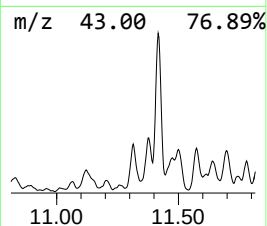
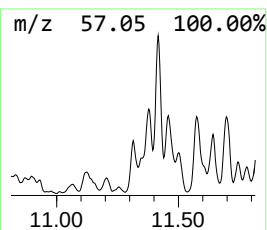
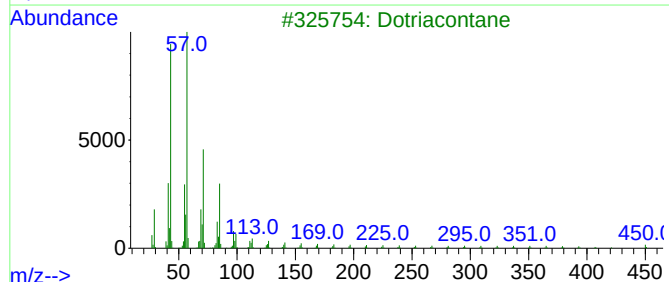
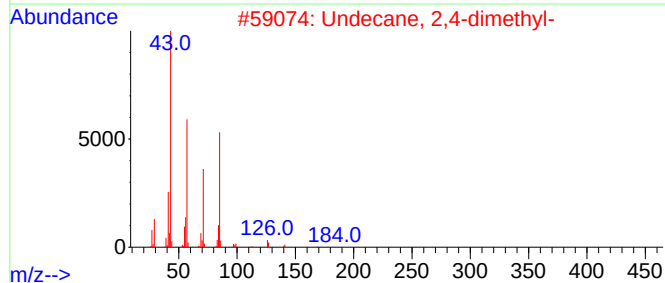
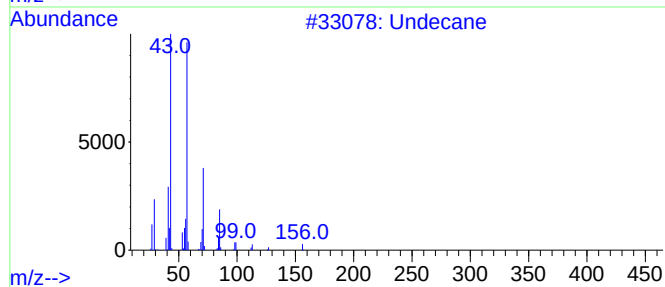
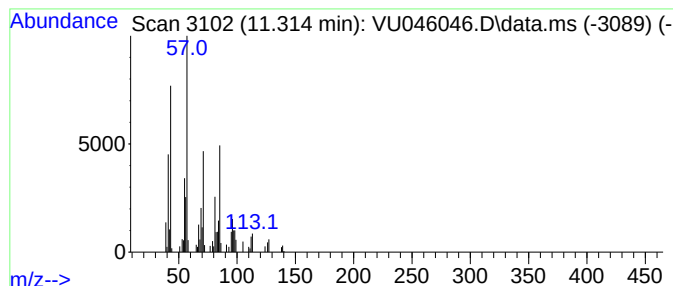
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Quant Title : VOC Analysis

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.314	5.59 ug/L	73982	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane	156	C11H24	001120-21-4	50
2			Undecane, 2,4-dimethyl-	184	C13H28	017312-80-0	50
3			Dotriacontane	451	C32H66	000544-85-4	47
4			Oxalic acid, monoamide, n-propyl...	341	C20H39NO3	1000309-25-8	47
5			Sulfurous acid, 2-ethylhexyl iso...	278	C14H30O3S	1000309-19-0	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120221\
Data File : VU046046.D
Acq On : 02 Dec 2021 12:50
Operator : SY/MD
Sample : M4868-07
Misc : 6.94g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGKP2

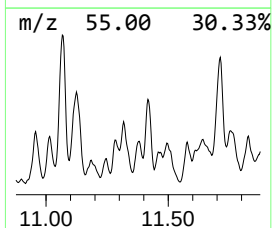
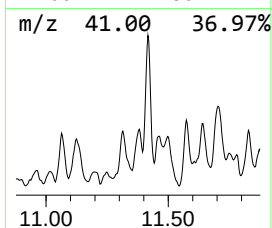
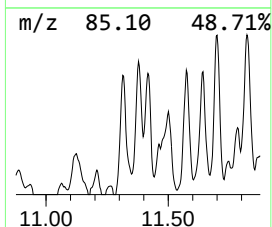
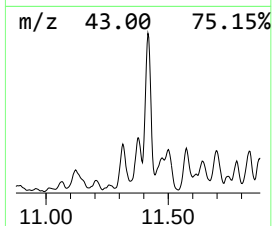
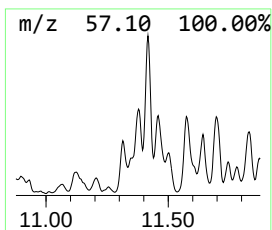
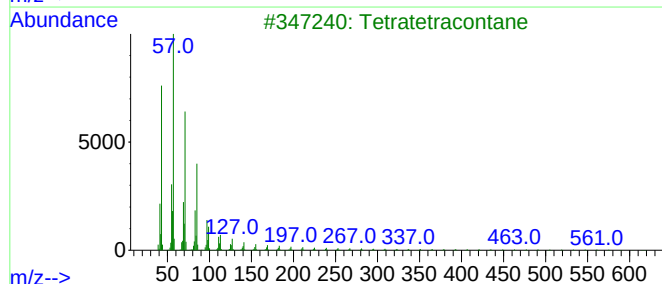
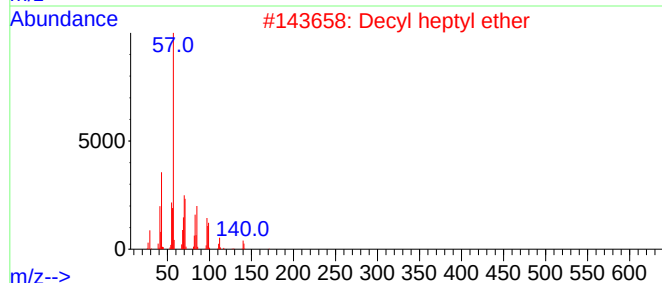
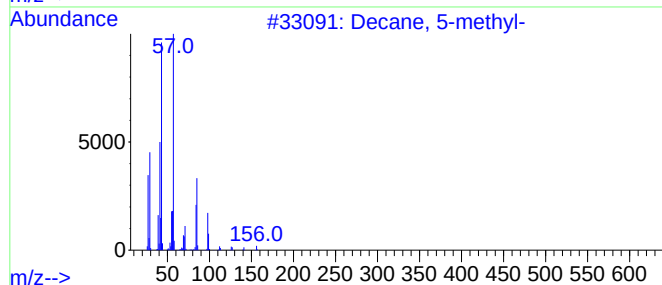
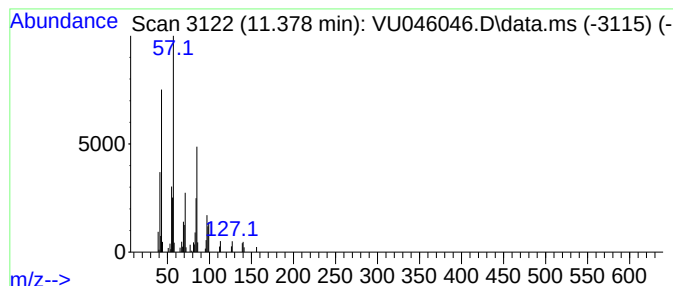
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Quant Title : VOC Analysis

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.378	3.16 ug/L	41770	1,4-Dichlorobenzene-d4	11.819

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Decane, 5-methyl-	156	C11H24	013151-35-4	64
2		Decyl heptyl ether	256	C17H36O	1000406-39-1	53
3		Tetratetracontane	619	C44H90	007098-22-8	53
4		Heptane, 2,2,3,3,5,6,6-heptamethyl-	198	C14H30	007225-67-4	50
5		Heptyl tetradecyl ether	312	C21H44O	1000406-39-3	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120221\
Data File : VU046046.D
Acq On : 02 Dec 2021 12:50
Operator : SY/MD
Sample : M4868-07
Misc : 6.94g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGKP2

