

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV010522\
 Data File : VV024282.D
 Acq On : 05 Jan 2022 16:56
 Operator : SY/MD
 Sample : N1025-10
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 BGLH4

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_V\Method\SFAMVLM123121WMA.M
 Title : VOC Analysis

Signal : TIC: VV024282.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.108	17	21	36	rVB2	4841	5977	0.09%	0.013%
2	1.310	74	84	96	rVB2	241375	267030	4.21%	0.572%
3	1.416	111	117	128	rVB	25989	28546	0.45%	0.061%
4	1.568	156	164	175	rBV	111429	123429	1.95%	0.264%
5	1.635	181	185	191	rVB5	4381	4645	0.07%	0.010%
6	1.709	201	208	218	rVB	19371	24164	0.38%	0.052%
7	1.950	273	283	300	rBV	1052762	1404227	22.15%	3.006%
8	2.108	321	332	343	rBV	241970	357412	5.64%	0.765%
9	2.172	343	352	381	rVB	111963	261485	4.12%	0.560%
10	2.346	386	406	407	rBV3	9486	23413	0.37%	0.050%
11	2.760	526	535	550	rVB	17104	31090	0.49%	0.067%
12	3.040	615	622	635	rVB5	2382	4518	0.07%	0.010%
13	3.188	651	668	690	rBV	59910	125743	1.98%	0.269%
14	3.760	834	846	860	rBV7	3831	9793	0.15%	0.021%
15	3.905	872	891	944	rBV2	997116	2624098	41.39%	5.616%
16	4.346	1009	1028	1054	rBV	156050	393771	6.21%	0.843%
17	4.677	1117	1131	1144	rVB5	4628	11390	0.18%	0.024%
18	4.773	1148	1161	1163	rBV4	2334	3563	0.06%	0.008%
19	5.043	1227	1245	1252	rBV2	346275	861316	13.59%	1.843%
20	5.095	1253	1261	1293	rVB	859529	1860570	29.35%	3.982%
21	5.522	1378	1394	1410	rBV3	26292	61697	0.97%	0.132%
22	5.612	1410	1422	1452	rVB	292307	587700	9.27%	1.258%
23	5.915	1507	1516	1535	rVB7	11948	26741	0.42%	0.057%
24	6.066	1550	1563	1591	rBV	182189	392753	6.20%	0.841%
25	6.185	1591	1600	1601	rBV3	7262	8699	0.14%	0.019%
26	6.982	1837	1848	1870	rBV	114259	206517	3.26%	0.442%
27	7.227	1915	1924	1937	rBV3	19043	35169	0.55%	0.075%
28	7.313	1939	1951	1961	rVV	425496	730466	11.52%	1.563%
29	7.384	1961	1973	1997	rVB	3812119	6339364	100.00%	13.568%
30	7.616	2036	2045	2063	rBV	83038	138477	2.18%	0.296%
31	7.702	2065	2072	2083	rVB4	3723	6148	0.10%	0.013%
32	7.937	2136	2145	2152	rBV4	5681	9896	0.16%	0.021%
33	8.005	2161	2166	2179	rVB4	6264	10545	0.17%	0.023%
34	8.082	2179	2190	2206	rBV2	244575	424406	6.69%	0.908%
35	8.722	2382	2389	2397	rVB6	2224	3211	0.05%	0.007%
36	8.850	2415	2429	2460	rBV2	447557	907271	14.31%	1.942%

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Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM123121WMA.M
 Title : VOC Analysis

37	9.008	2465	2478	2503	rBV	2403539	3728265	58.81%	7.980%
38	9.133	2504	2517	2539	rVV	3355152	5508368	86.89%	11.790%
39	9.236	2540	2549	2571	rVB	205315	326357	5.15%	0.699%
40	9.378	2584	2593	2610	rBV3	37745	68148	1.07%	0.146%
41	9.538	2631	2643	2674	rVB	2154650	3366638	53.11%	7.206%
42	9.673	2677	2685	2690	rBV5	6950	10501	0.17%	0.022%
43	9.706	2692	2695	2702	rVB4	4889	5114	0.08%	0.011%
44	9.808	2717	2727	2735	rBV4	3794	8214	0.13%	0.018%
45	9.853	2737	2741	2751	rVB6	5356	8553	0.13%	0.018%
46	9.931	2753	2765	2777	rBV	335294	522909	8.25%	1.119%
47	10.001	2777	2787	2803	rVB	168995	276172	4.36%	0.591%
48	10.214	2841	2853	2868	rBV	275876	428424	6.76%	0.917%
49	10.352	2883	2896	2912	rBV	259314	423107	6.67%	0.906%
50	10.445	2913	2925	2929	rBV	415167	656789	10.36%	1.406%
51	10.535	2944	2953	2966	rVB	257526	387043	6.11%	0.828%
52	10.693	2995	3002	3005	rBV3	6327	7453	0.12%	0.016%
53	10.734	3005	3015	3040	rVB	388045	652057	10.29%	1.396%
54	10.860	3047	3054	3059	rBV2	10787	15303	0.24%	0.033%
55	10.911	3059	3070	3087	rVB	2157523	3164811	49.92%	6.774%
56	10.988	3088	3094	3102	rVB3	10946	14221	0.22%	0.030%
57	11.040	3103	3110	3113	rBV3	13358	16776	0.26%	0.036%
58	11.085	3119	3124	3136	rVB2	30564	44619	0.70%	0.095%
59	11.191	3147	3157	3164	rBV3	38811	62205	0.98%	0.133%
60	11.246	3164	3174	3191	rVB	536038	838373	13.22%	1.794%
61	11.332	3191	3201	3214	rBV	507803	747158	11.79%	1.599%
62	11.525	3250	3261	3273	rBV	268205	472143	7.45%	1.011%
63	11.625	3281	3292	3316	rVB2	593277	1224184	19.31%	2.620%
64	11.744	3319	3329	3344	rBV	105974	166557	2.63%	0.356%
65	11.818	3344	3352	3366	rVB	36489	57697	0.91%	0.123%
66	11.908	3368	3380	3385	rBV	72131	116355	1.84%	0.249%
67	12.014	3404	3413	3432	rBV2	100415	174735	2.76%	0.374%
68	12.114	3436	3444	3457	rBV2	38642	77531	1.22%	0.166%
69	12.271	3481	3493	3499	rBV8	16229	37319	0.59%	0.080%
70	12.313	3501	3506	3512	rVB2	23151	29071	0.46%	0.062%
71	12.413	3530	3537	3544	rVB	51026	71794	1.13%	0.154%
72	12.464	3544	3553	3564	rVB	107548	159766	2.52%	0.342%
73	12.554	3573	3581	3588	rBV8	8598	13663	0.22%	0.029%
74	12.644	3602	3609	3616	rBV4	7105	9482	0.15%	0.020%
75	12.725	3625	3634	3643	rVB2	39881	57704	0.91%	0.124%
76	12.879	3664	3682	3695	rBV2	139511	237946	3.75%	0.509%

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 Title : VOC Analysis

77	12.998	3715	3719	3724	rVB2	4744	4650	0.07%	0.010%
78	13.046	3725	3734	3745	rBV4	46806	77040	1.22%	0.165%
79	13.165	3762	3771	3777	rBV8	11604	18032	0.28%	0.039%
80	13.213	3780	3786	3792	rVV4	9319	12488	0.20%	0.027%
81	13.262	3793	3801	3810	rBV3	52754	83198	1.31%	0.178%
82	13.500	3863	3875	3900	rVB	2206876	3399116	53.62%	7.275%
83	13.619	3903	3912	3926	rVB	56347	93793	1.48%	0.201%
84	13.705	3930	3939	3945	rBV7	11667	23006	0.36%	0.049%
85	13.831	3970	3978	3981	rBV6	6103	8195	0.13%	0.018%
86	13.860	3983	3987	3995	rVV	17009	23281	0.37%	0.050%
87	13.908	3995	4002	4017	rVB3	13614	23722	0.37%	0.051%
88	14.094	4049	4060	4072	rBV6	14760	34463	0.54%	0.074%
89	14.229	4096	4102	4114	rVB3	7734	10330	0.16%	0.022%
90	14.294	4114	4122	4132	rVB3	11088	17460	0.28%	0.037%
91	14.429	4156	4164	4178	rBV5	10299	18143	0.29%	0.039%
92	14.631	4216	4227	4240	rBV	95223	149782	2.36%	0.321%
93	14.815	4273	4284	4297	rBV	82064	132317	2.09%	0.283%
94	15.361	4442	4454	4463	rBV3	17476	29222	0.46%	0.063%
95	15.506	4492	4499	4507	rBV5	4068	6152	0.10%	0.013%
96	15.554	4507	4514	4520	rBV3	5027	7105	0.11%	0.015%
97	15.667	4538	4549	4561	rVV4	11176	19384	0.31%	0.041%
98	15.811	4585	4594	4599	rVV6	5197	8392	0.13%	0.018%
99	15.850	4601	4606	4614	rVB3	6415	8864	0.14%	0.019%
100	16.290	4736	4743	4754	rBV9	2722	4956	0.08%	0.011%

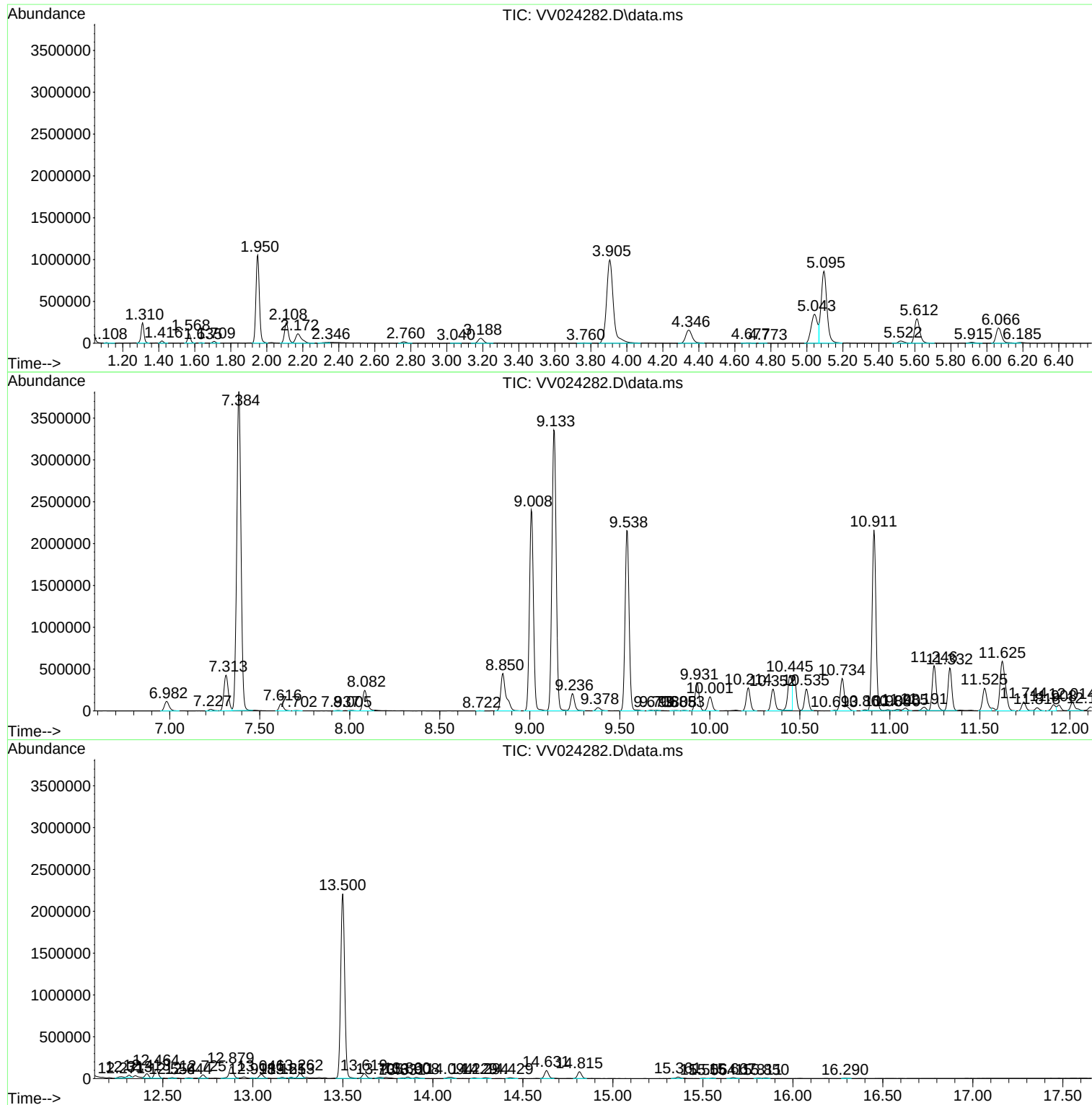
Sum of corrected areas: 46721856

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM123121WMA.M
Quant Title : VOC Analysis