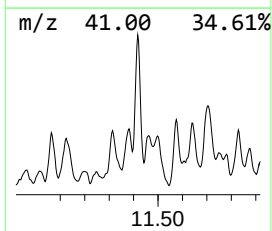
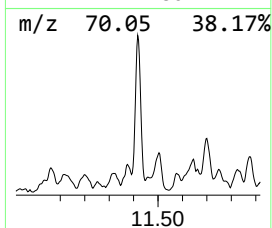
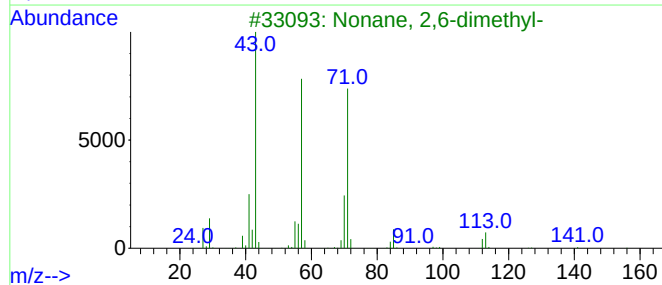
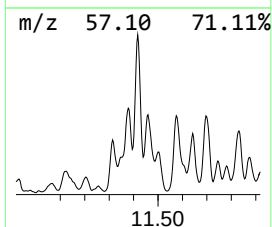
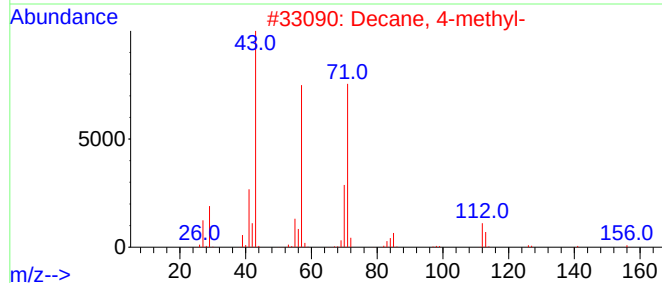
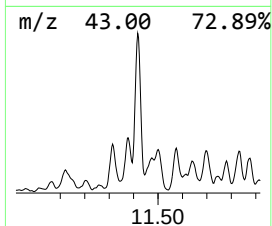
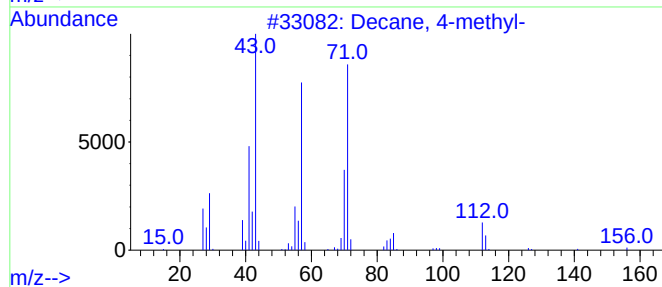
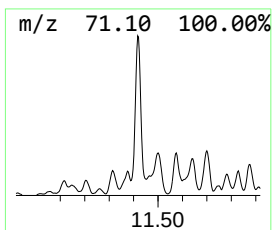
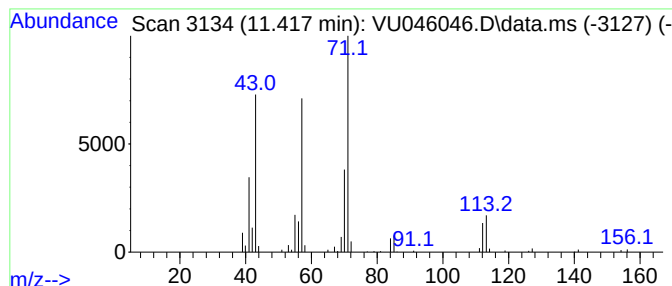
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Quant Title : VOC Analysis

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.417	9.15 ug/L	121086	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 4-methyl-	156	C11H24	002847-72-5	91
2			Decane, 4-methyl-	156	C11H24	002847-72-5	91
3			Nonane, 2,6-dimethyl-	156	C11H24	017302-28-2	91
4			Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	78
5			4-Methyl-3-heptanol, pentafluoro...	276	C11H17F5O2	1000365-51-7	78



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120221\
Data File : VU046046.D
Acq On : 02 Dec 2021 12:50
Operator : SY/MD
Sample : M4868-07
Misc : 6.94g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGKP2

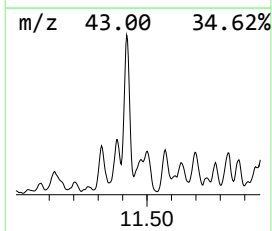
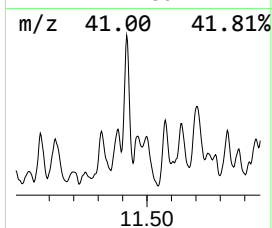
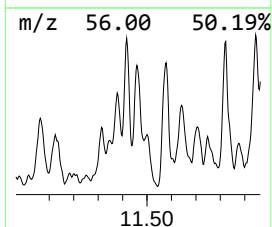
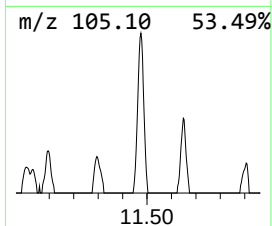
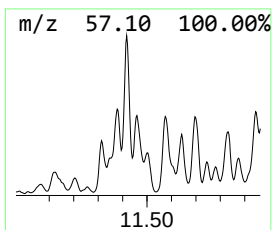
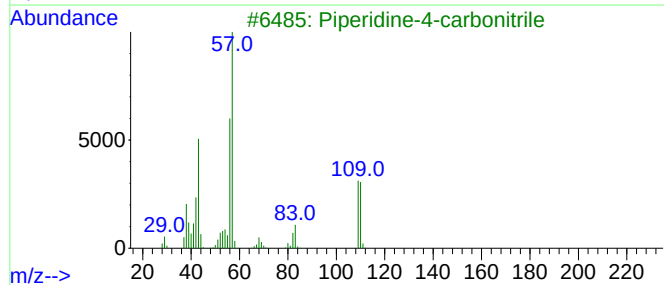
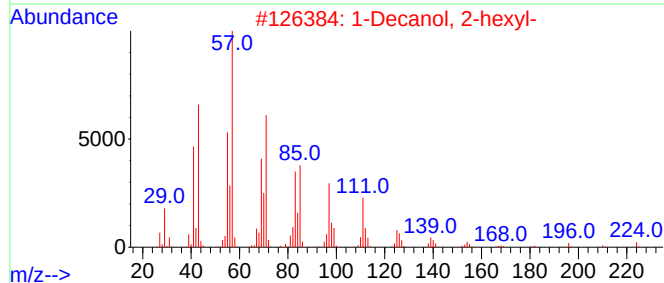
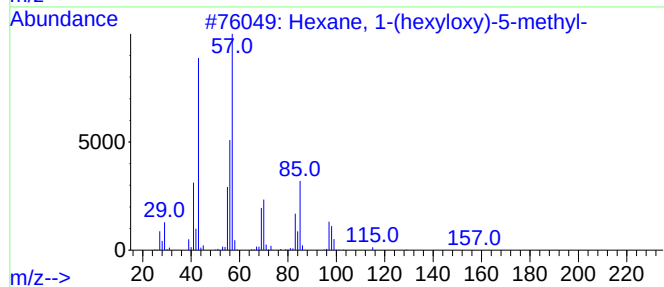
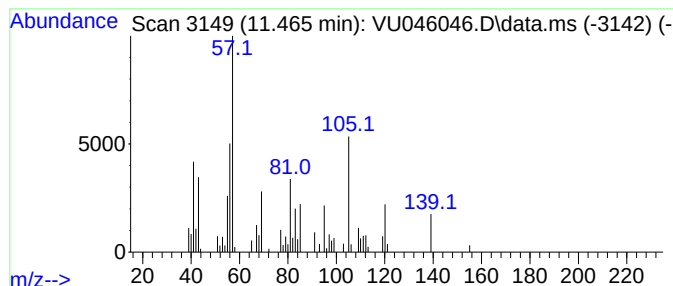
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Quant Title : VOC Analysis

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

Peak Number 13 unknown-01 Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.465	3.14 ug/L	41491	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 1-(hexyloxy)-5-methyl-	200	C13H28O	074421-19-5	38
2			1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	38
3			Piperidine-4-carbonitrile	110	C6H10N2	004395-98-6	38
4			1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	32
5			1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	22



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120221\
Data File : VU046046.D
Acq On : 02 Dec 2021 12:50
Operator : SY/MD
Sample : M4868-07
Misc : 6.94g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGKP2