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV010522\
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Instrument :
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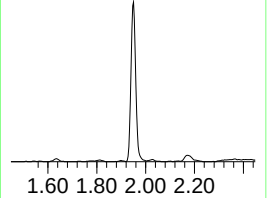
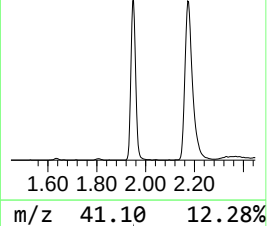
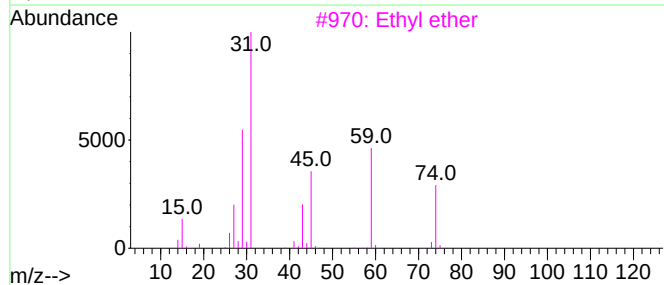
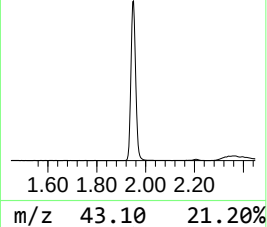
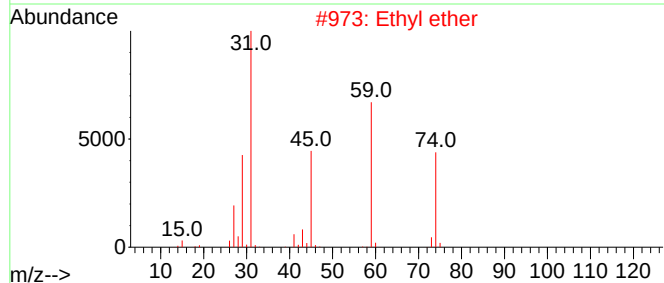
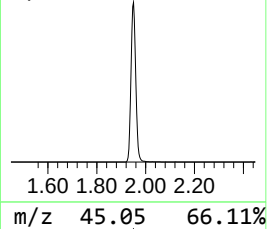
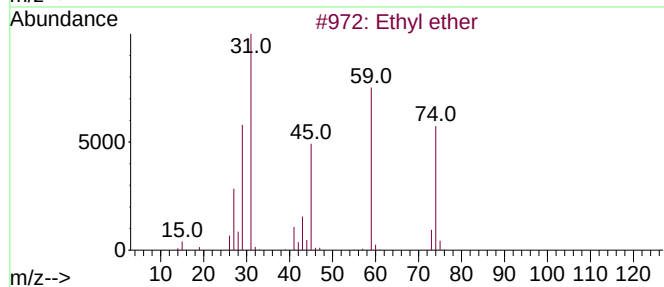
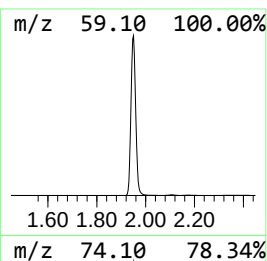
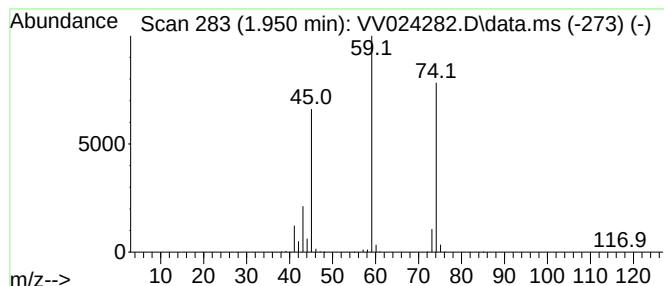
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM123121WMA.M
Quant Title : VOC Analysis

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

Peak Number 1 Ethyl ether Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.950	119.47 ug/L	1404230	1,4-Difluorobenzene	5.612

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethyl ether			74	C4H10O	000060-29-7	90
2	Ethyl ether			74	C4H10O	000060-29-7	90
3	Ethyl ether			74	C4H10O	000060-29-7	83
4	Ethane, 1,2-diethoxy-			118	C6H14O2	000629-14-1	72
5	Ethyl ether			74	C4H10O	000060-29-7	72



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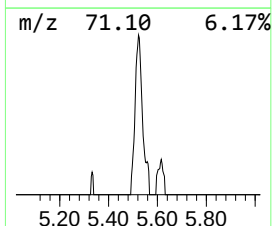
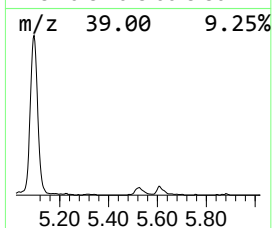
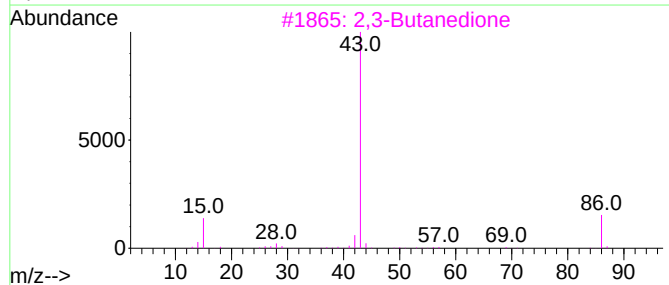
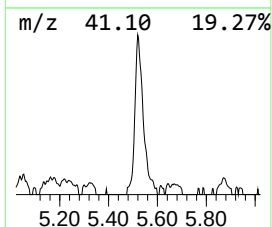
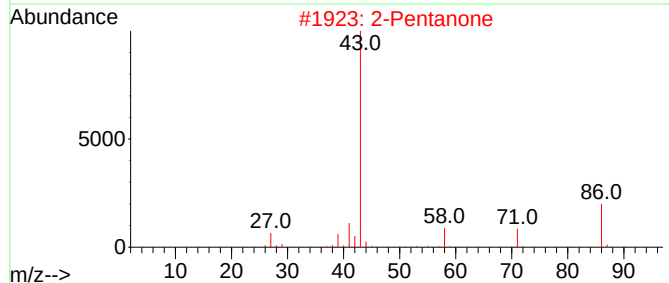
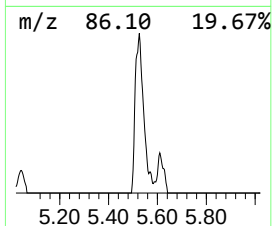
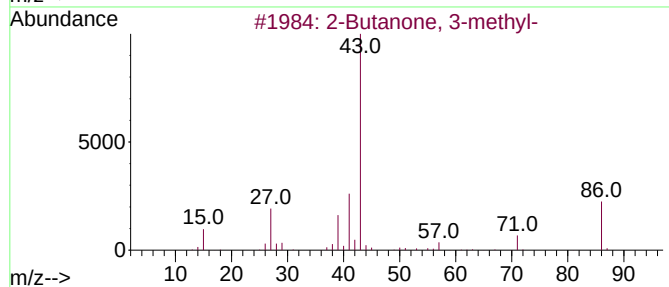
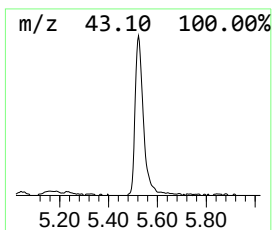
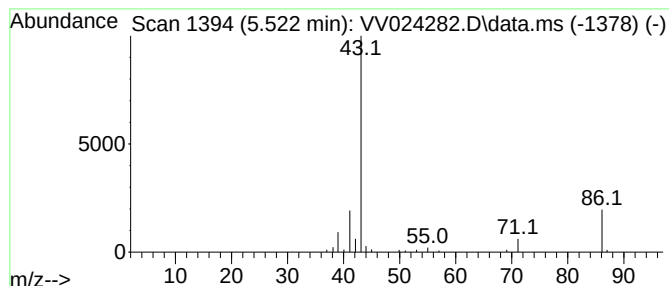
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM123121WMA.M
Quant Title : VOC Analysis

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

Peak Number 2 2-Butanone, 3-methyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.522	5.25 ug/L	61697	1,4-Difluorobenzene	5.612

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanone, 3-methyl-			86	C5H10O	000563-80-4	80
2	2-Pentanone			86	C5H10O	000107-87-9	72
3	2,3-Butanedione			86	C4H6O2	000431-03-8	72
4	2-Butanone, 3-methyl-			86	C5H10O	000563-80-4	59
5	2-Butanone, 3-methyl-			86	C5H10O	000563-80-4	59



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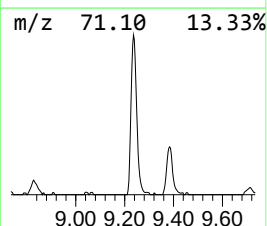
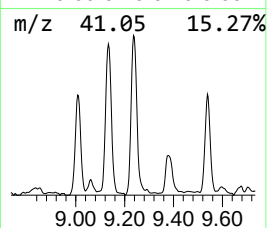
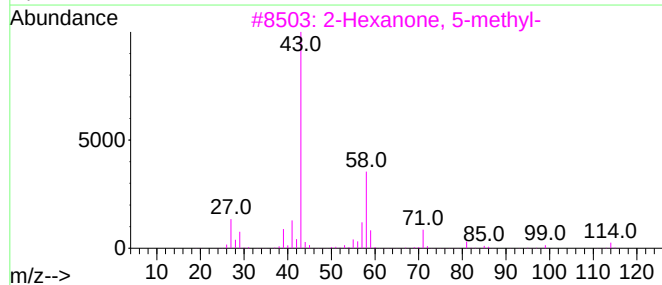
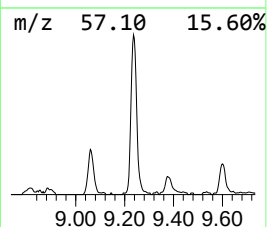
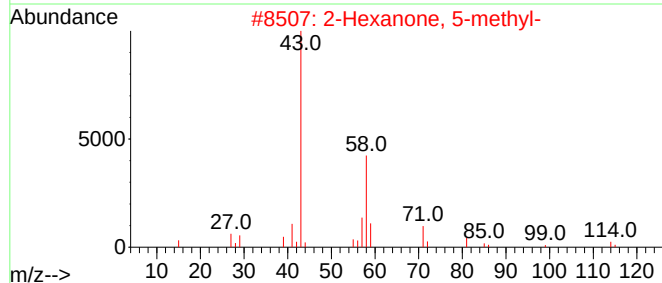
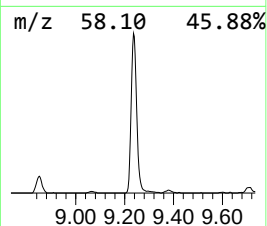
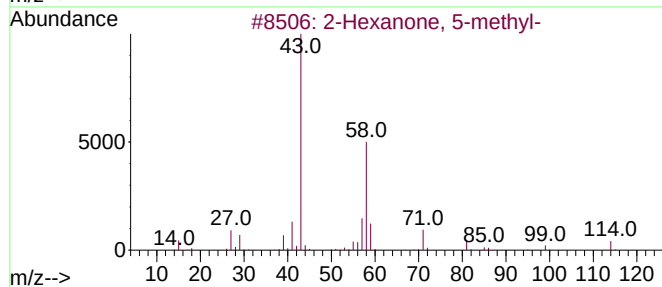
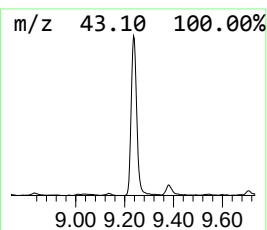
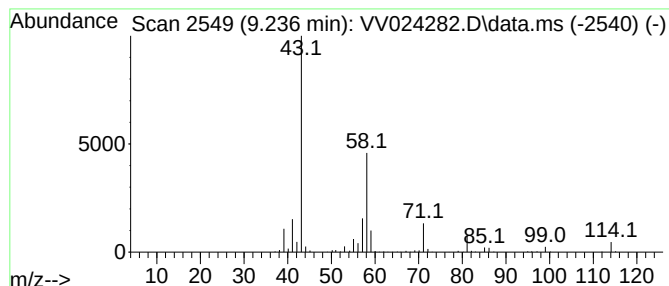
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM123121WMA.M
Quant Title : VOC Analysis

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

Peak Number 4 2-Hexanone, 5-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.236	17.99 ug/L	326357	Chlorobenzene-d5	8.850

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanone, 5-methyl-			114	C7H14O	000110-12-3	91
2	2-Hexanone, 5-methyl-			114	C7H14O	000110-12-3	91
3	2-Hexanone, 5-methyl-			114	C7H14O	000110-12-3	91
4	2-Hexanone, 5-methyl-			114	C7H14O	000110-12-3	91
5	2-Hexanone, 5-methyl-			114	C7H14O	000110-12-3	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV010522\
Data File : VV024282.D
Acq On : 05 Jan 2022 16:56
Operator : SY/MD
Sample : N1025-10
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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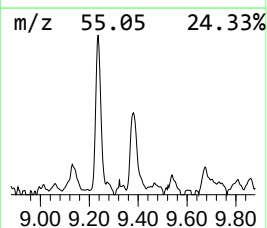
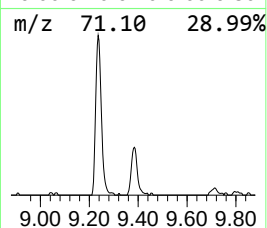
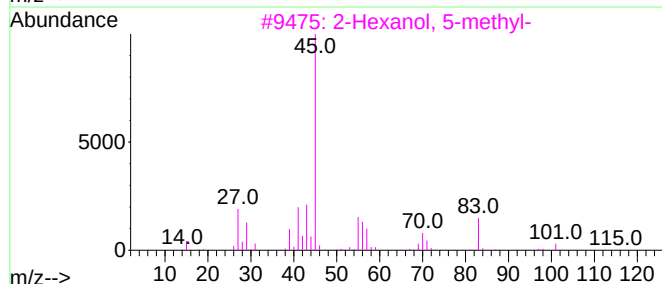
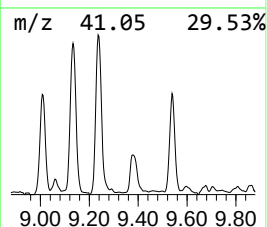
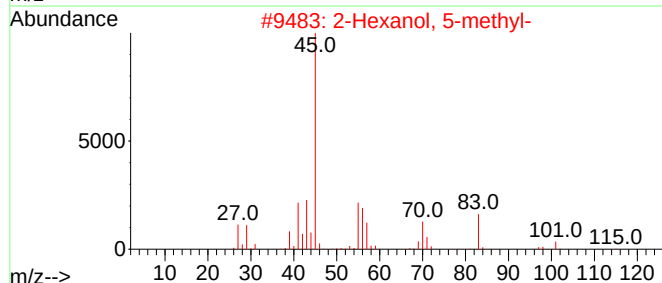
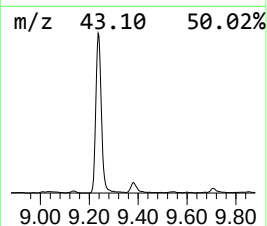
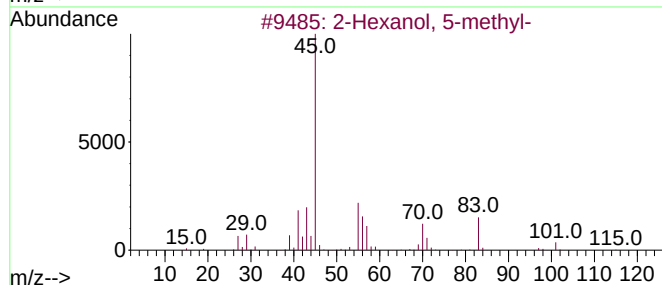
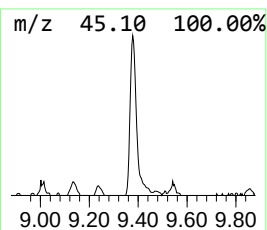
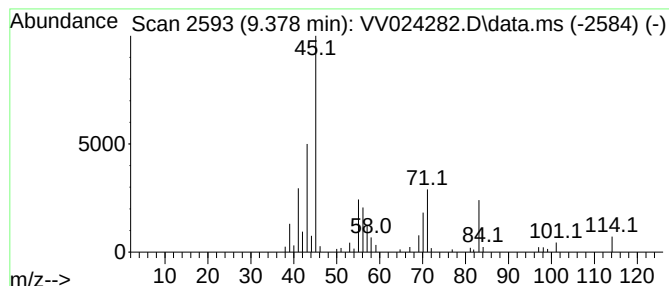
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Quant Title : VOC Analysis

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

Peak Number 5 2-Hexanol, 5-methyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.378	3.76 ug/L	68148	Chlorobenzene-d5	8.850

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanol, 5-methyl-			116	C7H16O	000627-59-8	78
2	2-Hexanol, 5-methyl-			116	C7H16O	000627-59-8	72
3	2-Hexanol, 5-methyl-			116	C7H16O	000627-59-8	64
4	(+)-5-Methyl-2-hexanol			116	C7H16O	111768-09-3	56
5	2-Decanol			158	C10H22O	001120-06-5	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV010522\
Data File : VV024282.D
Acq On : 05 Jan 2022 16:56
Operator : SY/MD
Sample : N1025-10
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLH4

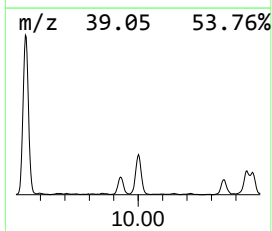
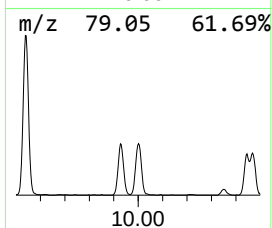
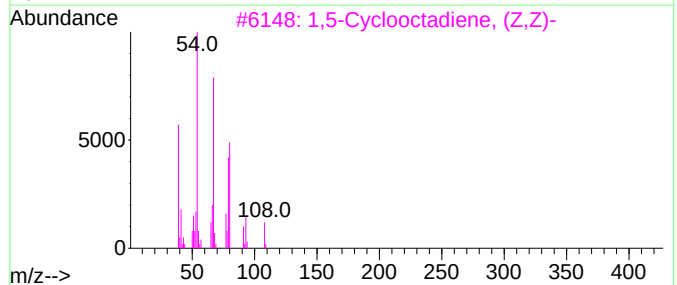
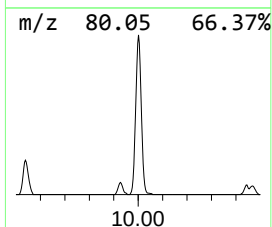
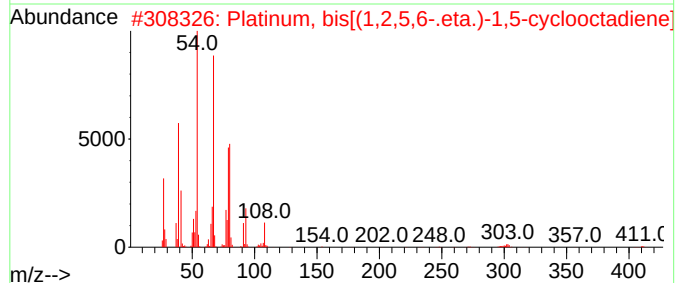
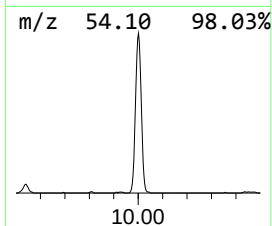
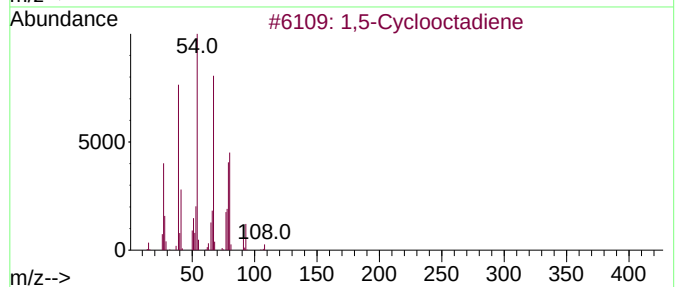
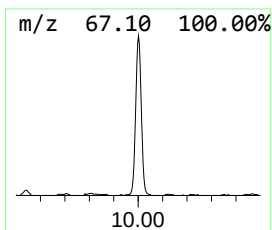
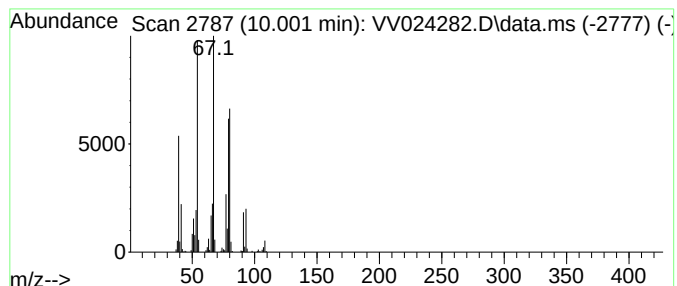
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Quant Title : VOC Analysis

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

Peak Number 6 1,5-Cyclooctadiene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.001	15.22 ug/L	276172	Chlorobenzene-d5	8.850

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,5-Cyclooctadiene	108	C8H12	000111-78-4	94
2			Platinum, bis[(1,2,5,6-.eta.)-1,...	411	C16H24Pt	012130-66-4	83
3			1,5-Cyclooctadiene, (Z,Z)-	108	C8H12	001552-12-1	83
4			1,5-Cyclooctadiene, (E,Z)-	108	C8H12	005259-71-2	78
5			1,5-Cyclooctadiene	108	C8H12	000111-78-4	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV010522\
Data File : VV024282.D
Acq On : 05 Jan 2022 16:56
Operator : SY/MD
Sample : N1025-10
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLH4

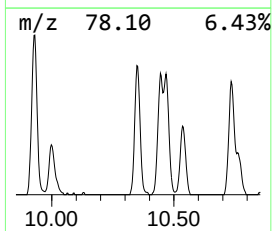
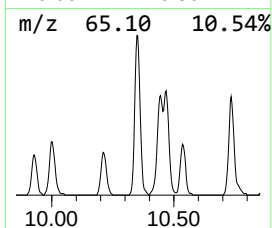
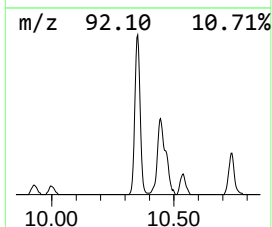
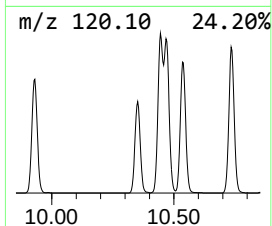
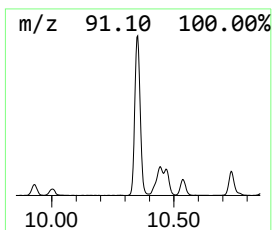
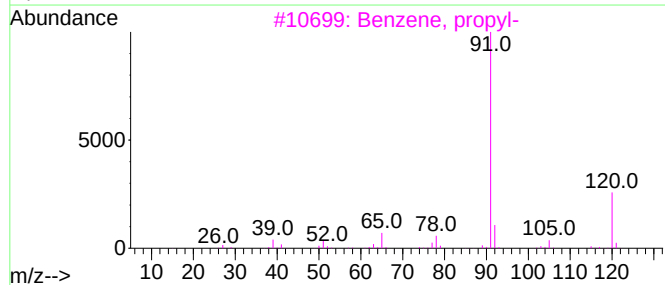
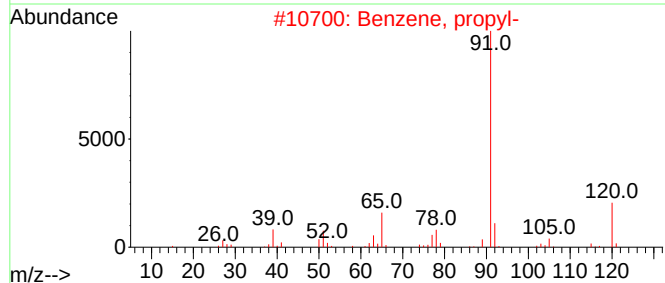
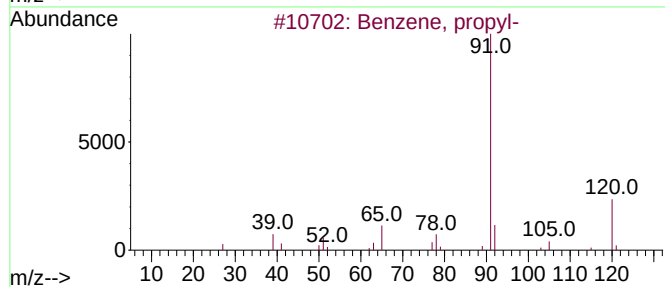
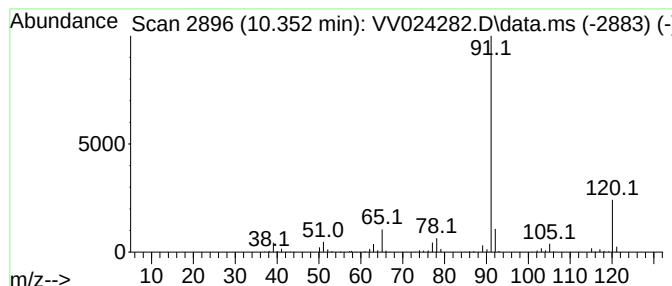
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Quant Title : VOC Analysis

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

Peak Number 7 Benzene, propyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.352	25.23 ug/L	423107	1,4-Dichlorobenzene-d4	11.246

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	93
2			Benzene, propyl-	120	C9H12	000103-65-1	90
3			Benzene, propyl-	120	C9H12	000103-65-1	90
4			Benzene, propyl-	120	C9H12	000103-65-1	90
5			Phenethylamine, N-benzyl-p-chloro-	245	C15H16ClN	013622-43-0	78



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV010522\
Data File : VV024282.D
Acq On : 05 Jan 2022 16:56
Operator : SY/MD
Sample : N1025-10
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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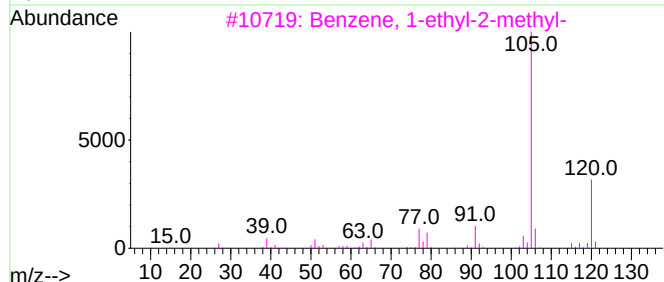
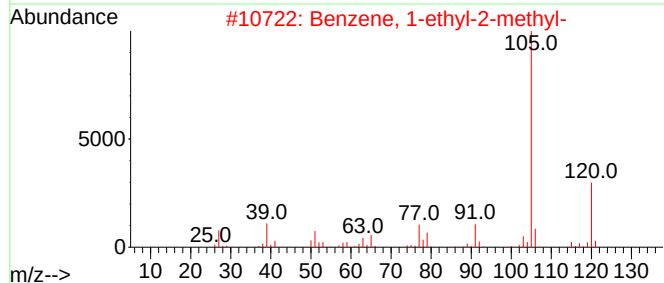
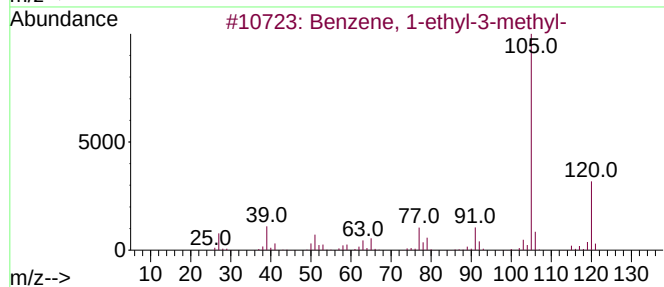
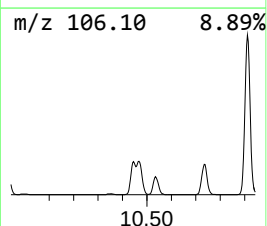
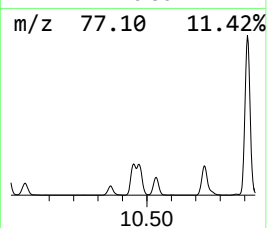
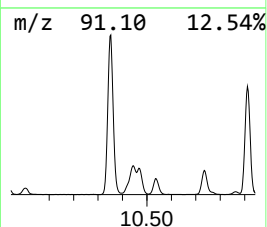
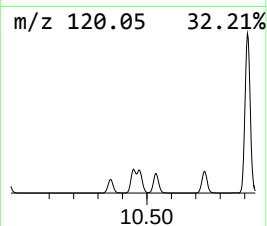
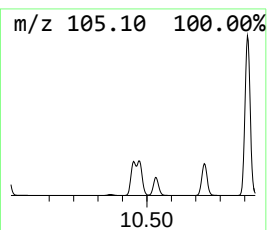
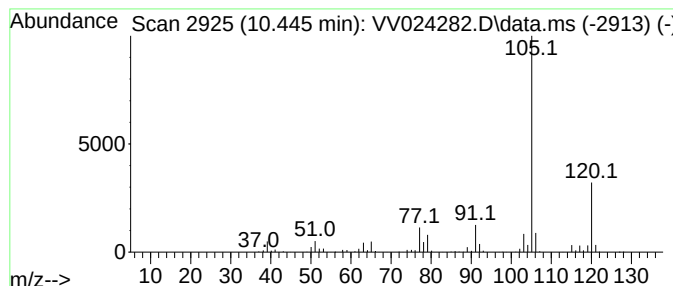
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Quant Title : VOC Analysis

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

Peak Number 8 Benzene, 1-ethyl-3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.445	39.17 ug/L	656789	1,4-Dichlorobenzene-d4	11.246