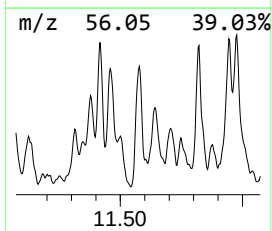
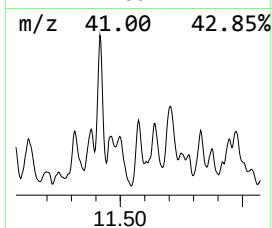
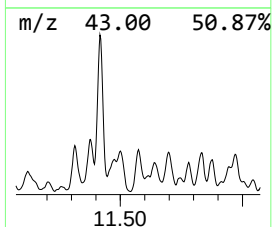
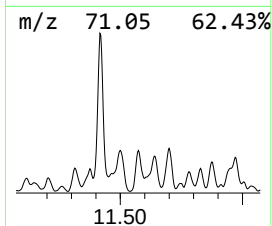
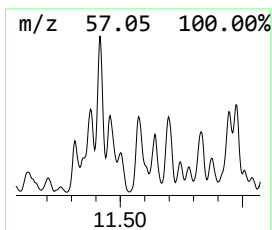
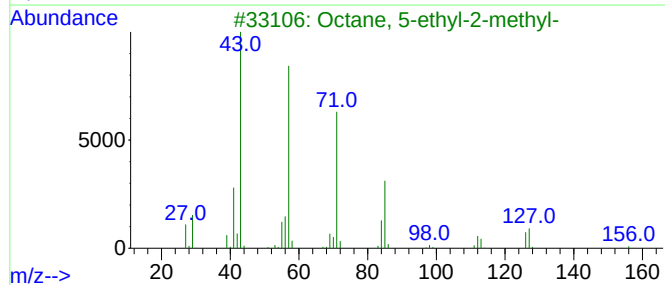
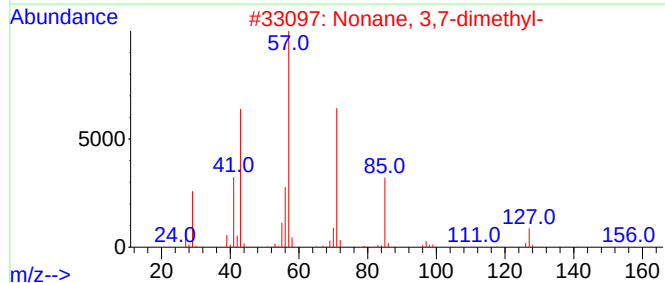
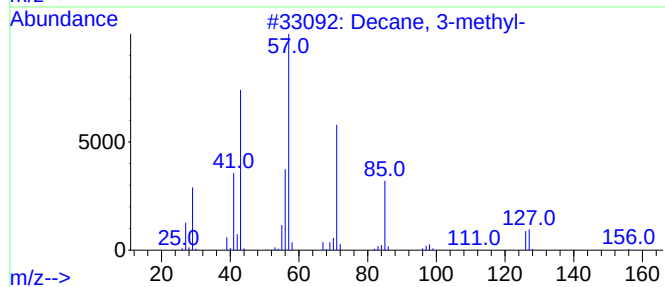
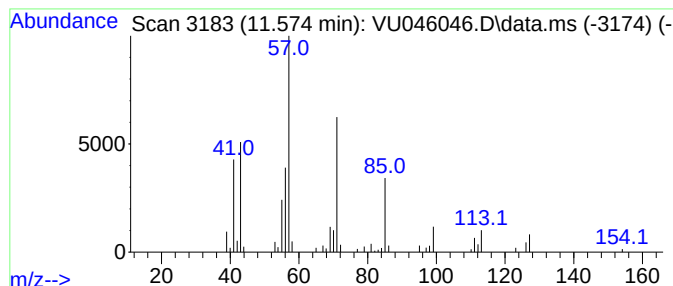
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Quant Title : VOC Analysis

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

Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.574	5.34 ug/L	70615	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 3-methyl-	156	C11H24	013151-34-3	64
2			Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	50
3			Octane, 5-ethyl-2-methyl-	156	C11H24	062016-18-6	50
4			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	50
5			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120221\
Data File : VU046046.D
Acq On : 02 Dec 2021 12:50
Operator : SY/MD
Sample : M4868-07
Misc : 6.94g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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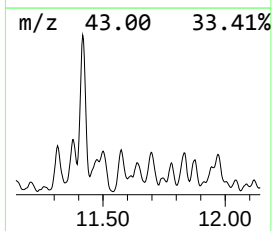
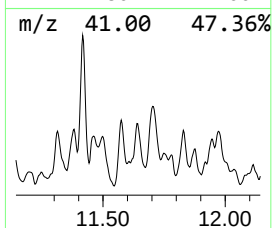
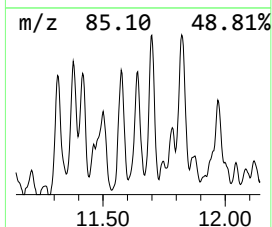
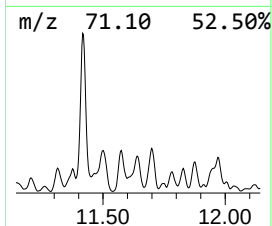
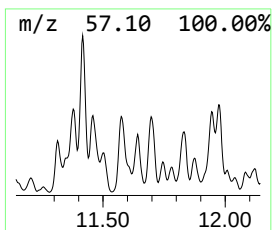
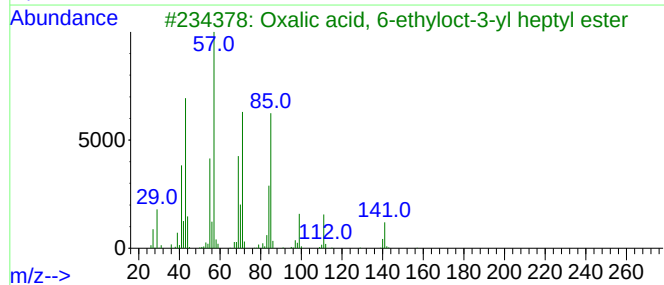
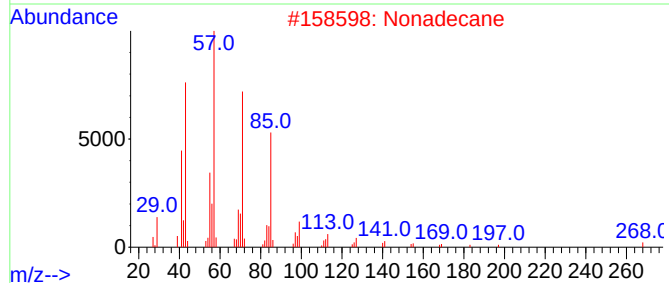
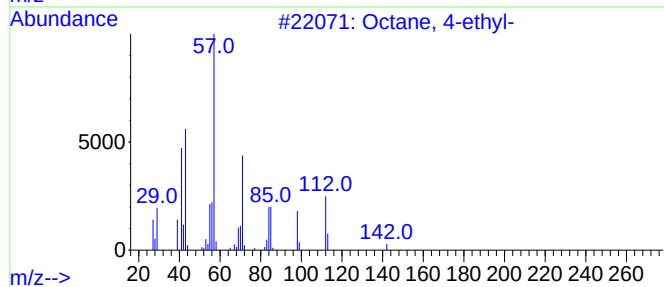
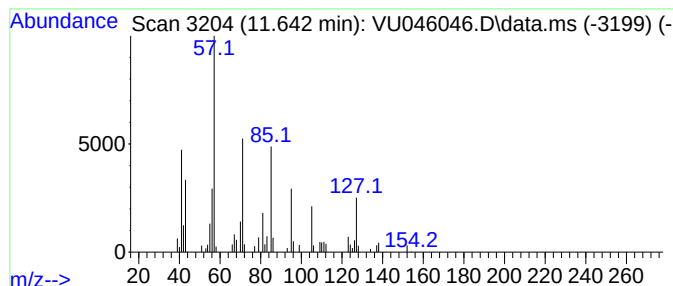
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Quant Title : VOC Analysis

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

Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.642	3.17 ug/L	41949	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 4-ethyl-	142	C10H22	015869-86-0	38
2			Nonadecane	268	C19H40	000629-92-5	37
3			Oxalic acid, 6-ethyloct-3-yl hep...	328	C19H36O4	1000309-34-5	37
4			Nonadecane	268	C19H40	000629-92-5	37
5			Sulfurous acid, 2-ethylhexyl iso...	278	C14H30O3S	1000309-19-0	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120221\
Data File : VU046046.D
Acq On : 02 Dec 2021 12:50
Operator : SY/MD
Sample : M4868-07
Misc : 6.94g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGKP2