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
5			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV010522\
Data File : VV024282.D
Acq On : 05 Jan 2022 16:56
Operator : SY/MD
Sample : N1025-10
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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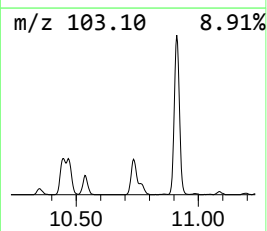
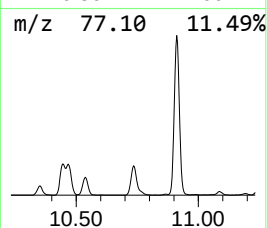
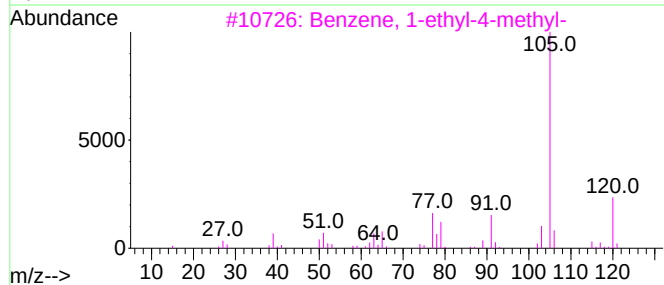
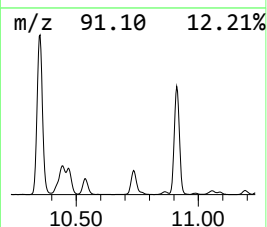
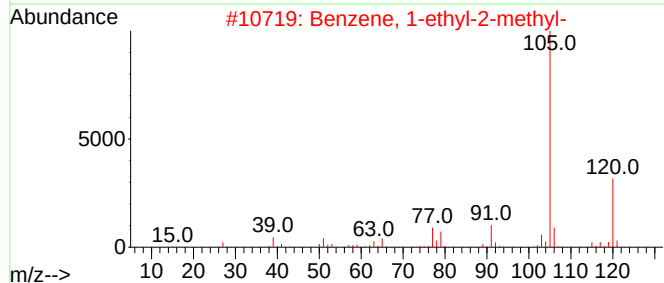
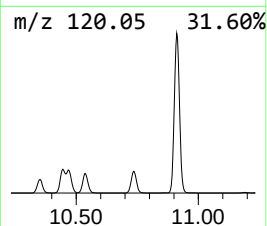
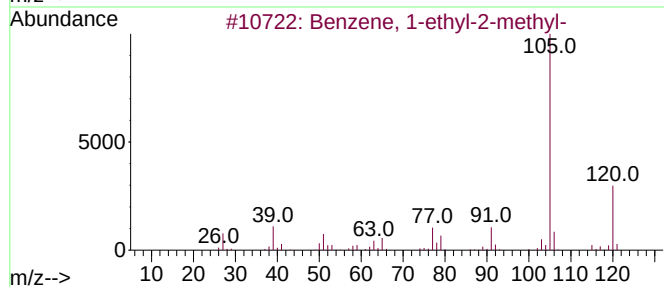
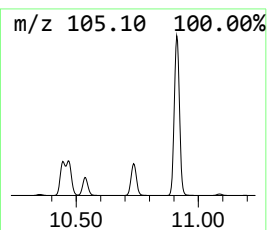
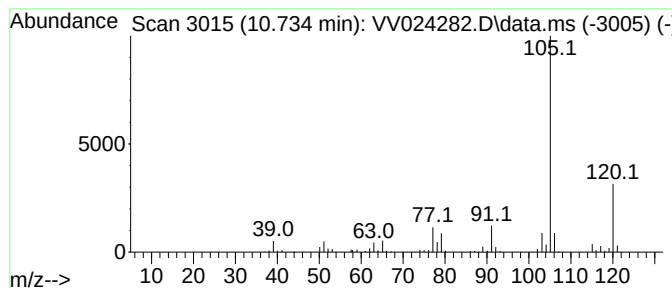
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Quant Title : VOC Analysis

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

Peak Number 9 Benzene, 1-ethyl-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.734	38.89 ug/L	652057	1,4-Dichlorobenzene-d4	11.246

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
5			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV010522\
Data File : VV024282.D
Acq On : 05 Jan 2022 16:56
Operator : SY/MD
Sample : N1025-10
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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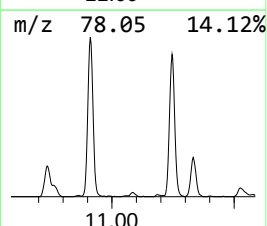
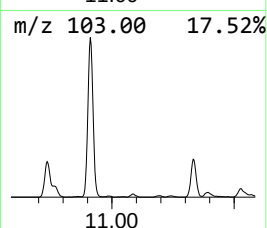
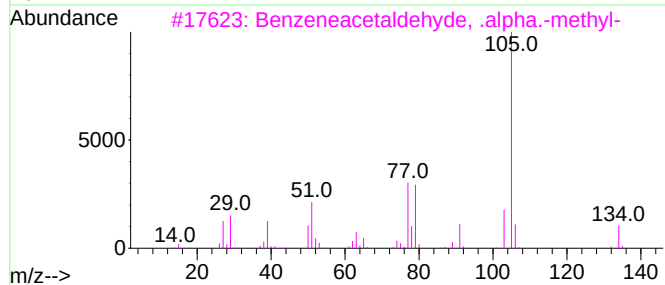
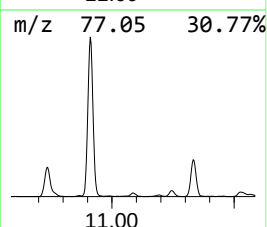
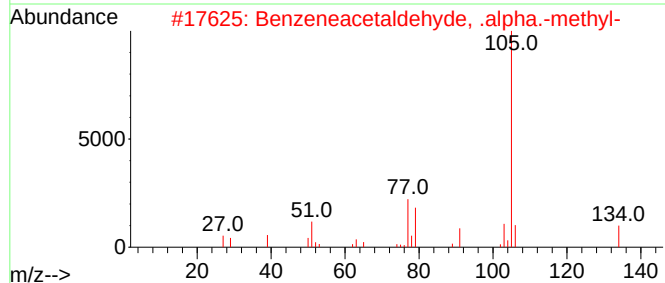
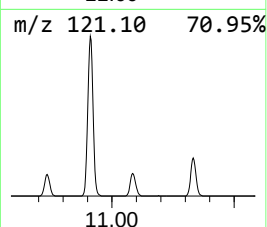
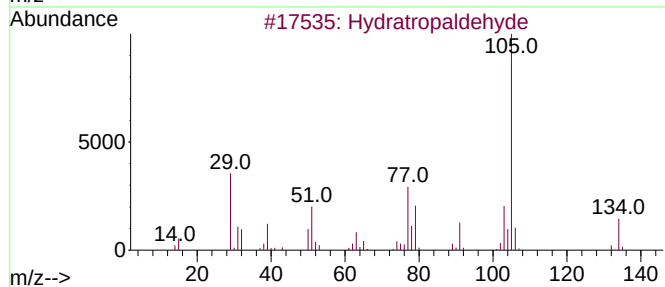
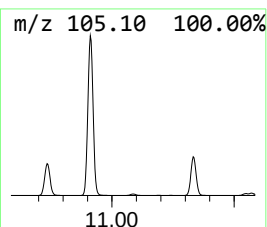
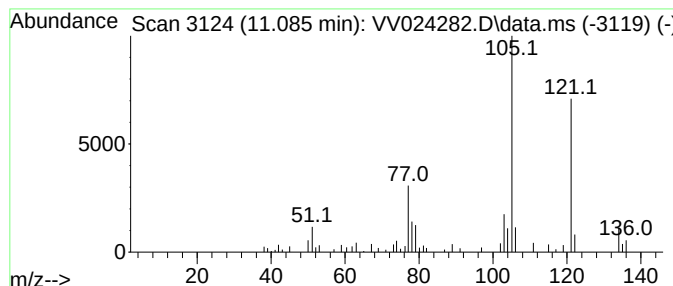
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Quant Title : VOC Analysis

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

Peak Number 10 Hydratropaldehyde Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.085	2.66 ug/L	44619	1,4-Dichlorobenzene-d4	11.246

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hydratropaldehyde	134	C9H10O	034713-70-7	64
2			Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	58
3			Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	52
4			Benzene, (1-nitroethyl)-	151	C8H9NO2	007214-61-1	50
5			1-Propanone, 1-phenyl-	134	C9H10O	000093-55-0	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV010522\
Data File : VV024282.D
Acq On : 05 Jan 2022 16:56
Operator : SY/MD
Sample : N1025-10
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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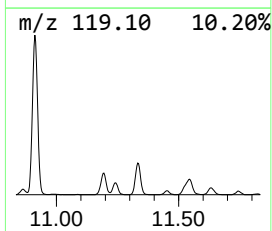
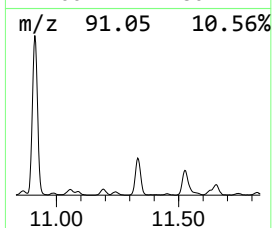
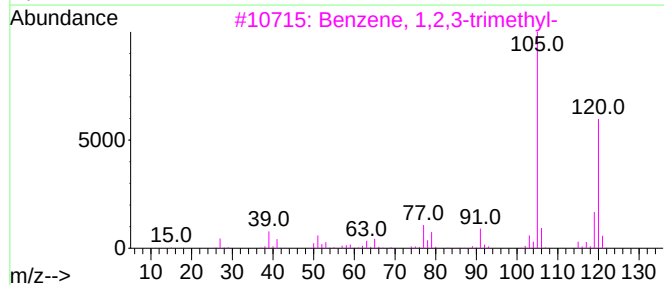
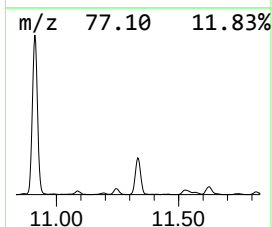
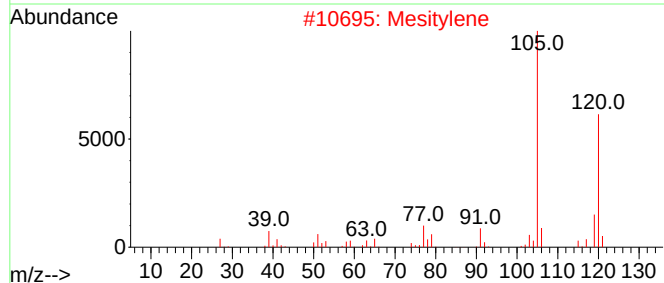
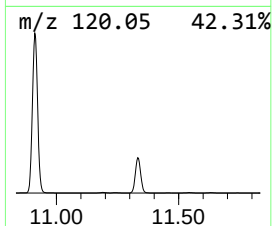
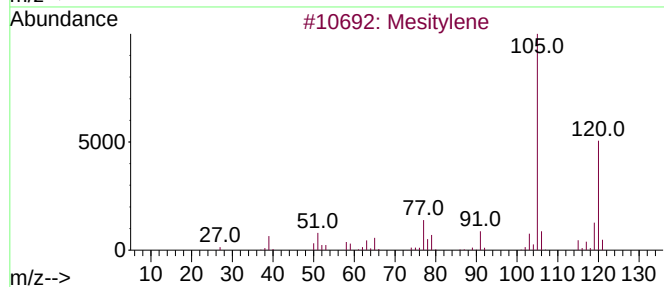
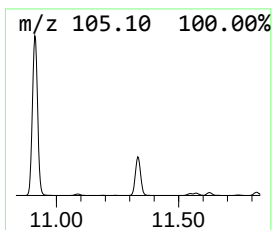
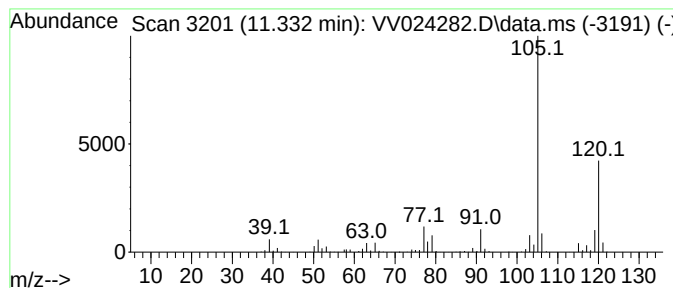
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Quant Title : VOC Analysis

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

Peak Number 11 Benzene, 1,2,3-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.333	44.56 ug/L	747158	1,4-Dichlorobenzene-d4	11.246

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	97
2			Mesitylene	120	C9H12	000108-67-8	97
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV010522\
Data File : VV024282.D
Acq On : 05 Jan 2022 16:56
Operator : SY/MD
Sample : N1025-10
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLH4

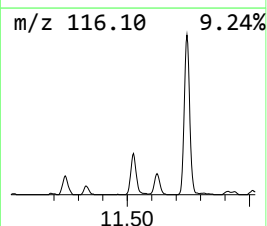
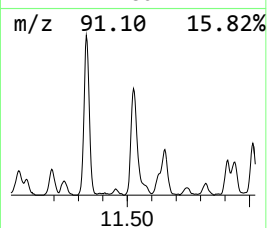
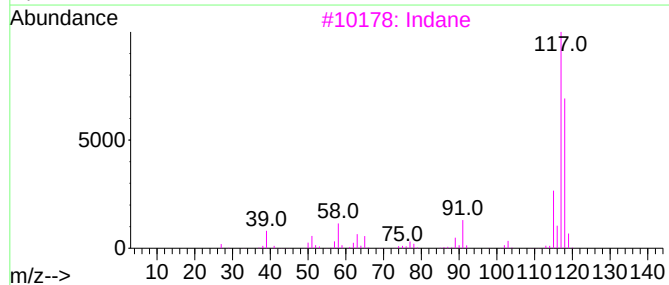
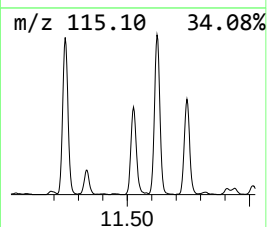
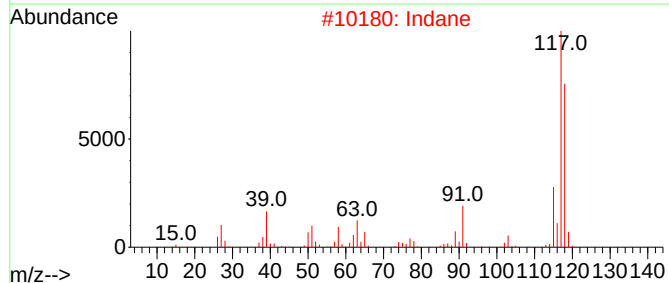
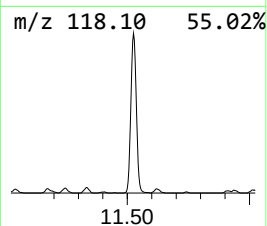
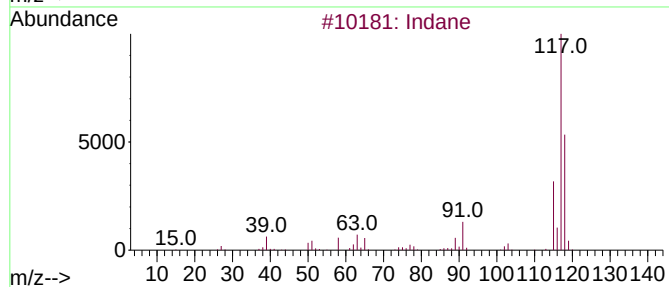
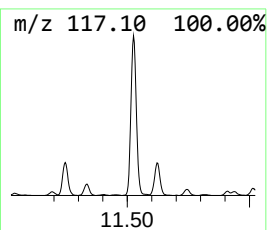
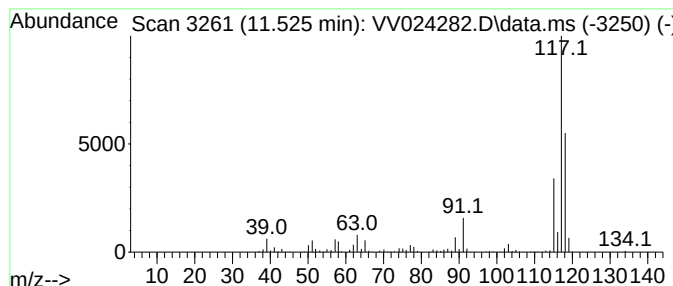
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Quant Title : VOC Analysis

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

Peak Number 12 Indane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.525	28.16 ug/L	472143	1,4-Dichlorobenzene-d4	11.246

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane			118	C9H10	000496-11-7	95
2	Indane			118	C9H10	000496-11-7	90
3	Indane			118	C9H10	000496-11-7	87
4	Benzene, cyclopropyl-			118	C9H10	000873-49-4	83
5	Benzene, 1-ethenyl-4-methyl-			118	C9H10	000622-97-9	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV010522\
Data File : VV024282.D
Acq On : 05 Jan 2022 16:56
Operator : SY/MD
Sample : N1025-10
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLH4

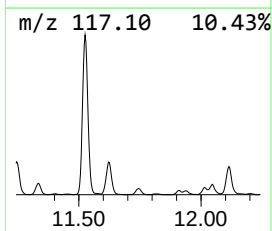
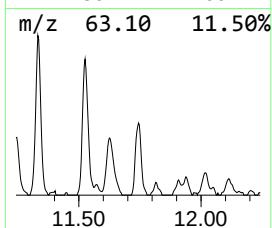
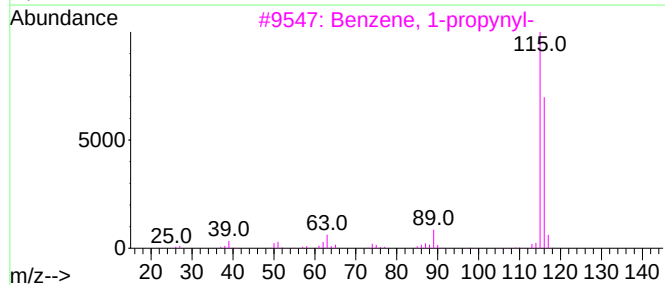
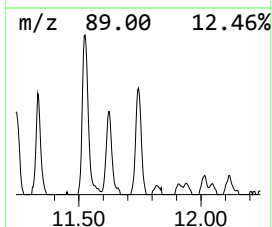
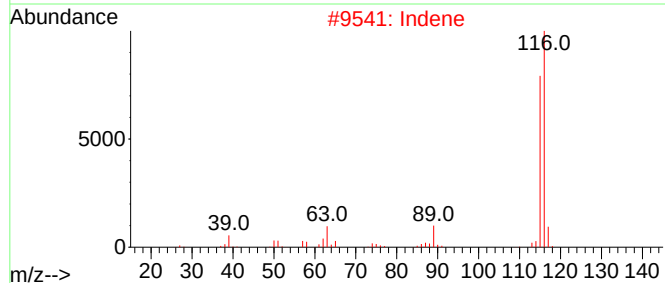
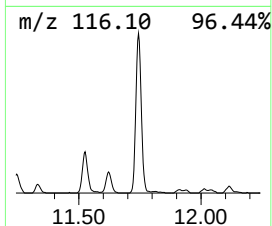
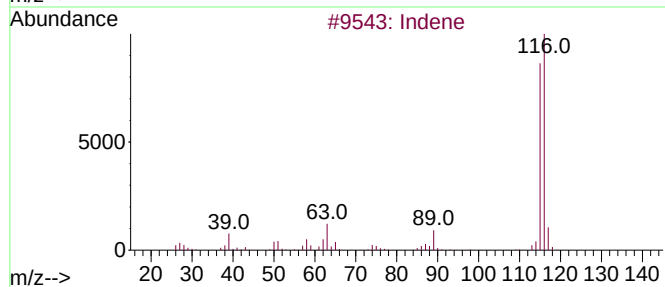
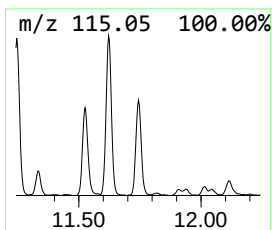
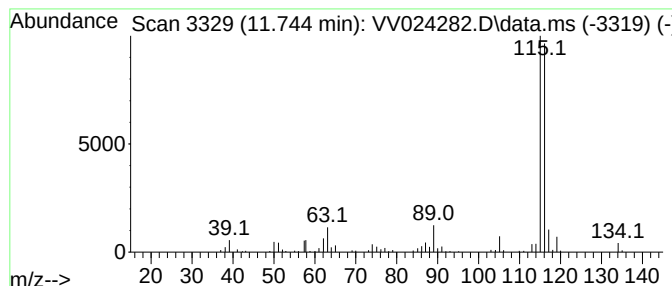
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Quant Title : VOC Analysis

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

Peak Number 13 Indene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.744	9.93 ug/L	166557	1,4-Dichlorobenzene-d4	11.246

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indene	116	C9H8	000095-13-6	97
2			Indene	116	C9H8	000095-13-6	95
3			Benzene, 1-propynyl-	116	C9H8	000673-32-5	94
4			3-Methylphenylacetylene	116	C9H8	000766-82-5	94
5			Benzene, 1-propynyl-	116	C9H8	000673-32-5	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV010522\
Data File : VV024282.D
Acq On : 05 Jan 2022 16:56
Operator : SY/MD
Sample : N1025-10
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLH4

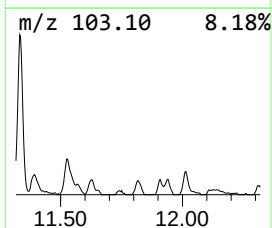
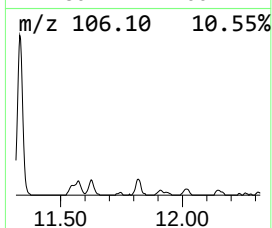
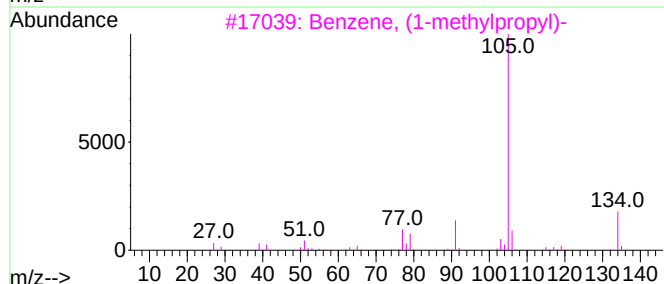
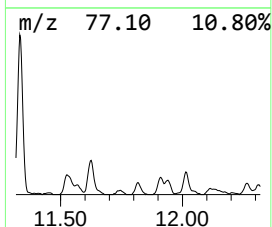
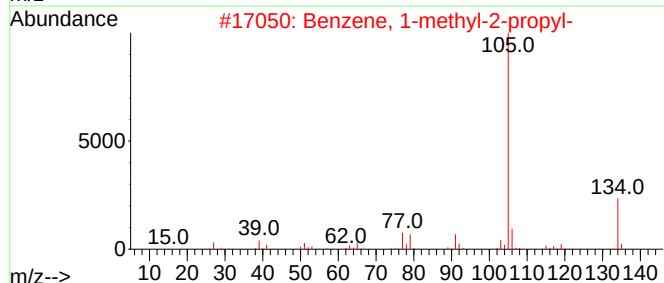
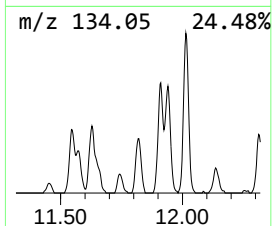
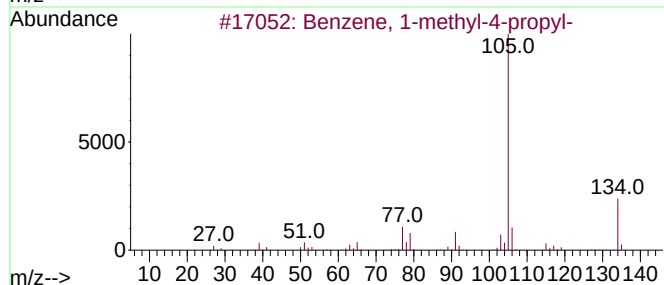
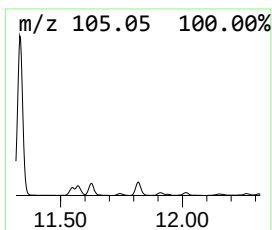
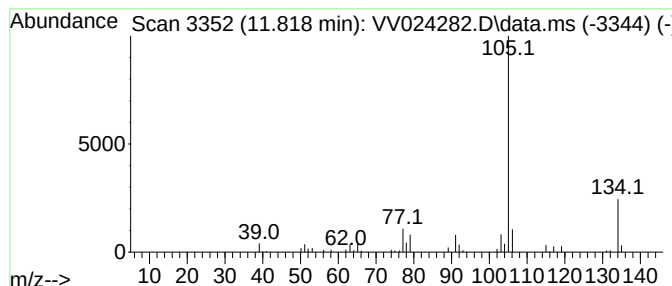
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Quant Title : VOC Analysis

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

Peak Number 14 Benzene, 1-methyl-4-propyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.818	3.44 ug/L	57697	1,4-Dichlorobenzene-d4	11.246

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	94
2			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	94
3			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	91
4			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
5			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV010522\
Data File : VV024282.D
Acq On : 05 Jan 2022 16:56
Operator : SY/MD
Sample : N1025-10
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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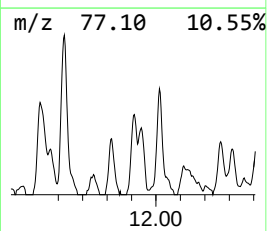
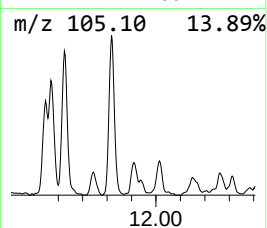
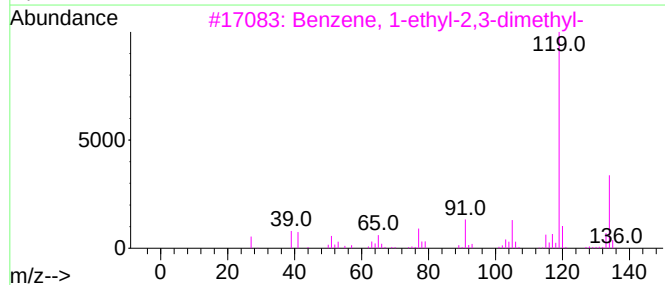
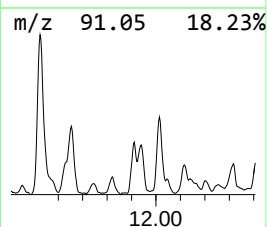
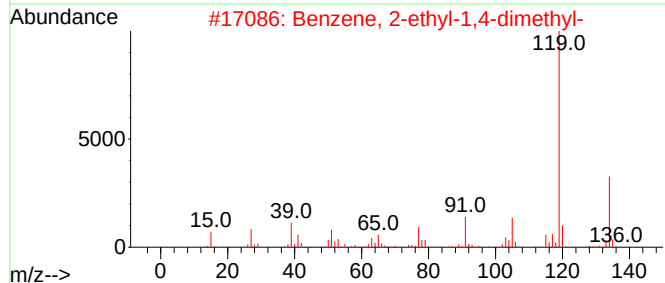
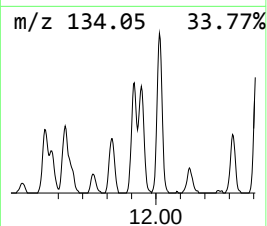
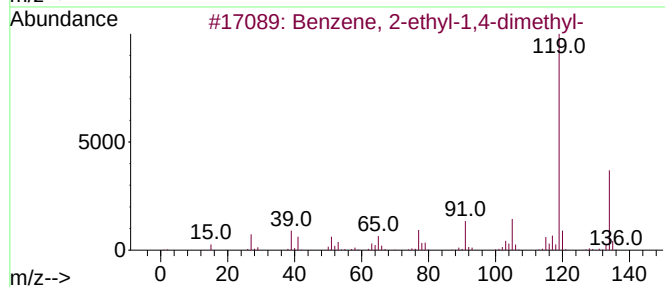
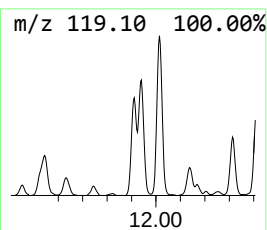
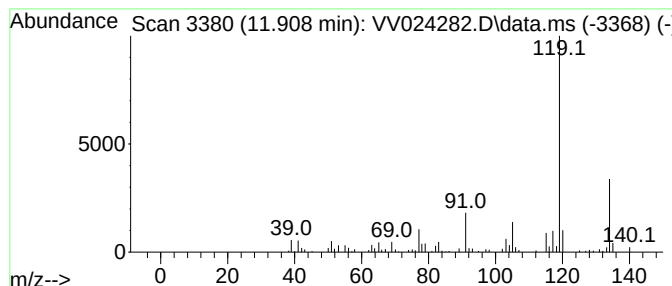
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Quant Title : VOC Analysis