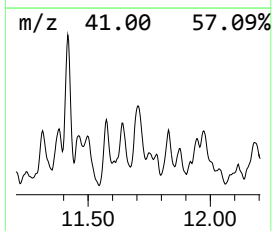
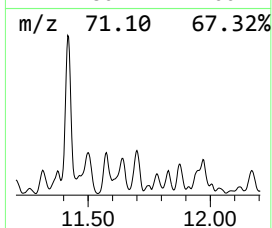
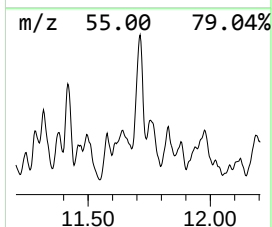
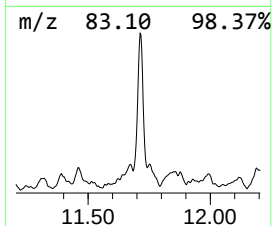
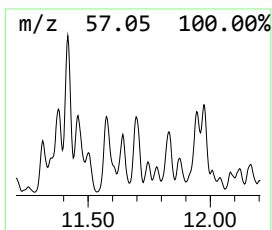
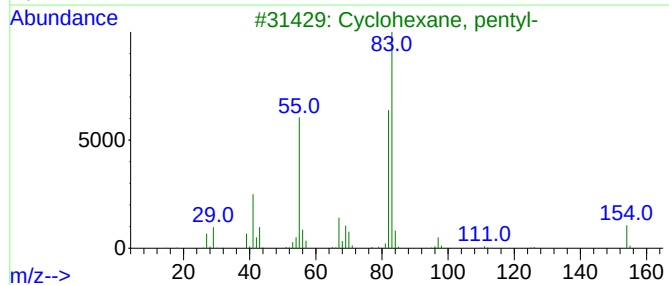
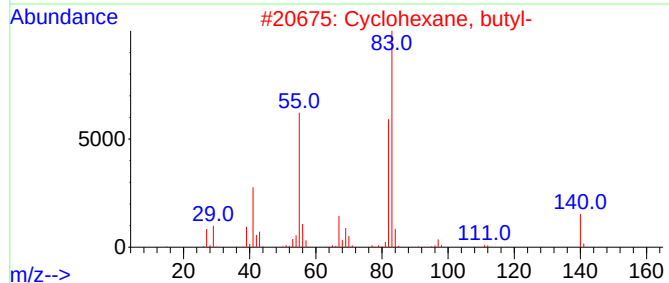
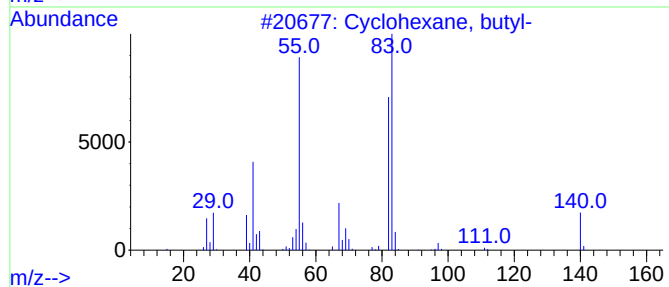
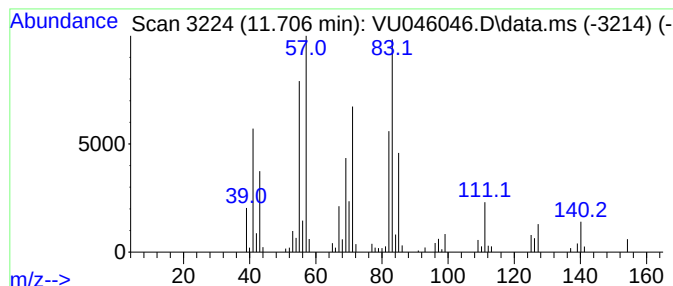
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Quant Title : VOC Analysis

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

Peak Number 16 (DEL) Alkane: Cyclic11.706 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.706	7.29 ug/L	96430	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, butyl-	140	C10H20	001678-93-9	50
2			Cyclohexane, butyl-	140	C10H20	001678-93-9	46
3			Cyclohexane, pentyl-	154	C11H22	004292-92-6	45
4			Cyclohexane, pentyl-	154	C11H22	004292-92-6	41
5			Cyclohexane, octyl-	196	C14H28	001795-15-9	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120221\
Data File : VU046046.D
Acq On : 02 Dec 2021 12:50
Operator : SY/MD
Sample : M4868-07
Misc : 6.94g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGKP2

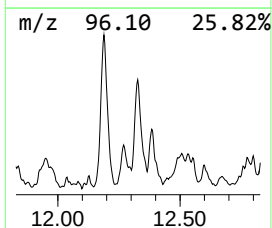
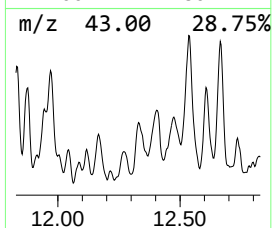
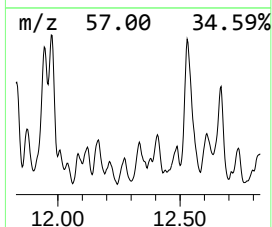
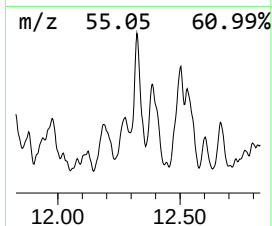
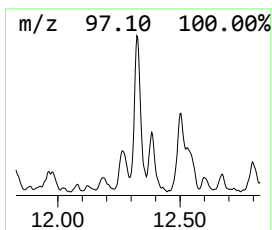
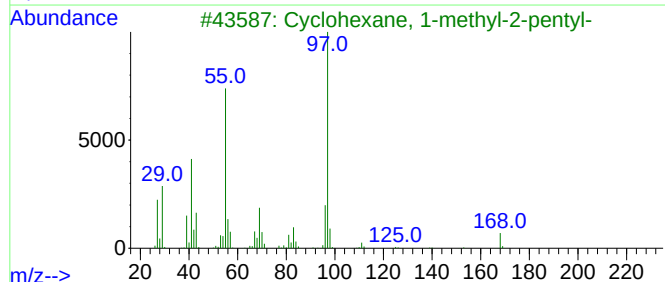
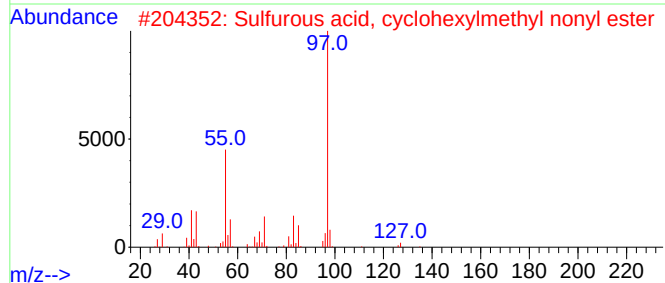
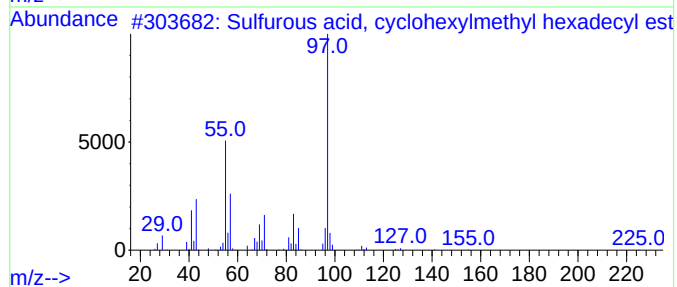
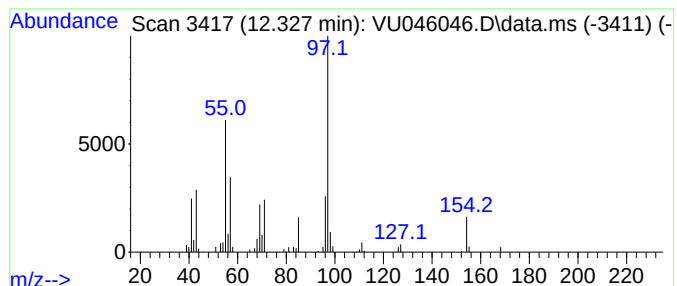
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Quant Title : VOC Analysis

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

Peak Number 17 Sulfurous acid, cyclohexylm... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.327	3.41 ug/L	45117	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, cyclohexylmethyl...	402	C23H46O3S	1000309-22-4	64
2			Sulfurous acid, cyclohexylmethyl...	304	C16H32O3S	1000309-21-8	59
3			Cyclohexane, 1-methyl-2-pentyl-	168	C12H24	054411-01-7	53
4			Cyclohexane, 1-methyl-2-pentyl-	168	C12H24	054411-01-7	53
5			Sulfurous acid, cyclohexylmethyl...	416	C24H48O3S	1000309-22-5	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120221\
Data File : VU046046.D
Acq On : 02 Dec 2021 12:50
Operator : SY/MD
Sample : M4868-07
Misc : 6.94g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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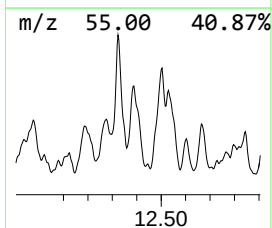
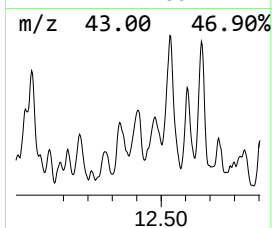
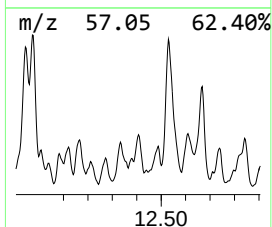
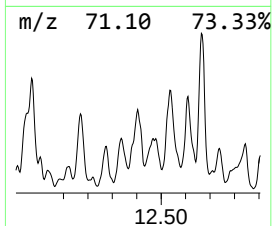
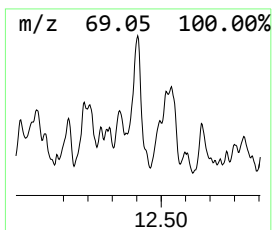
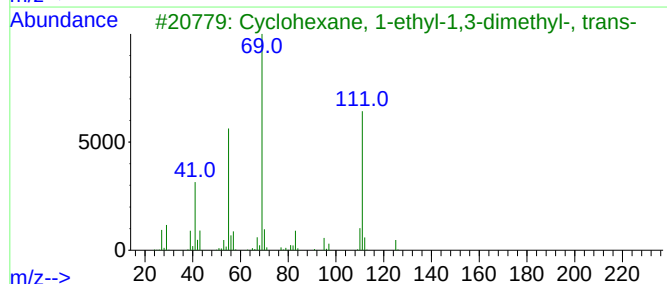
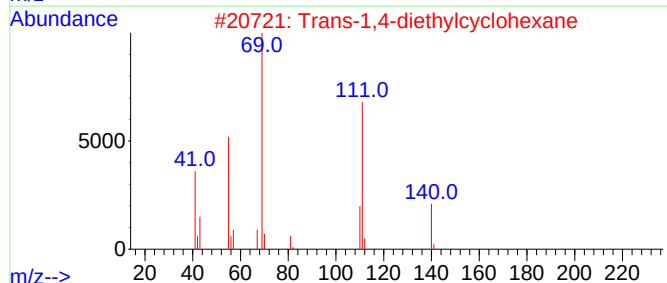
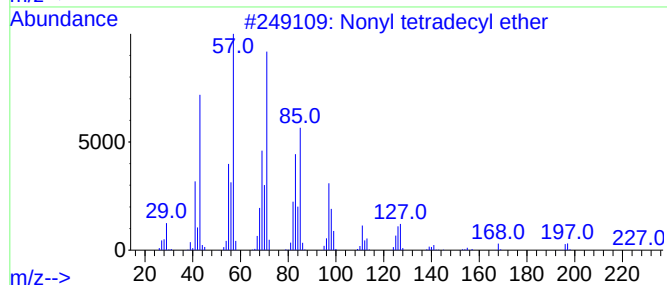
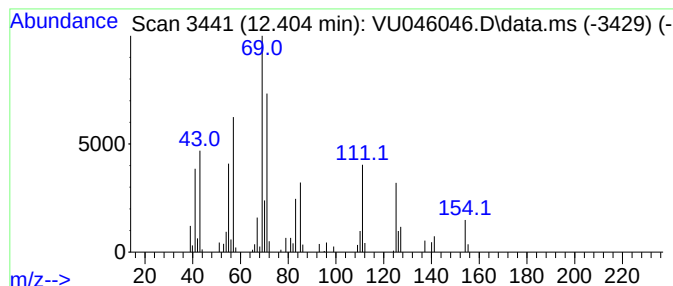
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Quant Title : VOC Analysis

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

Peak Number 18 unknown-02 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.404	5.55 ug/L	73462	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonyl tetradecyl ether	340	C23H48O	1000406-37-6	38
2			Trans-1,4-diethylcyclohexane	140	C10H20	013990-93-7	38
3			Cyclohexane, 1-ethyl-1,3-dimethy...	140	C10H20	062238-29-3	35
4			Terpinen-4-ol	154	C10H18O	000562-74-3	35
5			3-Cyclohexen-1-ol, 4-methyl-1-(1...	154	C10H18O	020126-76-5	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120221\
Data File : VU046046.D
Acq On : 02 Dec 2021 12:50
Operator : SY/MD
Sample : M4868-07
Misc : 6.94g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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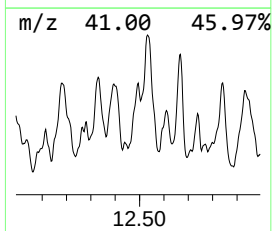
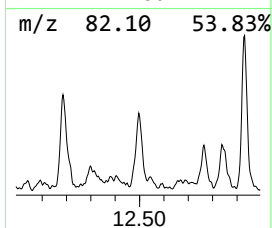
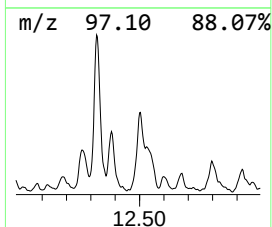
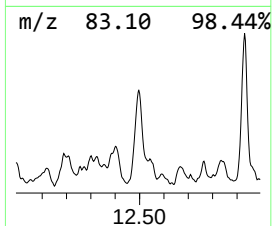
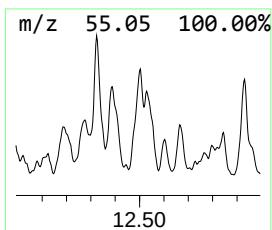
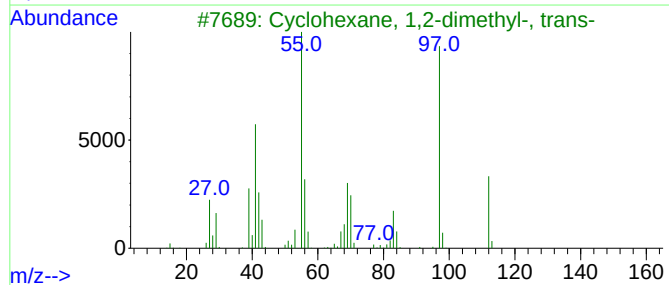
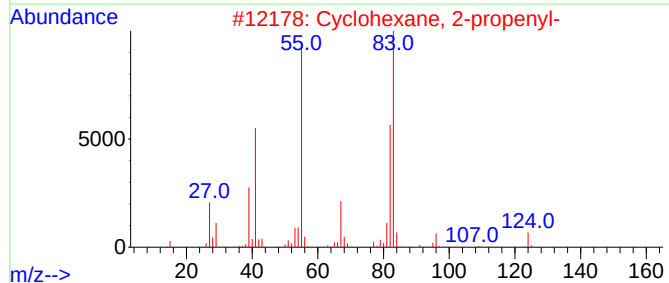
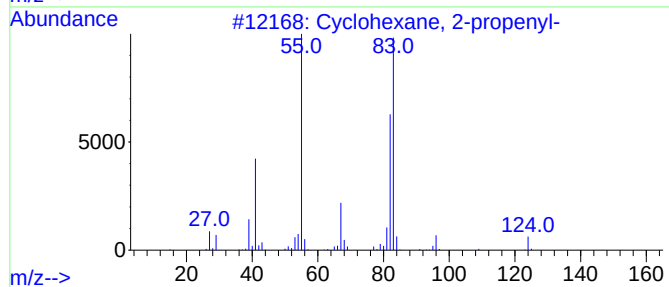
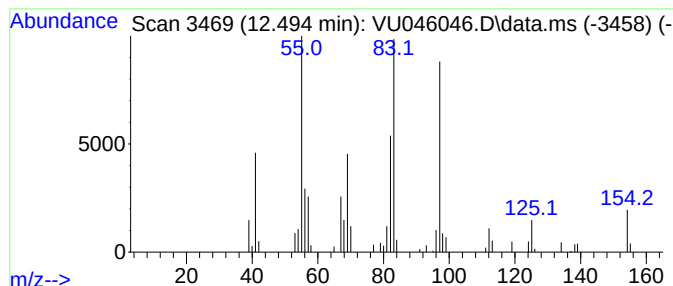
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Quant Title : VOC Analysis

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

Peak Number 19 Cyclohexane, 2-propenyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.494	3.84 ug/L	50824	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 2-propenyl-	124	C9H16	002114-42-3	55
2			Cyclohexane, 2-propenyl-	124	C9H16	002114-42-3	45
3			Cyclohexane, 1,2-dimethyl-, trans-	112	C8H16	006876-23-9	43
4			(2-Methylbutyl)cyclohexane	154	C11H22	054105-77-0	43
5			Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120221\
Data File : VU046046.D
Acq On : 02 Dec 2021 12:50
Operator : SY/MD
Sample : M4868-07
Misc : 6.94g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGKP2