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

Peak Number 15 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.908	6.94 ug/L	116355	1,4-Dichlorobenzene-d4	11.246

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
3			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4			o-Cymene	134	C10H14	000527-84-4	95
5			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV010522\
Data File : VV024282.D
Acq On : 05 Jan 2022 16:56
Operator : SY/MD
Sample : N1025-10
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLH4

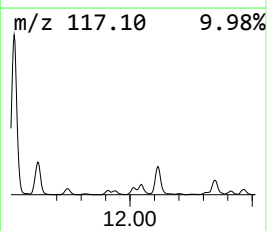
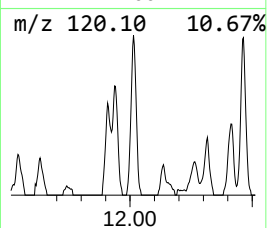
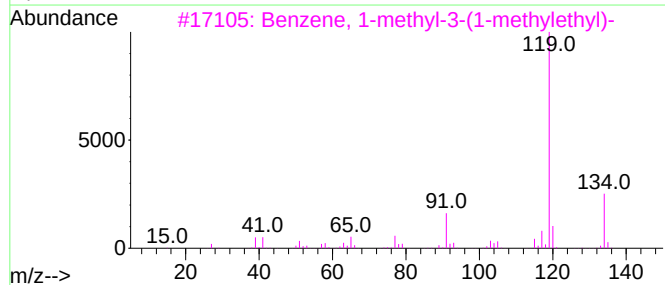
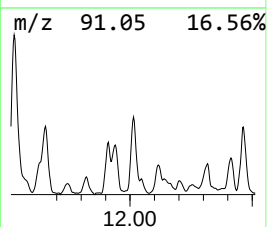
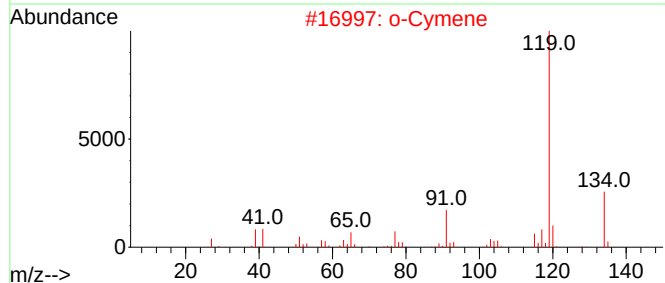
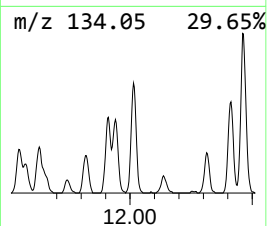
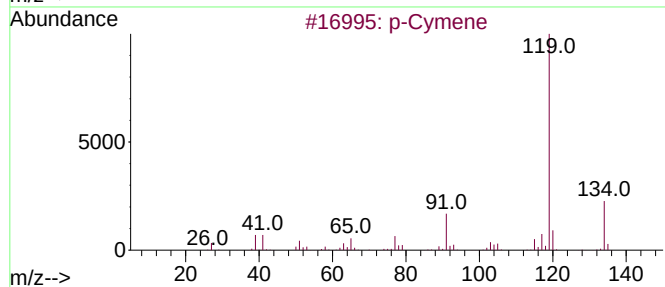
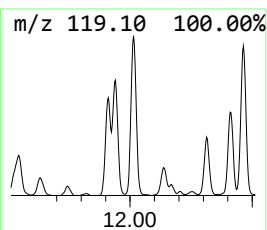
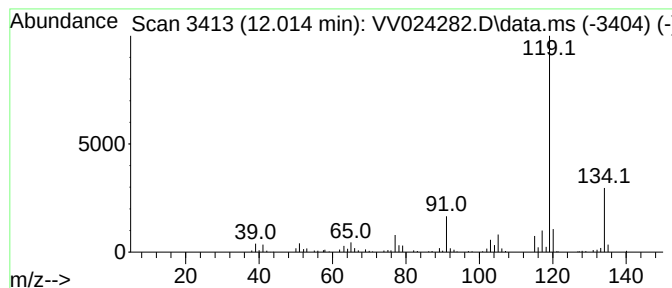
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Quant Title : VOC Analysis

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

Peak Number 16 p-Cymene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.014	10.42 ug/L	174735	1,4-Dichlorobenzene-d4	11.246

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			p-Cymene	134	C10H14	000099-87-6	97
2			o-Cymene	134	C10H14	000527-84-4	97
3			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	97
4			o-Cymene	134	C10H14	000527-84-4	97
5			o-Cymene	134	C10H14	000527-84-4	97



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV010522\
Data File : VV024282.D
Acq On : 05 Jan 2022 16:56
Operator : SY/MD
Sample : N1025-10
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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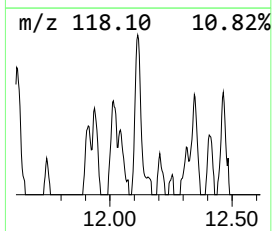
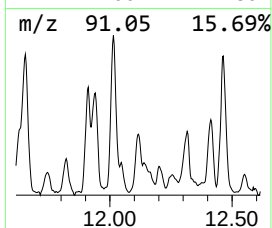
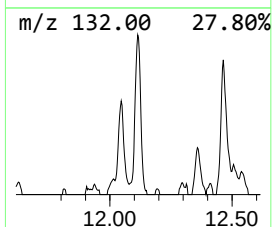
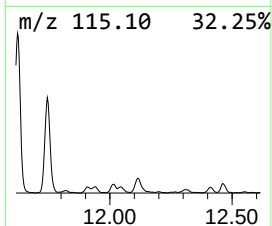
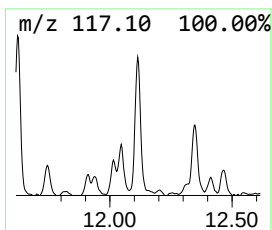
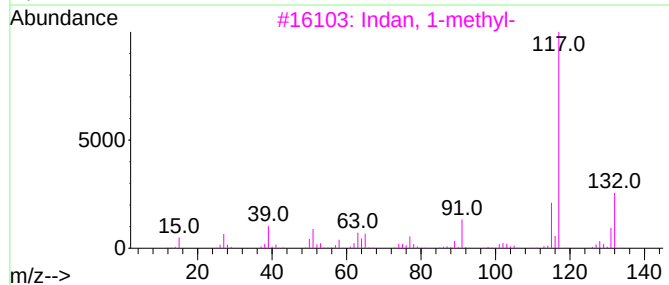
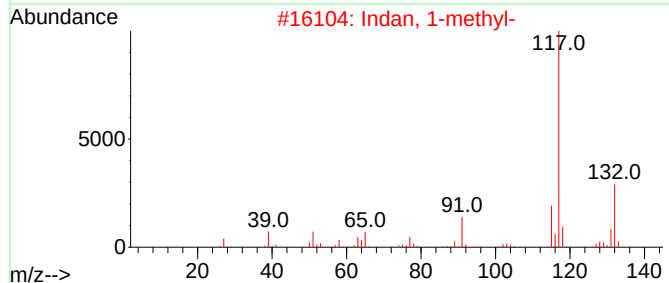
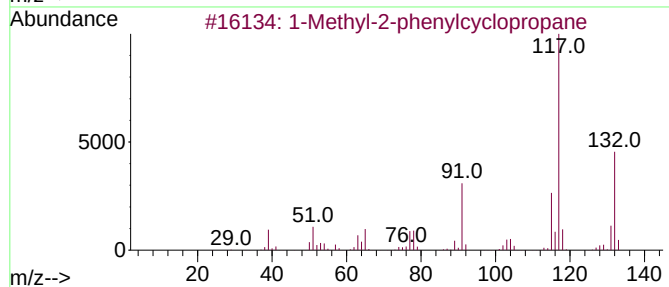
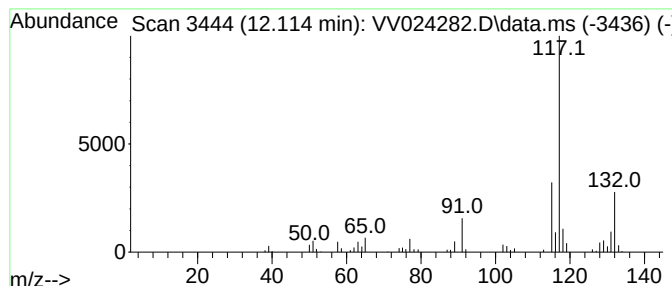
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Quant Title : VOC Analysis

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

Peak Number 17 1-Methyl-2-phenylcyclopropane Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.114	4.62 ug/L	77531	1,4-Dichlorobenzene-d4	11.246

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	90
2		Indan, 1-methyl-	132	C10H12	000767-58-8	87
3		Indan, 1-methyl-	132	C10H12	000767-58-8	87
4		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	87
5		1-Phenyl-1-butene	132	C10H12	000824-90-8	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV010522\
Data File : VV024282.D
Acq On : 05 Jan 2022 16:56
Operator : SY/MD
Sample : N1025-10
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

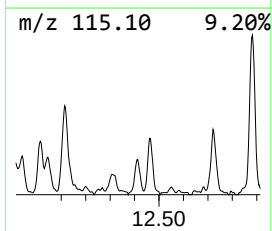
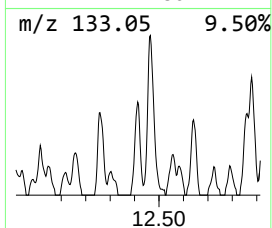
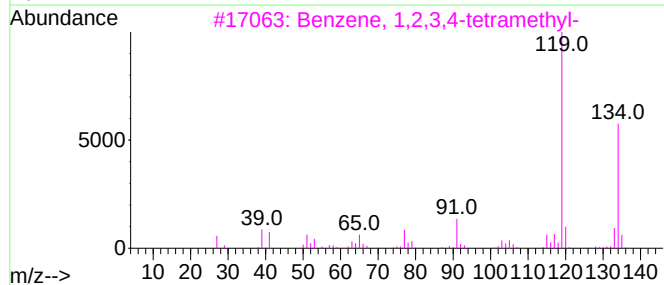
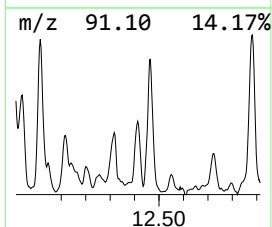
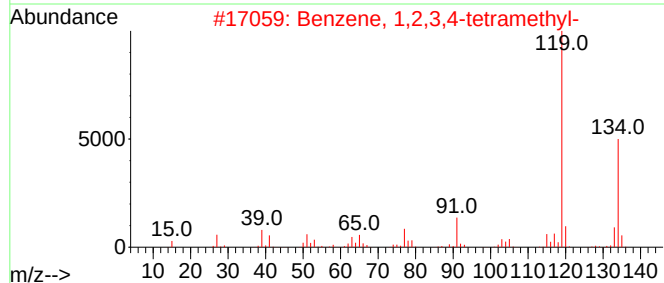
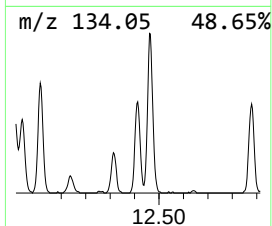
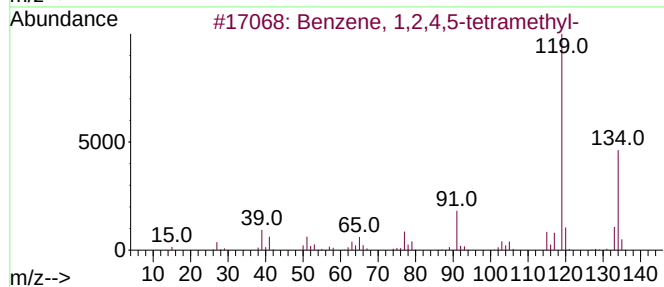
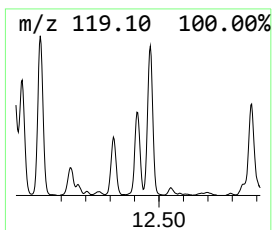
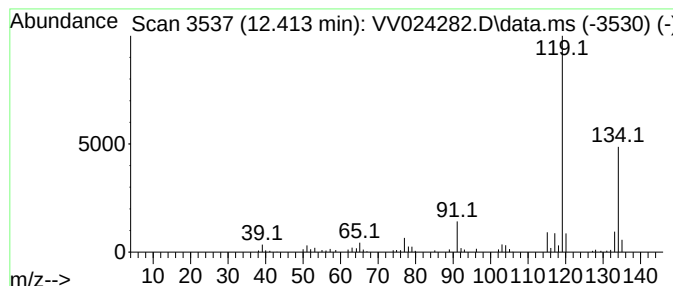
Instrument :
MSVOA_V
ClientSampleId :
BGLH4

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM123121WMA.M
Quant Title : VOC Analysis

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

Peak Number 18 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD			R.T.
12.413	4.28 ug/L	71794	1,4-Dichlorobenzene-d4			11.246
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-		134	C10H14	000095-93-2	95
2	Benzene, 1,2,3,4-tetramethyl-		134	C10H14	000488-23-3	95
3	Benzene, 1,2,3,4-tetramethyl-		134	C10H14	000488-23-3	94
4	Benzene, 1,2,3,4-tetramethyl-		134	C10H14	000488-23-3	94
5	Benzene, 1,2,3,5-tetramethyl-		134	C10H14	000527-53-7	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV010522\
Data File : VV024282.D
Acq On : 05 Jan 2022 16:56
Operator : SY/MD
Sample : N1025-10
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLH4