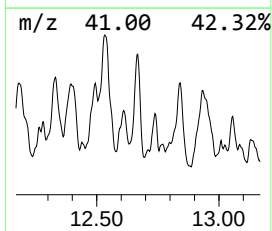
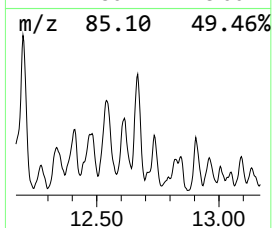
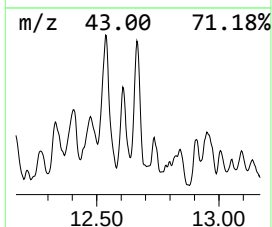
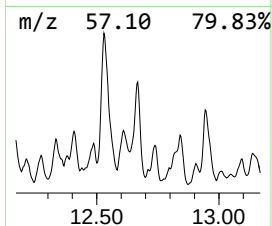
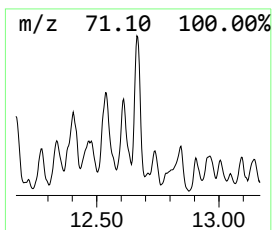
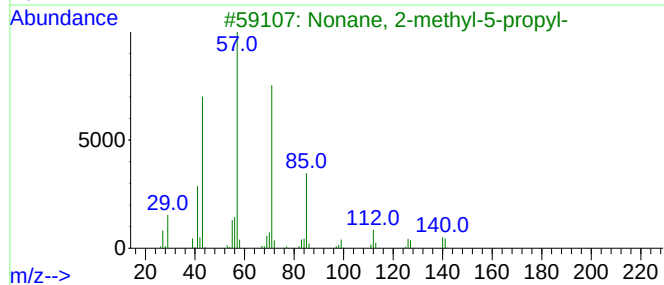
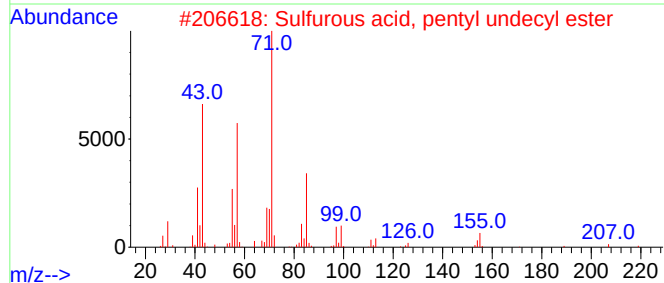
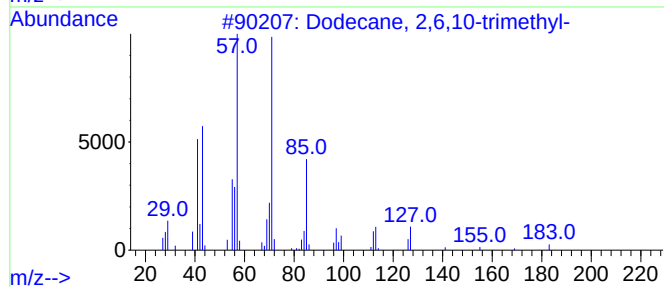
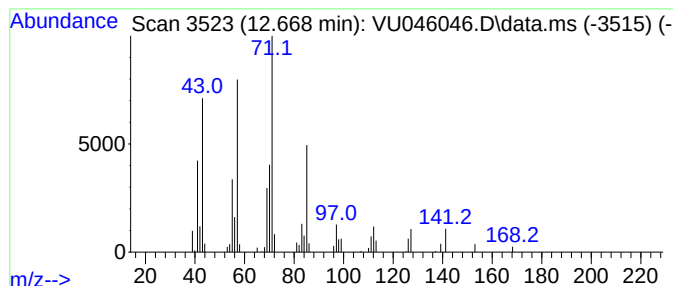
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Quant Title : VOC Analysis

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

Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.668	4.08 ug/L	53995	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	64
2			Sulfurous acid, pentyl undecyl e...	306	C16H34O3S	1000309-14-4	53
3			Nonane, 2-methyl-5-propyl-	184	C13H28	031081-17-1	53
4			Nonane, 3-methyl-5-propyl-	184	C13H28	031081-18-2	53
5			Heptacosane	380	C27H56	000593-49-7	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120221\
Data File : VU046046.D
Acq On : 02 Dec 2021 12:50
Operator : SY/MD
Sample : M4868-07
Misc : 6.94g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGKP2

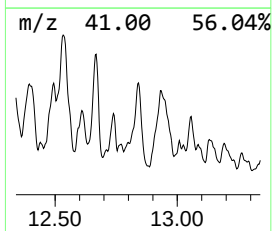
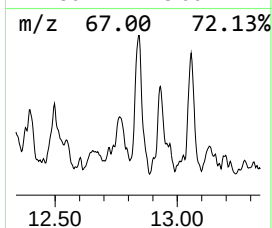
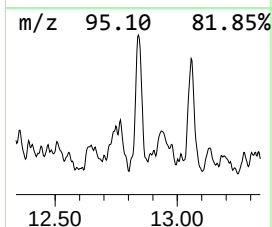
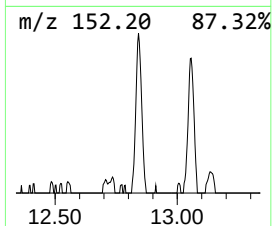
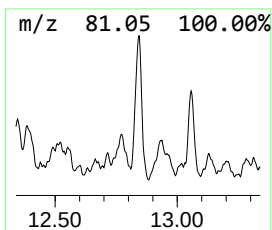
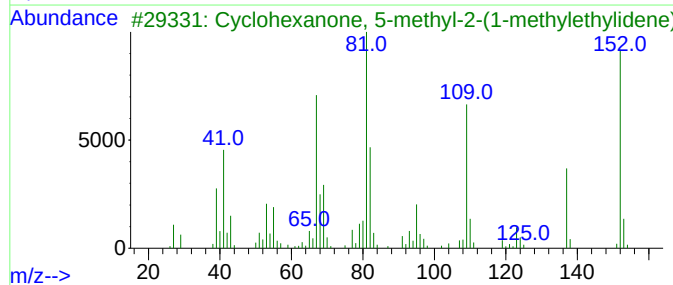
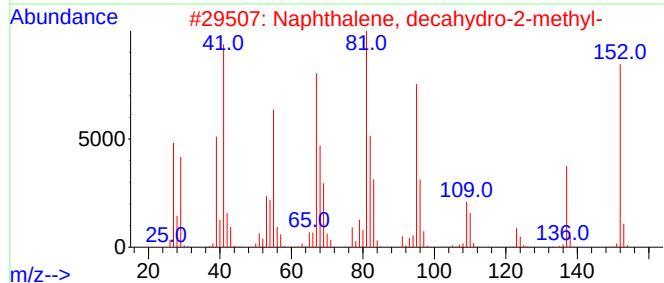
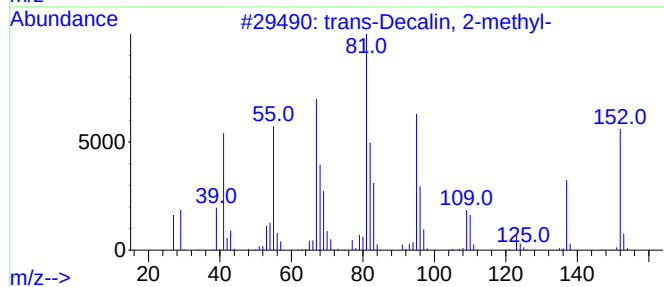
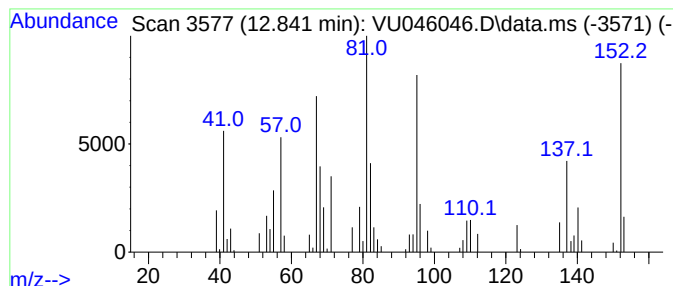
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Quant Title : VOC Analysis

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

Peak Number 21 trans-Decalin, 2-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.841	5.50 ug/L	72745	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	91
2			Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	90
3			Cyclohexanone, 5-methyl-2-(1-met...	152	C10H16O	015932-80-6	86
4			Cyclodecene, 1-methyl-	152	C11H20	066633-38-3	72
5			Cyclohexanone, 2-methyl-5-(1-met...	152	C10H16O	005948-04-9	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120221\
Data File : VU046046.D
Acq On : 02 Dec 2021 12:50
Operator : SY/MD
Sample : M4868-07
Misc : 6.94g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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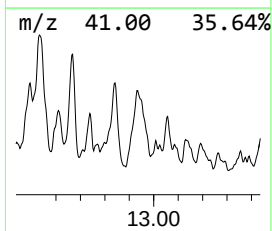
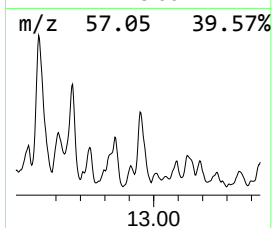
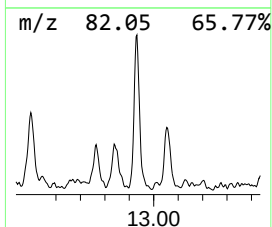
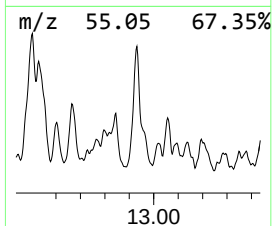
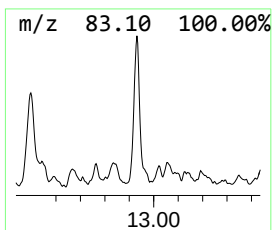
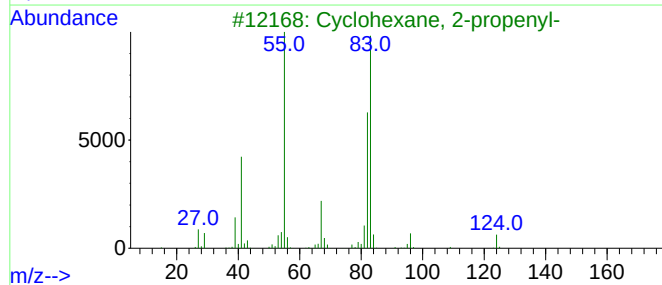
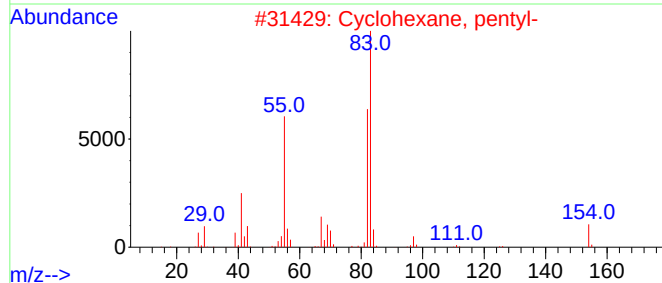
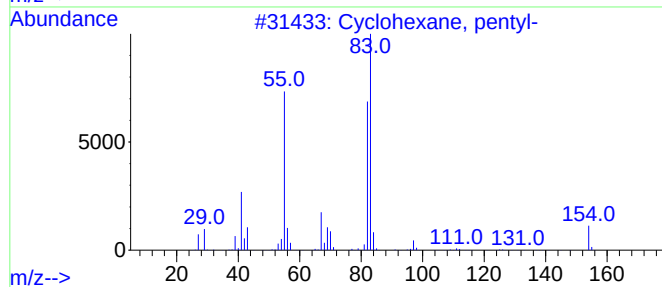
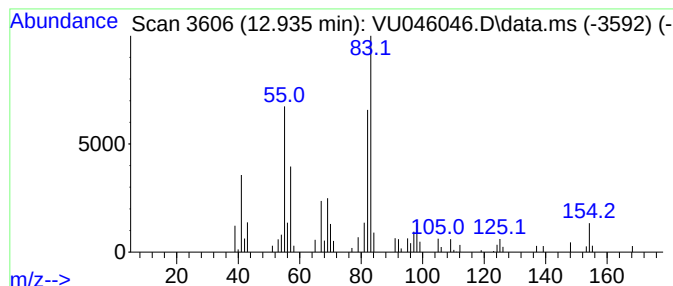
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Quant Title : VOC Analysis

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

Peak Number 22 (DEL) Alkane: Cyclic12.935 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.935	8.25 ug/L	109094	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, pentyl-	154	C11H22	004292-92-6	87
2			Cyclohexane, pentyl-	154	C11H22	004292-92-6	81
3			Cyclohexane, 2-propenyl-	124	C9H16	002114-42-3	72
4			Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	68
5			Cyclohexane, (4-methylpentyl)-	168	C12H24	061142-20-9	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120221\
Data File : VU046046.D
Acq On : 02 Dec 2021 12:50
Operator : SY/MD
Sample : M4868-07
Misc : 6.94g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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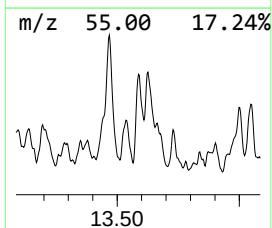
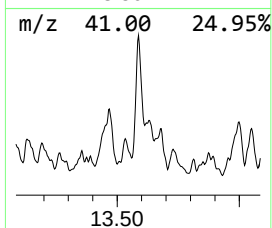
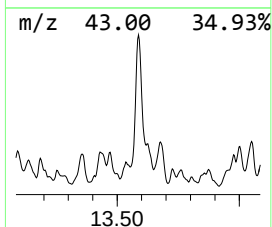
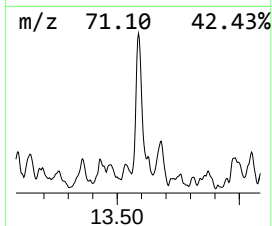
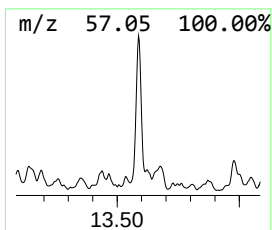
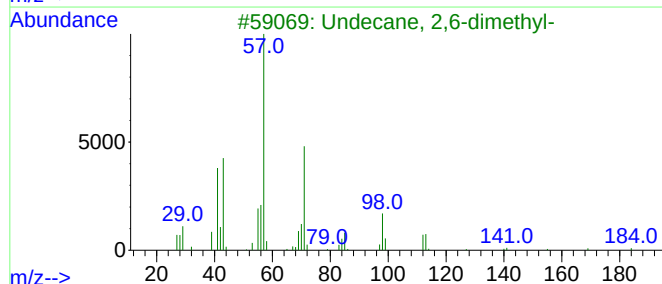
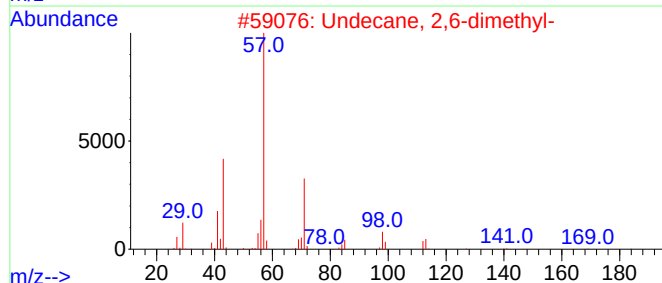
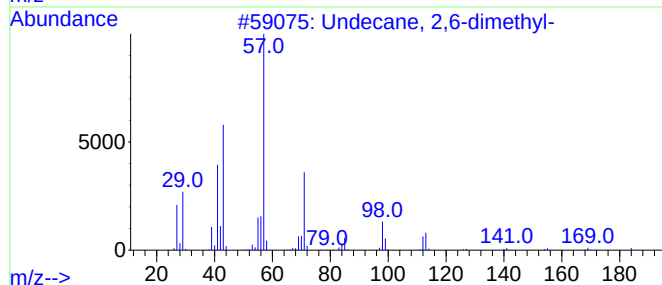
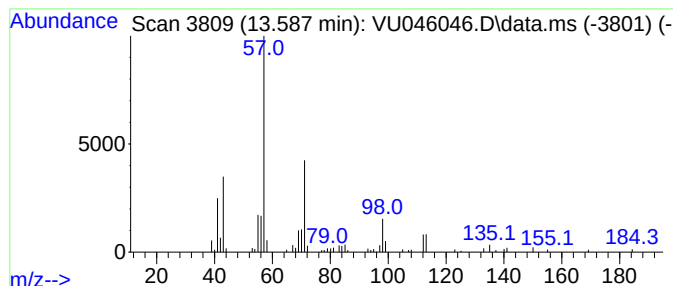
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Quant Title : VOC Analysis

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

Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.587	5.77 ug/L	76384	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	90
2			Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	90
3			Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	87
4			Dodecane, 6-methyl-	184	C13H28	006044-71-9	83
5			Undecane, 2,5-dimethyl-	184	C13H28	017301-22-3	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120221\
Data File : VU046046.D
Acq On : 02 Dec 2021 12:50
Operator : SY/MD
Sample : M4868-07
Misc : 6.94g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGKP2

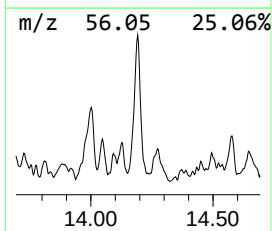
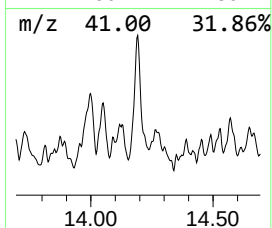
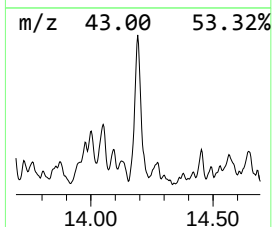
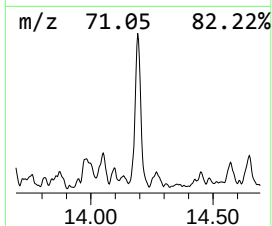
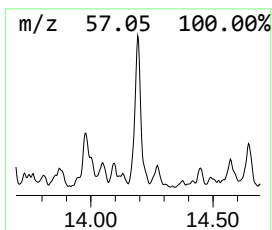
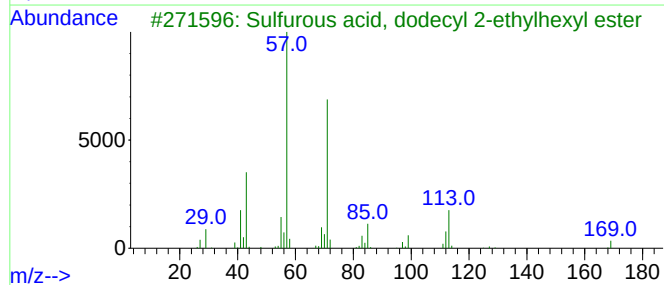
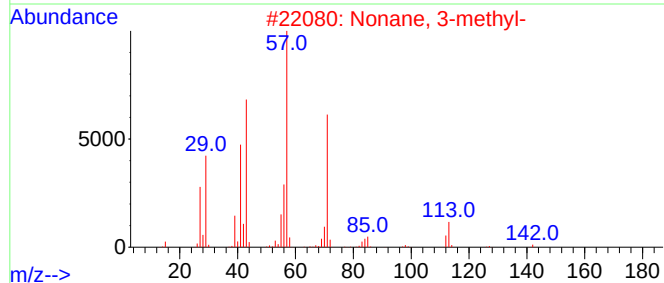
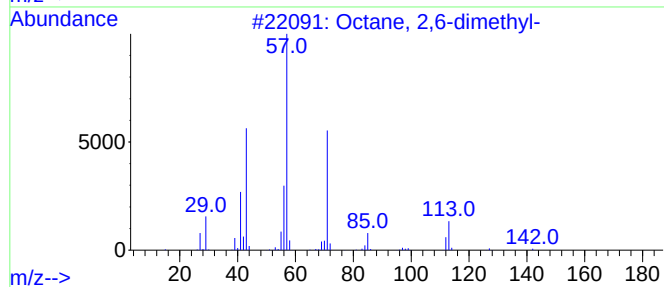
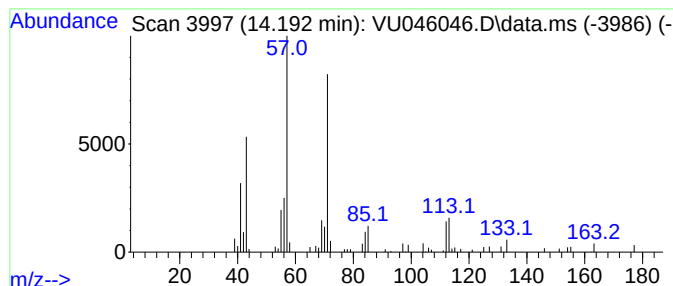
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