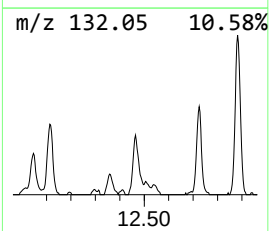
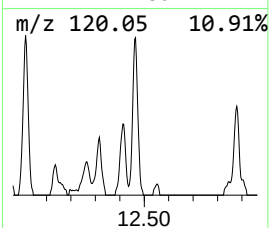
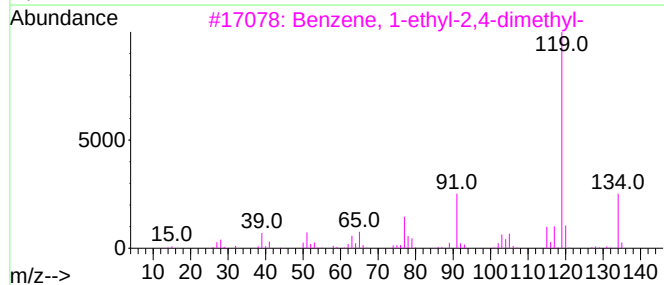
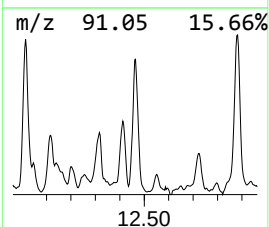
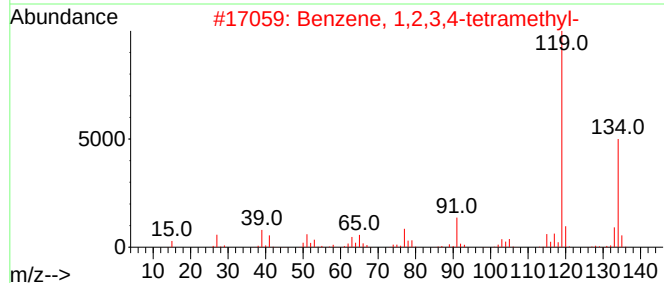
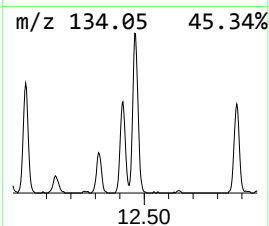
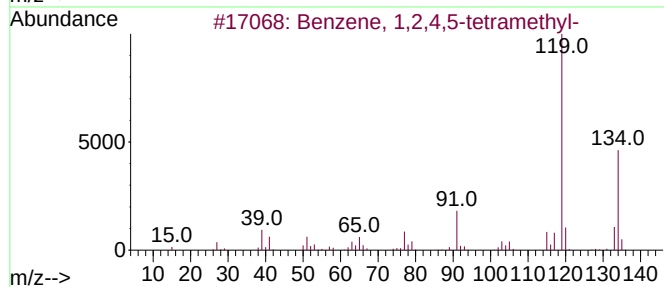
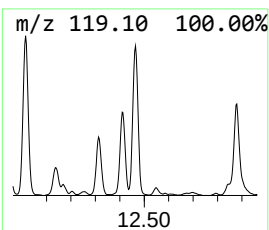
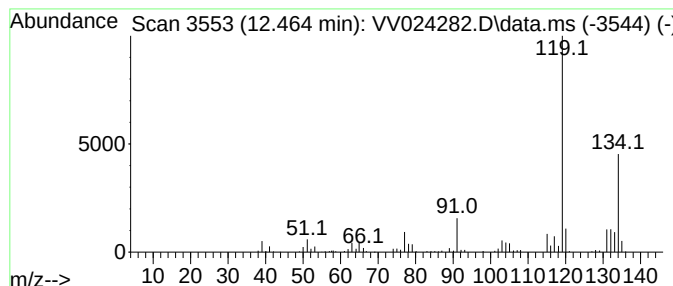
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Quant Title : VOC Analysis

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

Peak Number 19 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.464	9.53 ug/L	159766	1,4-Dichlorobenzene-d4	11.246

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
3			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
4			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	93
5			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV010522\
Data File : VV024282.D
Acq On : 05 Jan 2022 16:56
Operator : SY/MD
Sample : N1025-10
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLH4

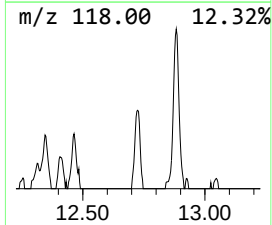
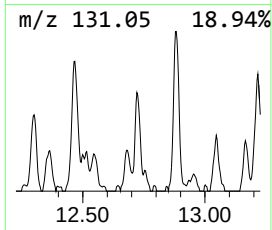
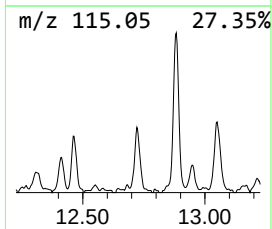
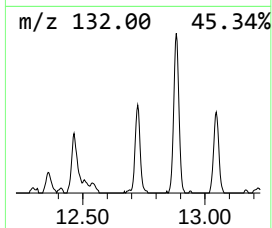
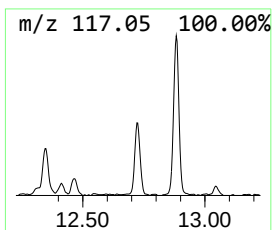
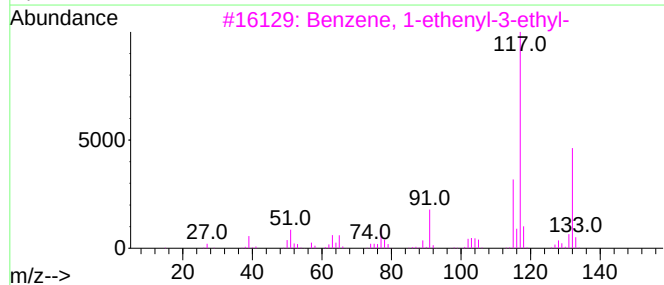
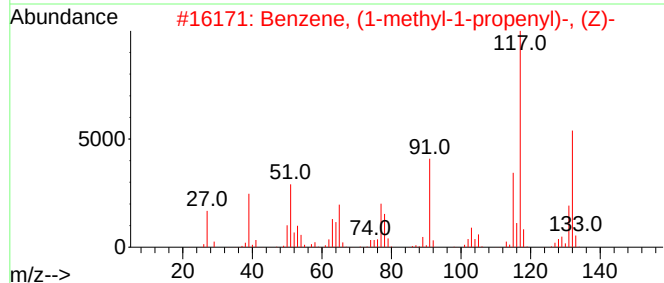
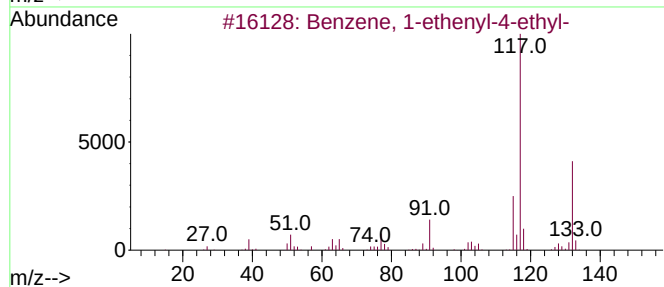
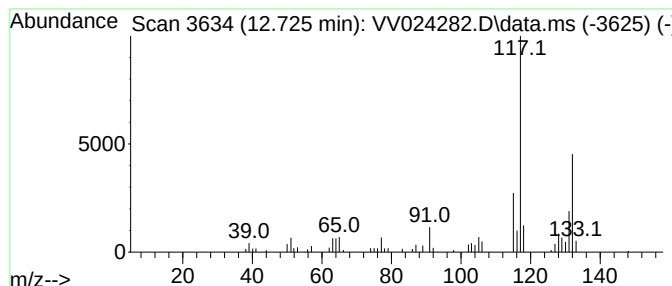
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Quant Title : VOC Analysis

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

Peak Number 20 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.725	3.44 ug/L	57704	1,4-Dichlorobenzene-d4	11.246

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	93
2		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	91
3		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
4		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
5		1-Phenyl-1-butene	132	C10H12	000824-90-8	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV010522\
Data File : VV024282.D
Acq On : 05 Jan 2022 16:56
Operator : SY/MD
Sample : N1025-10
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLH4

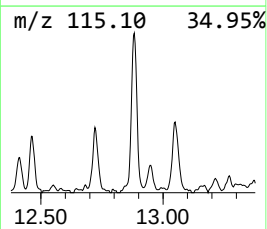
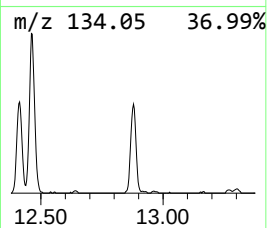
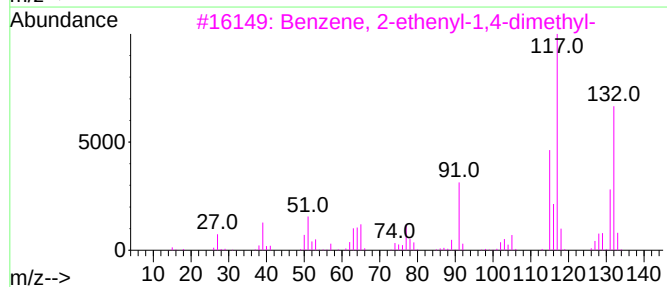
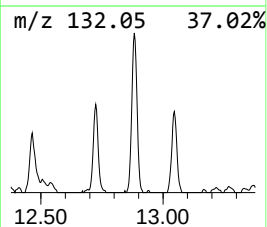
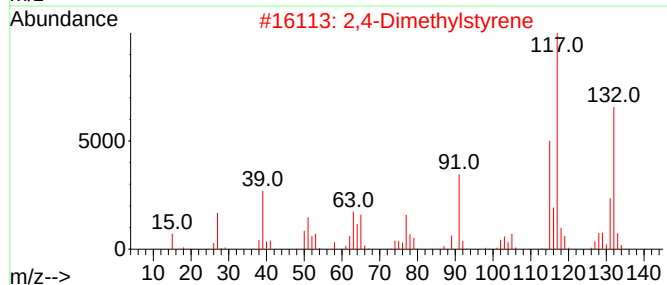
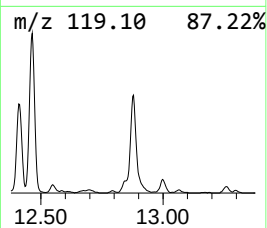
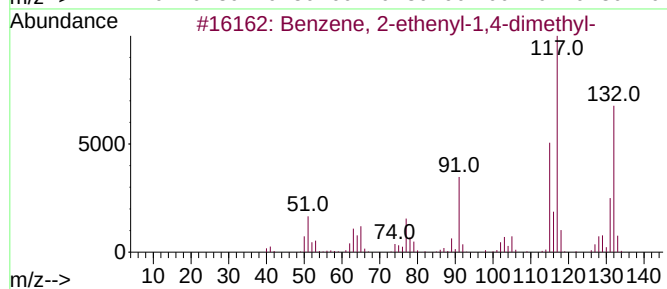
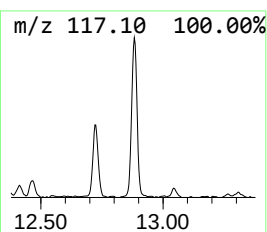
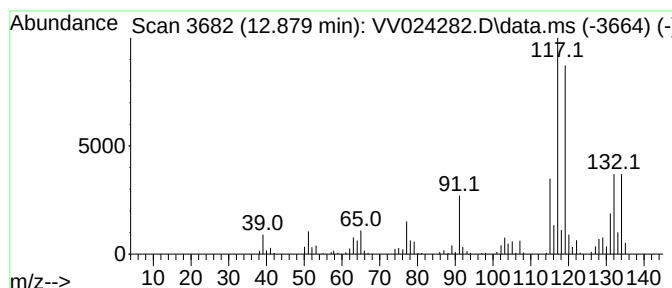
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Quant Title : VOC Analysis

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

Peak Number 21 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.879	14.19 ug/L	237946	1,4-Dichlorobenzene-d4	11.246

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	86
2			2,4-Dimethylstyrene	132	C10H12	002234-20-0	83
3			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	80
4			Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	70
5			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV010522\
Data File : VV024282.D
Acq On : 05 Jan 2022 16:56
Operator : SY/MD
Sample : N1025-10
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

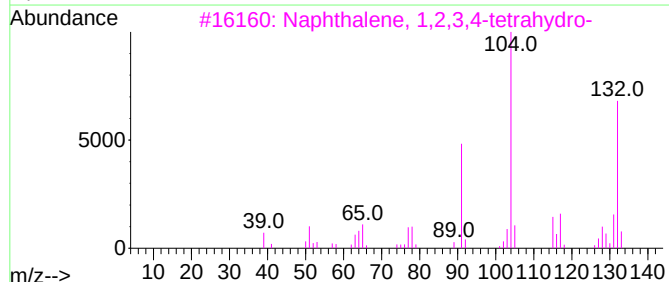
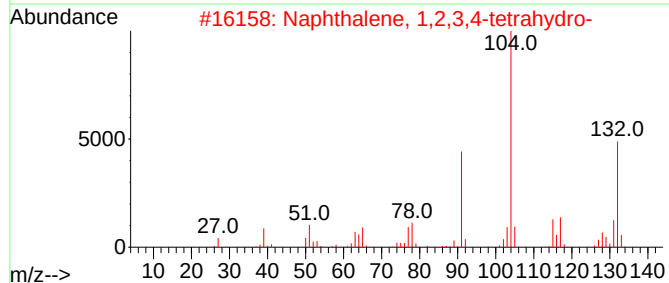
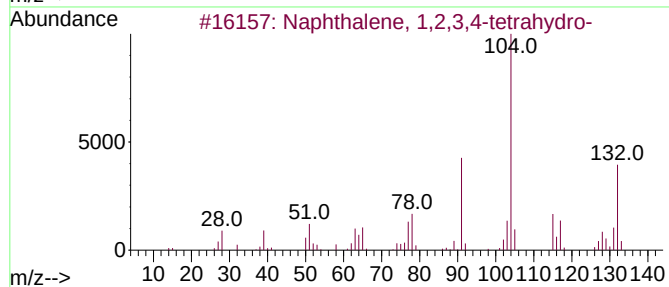
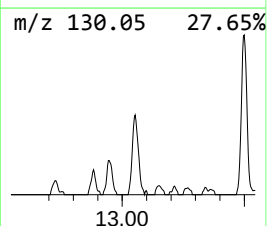
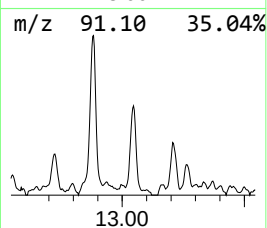
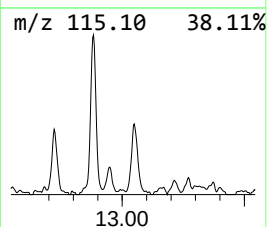
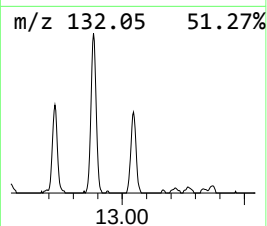
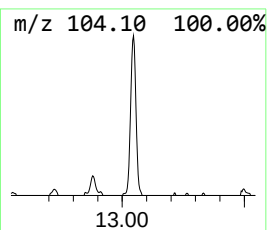
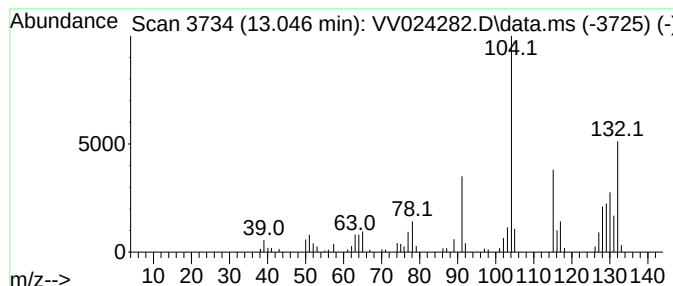
Instrument :
MSVOA_V
ClientSampleId :
BGLH4