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

Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.192	4.86 ug/L	64254	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	78
2			Nonane, 3-methyl-	142	C10H22	005911-04-6	78
3			Sulfurous acid, dodecyl 2-ethylh...	362	C20H42O3S	1000309-19-5	72
4			Octane, 2,3,7-trimethyl-	156	C11H24	062016-34-6	72
5			Sulfurous acid, butyl octyl ester	250	C12H26O3S	1000309-17-5	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120221\
Data File : VU046046.D
Acq On : 02 Dec 2021 12:50
Operator : SY/MD
Sample : M4868-07
Misc : 6.94g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGKP2

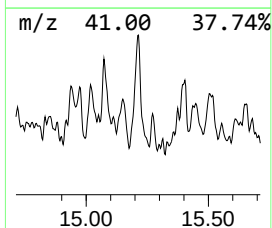
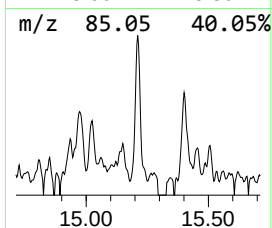
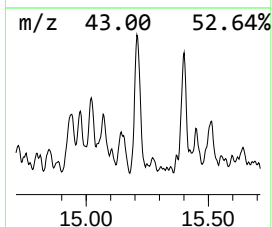
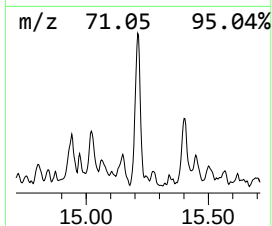
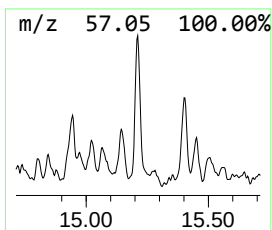
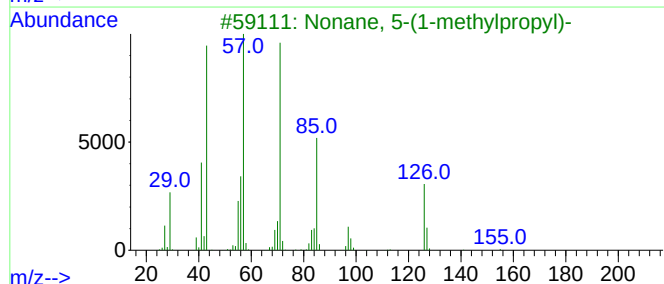
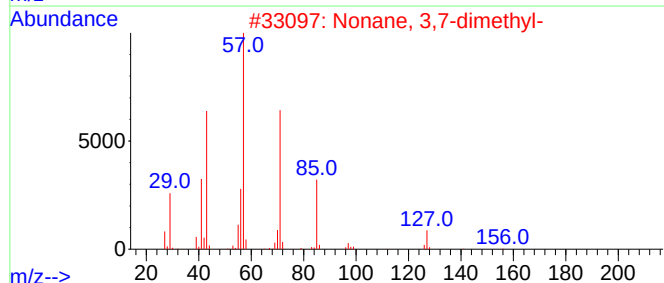
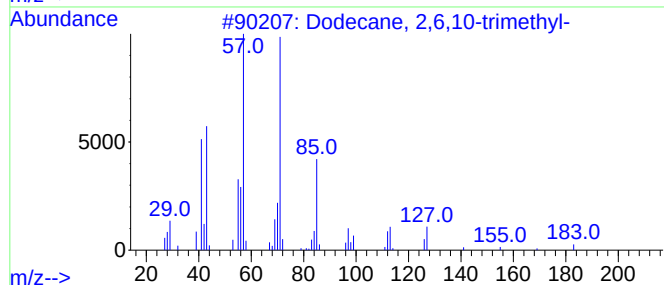
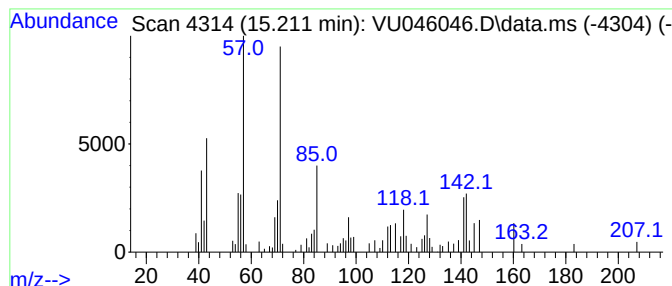
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

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

Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.211	3.80 ug/L	50223	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	60
2			Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	43
3			Nonane, 5-(1-methylpropyl)-	184	C13H28	062185-54-0	43
4			Hexadecyl octyl ether	354	C24H50O	1000406-38-6	38
5			Dodecyl octyl ether	298	C20H42O	1000406-38-4	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120221\
 Data File : VU046046.D
 Acq On : 02 Dec 2021 12:50
 Operator : SY/MD
 Sample : M4868-07
 Misc : 6.94g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BGKP2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
 Quant Title : VOC Analysis

TIC Library : C:\Database\NIST0.L
 TIC Integration Parameters: LSCINT.P

					--Internal Standard--			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
(DEL) Alkane: C...	10.025	5.8	ug/L	73808	2	9.427	633225	50.0
(DEL) Alkane: S...	10.253	10.8	ug/L	136695	2	9.427	633225	50.0
(DEL) Alkane: C...	10.356	7.1	ug/L	90088	2	9.427	633225	50.0
(DEL) Alkane: S...	10.398	4.0	ug/L	50880	2	9.427	633225	50.0
(DEL) Alkane: S...	10.558	8.2	ug/L	103412	2	9.427	633225	50.0
(DEL) Alkane: S...	10.626	2.6	ug/L	34962	3	11.819	661436	50.0
Sulfurous acid,...	11.066	4.1	ug/L	54016	3	11.819	661436	50.0
(DEL) Alkane: S...	11.124	5.7	ug/L	75793	3	11.819	661436	50.0
(DEL) Alkane: S...	11.314	5.6	ug/L	73982	3	11.819	661436	50.0
(DEL) Alkane: S...	11.378	3.2	ug/L	41770	3	11.819	661436	50.0
(DEL) Alkane: S...	11.417	9.2	ug/L	121086	3	11.819	661436	50.0
unknown-01	11.465	3.1	ug/L	41491	3	11.819	661436	50.0
(DEL) Alkane: S...	11.574	5.3	ug/L	70615	3	11.819	661436	50.0
(DEL) Alkane: S...	11.642	3.2	ug/L	41949	3	11.819	661436	50.0
(DEL) Alkane: C...	11.706	7.3	ug/L	96430	3	11.819	661436	50.0
Sulfurous acid,...	12.327	3.4	ug/L	45117	3	11.819	661436	50.0
unknown-02	12.404	5.5	ug/L	73462	3	11.819	661436	50.0
Cyclohexane, 2-...	12.494	3.8	ug/L	50824	3	11.819	661436	50.0
(DEL) Alkane: S...	12.668	4.1	ug/L	53995	3	11.819	661436	50.0
trans-Decalin, ...	12.841	5.5	ug/L	72745	3	11.819	661436	50.0
(DEL) Alkane: C...	12.935	8.3	ug/L	109094	3	11.819	661436	50.0
(DEL) Alkane: S...	13.587	5.8	ug/L	76384	3	11.819	661436	50.0
(DEL) Alkane: S...	14.192	4.9	ug/L	64254	3	11.819	661436	50.0
(DEL) Alkane: S...	15.211	3.8	ug/L	50223	3	11.819	661436	50.0