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM123121WMA.M
Quant Title : VOC Analysis

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

Peak Number 22 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.046	4.59 ug/L	77040	1,4-Dichlorobenzene-d4		11.246	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-		132	C10H12	000119-64-2	78
2	Naphthalene, 1,2,3,4-tetrahydro-		132	C10H12	000119-64-2	78
3	Naphthalene, 1,2,3,4-tetrahydro-		132	C10H12	000119-64-2	78
4	Naphthalene, 1,2,3,4-tetrahydro-		132	C10H12	000119-64-2	60
5	Naphthalene, 1,2,3,4-tetrahydro-		132	C10H12	000119-64-2	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV010522\
Data File : VV024282.D
Acq On : 05 Jan 2022 16:56
Operator : SY/MD
Sample : N1025-10
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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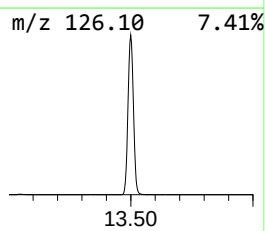
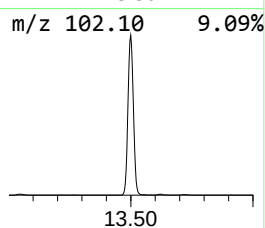
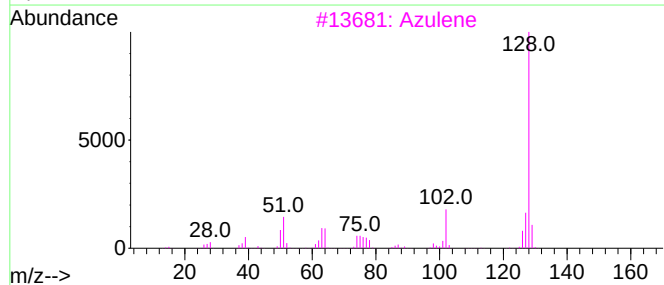
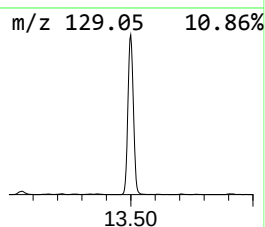
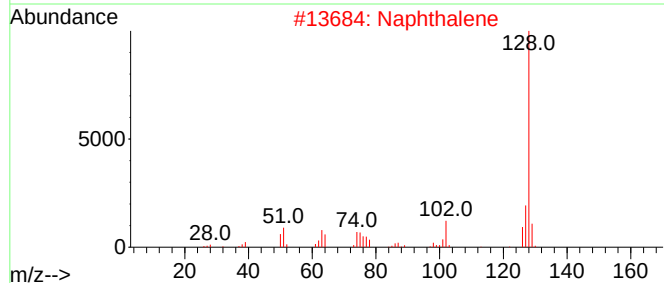
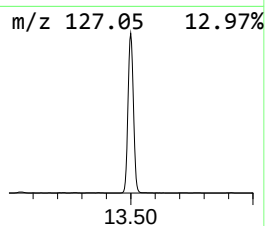
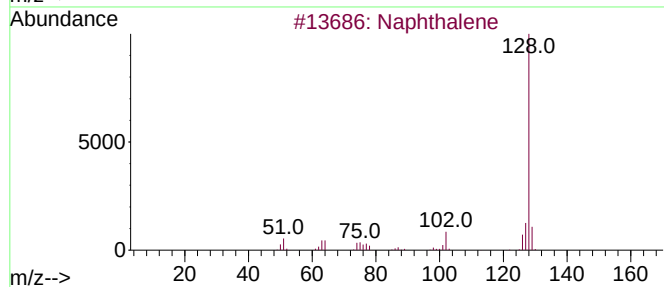
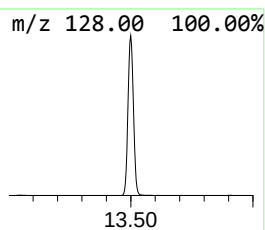
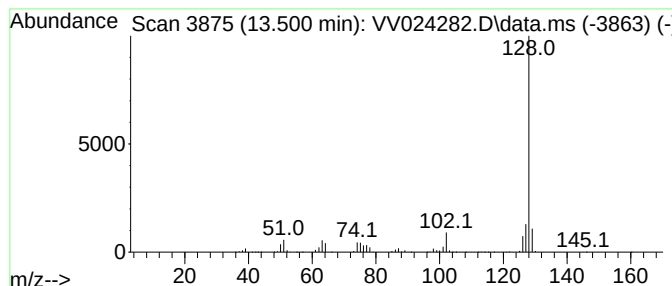
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Quant Title : VOC Analysis

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

Peak Number 23 Azulene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.500	202.72 ug/L	3399120	1,4-Dichlorobenzene-d4	11.246

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene	128	C10H8	000091-20-3	95
2			Naphthalene	128	C10H8	000091-20-3	95
3			Azulene	128	C10H8	000275-51-4	91
4			Azulene	128	C10H8	000275-51-4	91
5			Azulene	128	C10H8	000275-51-4	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV010522\
Data File : VV024282.D
Acq On : 05 Jan 2022 16:56
Operator : SY/MD
Sample : N1025-10
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLH4

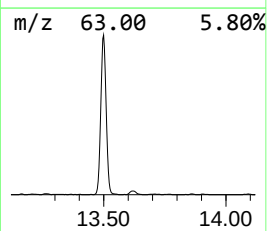
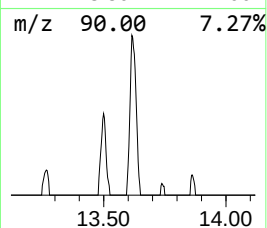
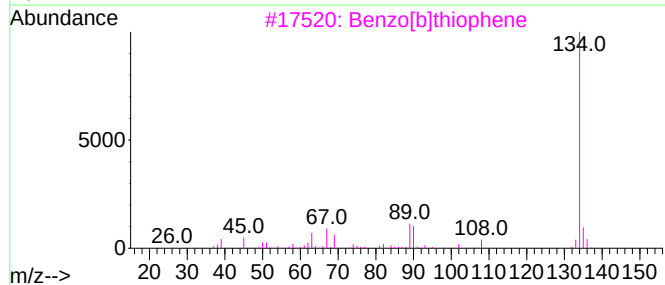
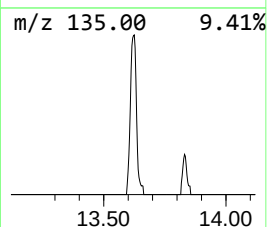
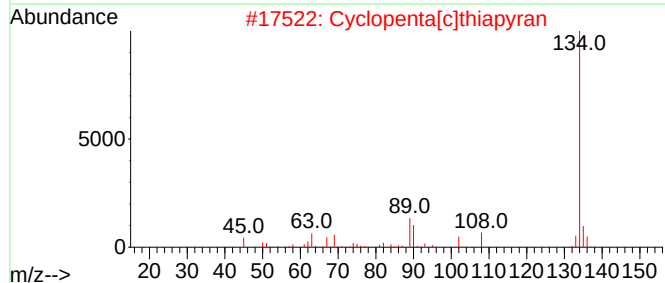
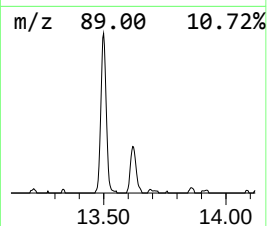
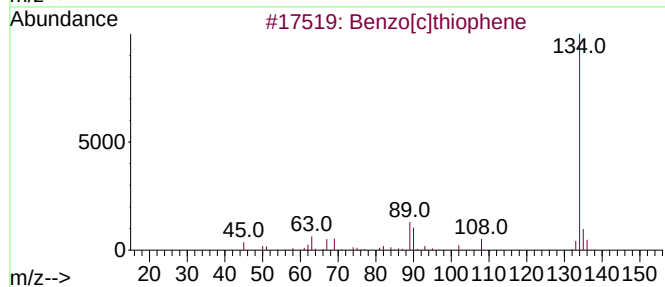
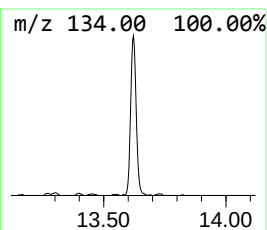
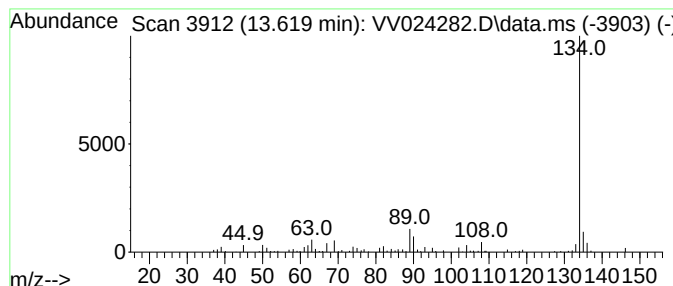
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Quant Title : VOC Analysis

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

Peak Number 24 Benzo[c]thiophene Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.619	5.59 ug/L	93793	1,4-Dichlorobenzene-d4	11.246

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[c]thiophene	134	C8H6S	000270-82-6	95
2			Cyclopenta[c]thiapyran	134	C8H6S	000270-63-3	94
3			Benzo[b]thiophene	134	C8H6S	000095-15-8	91
4			Benzo[b]thiophene	134	C8H6S	000095-15-8	87
5			Benzo[b]thiophene	134	C8H6S	000095-15-8	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV010522\
Data File : VV024282.D
Acq On : 05 Jan 2022 16:56
Operator : SY/MD
Sample : N1025-10
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

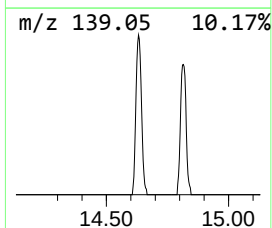
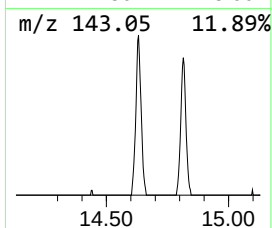
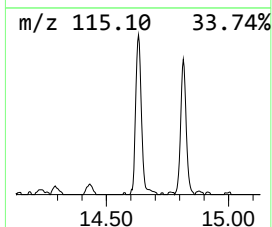
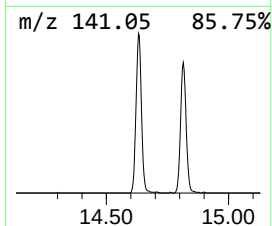
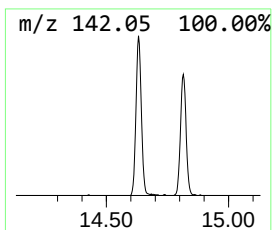
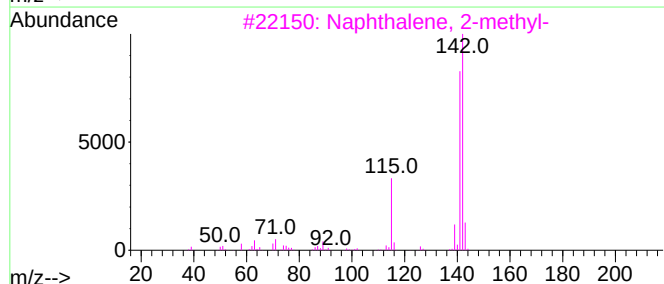
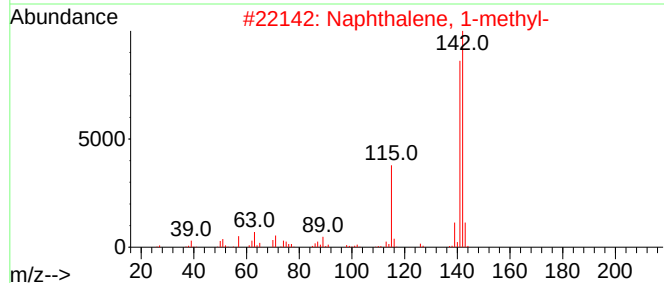
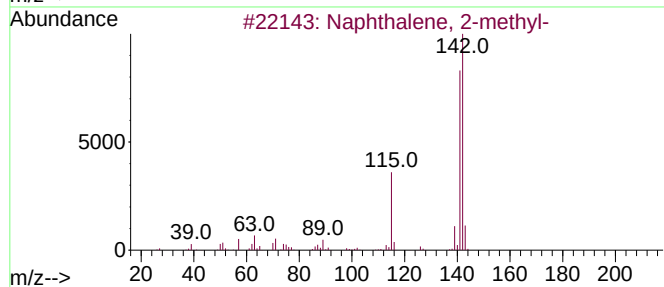
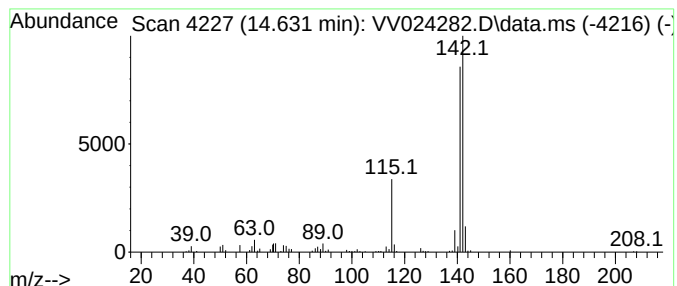
Instrument :
MSVOA_V
ClientSampleId :
BGLH4

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM123121WMA.M
Quant Title : VOC Analysis

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

Peak Number 25 Naphthalene, 2-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.631	8.93 ug/L	149782	1,4-Dichlorobenzene-d4	11.246	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
4	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	95
5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV010522\
Data File : VV024282.D
Acq On : 05 Jan 2022 16:56
Operator : SY/MD
Sample : N1025-10
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLH4

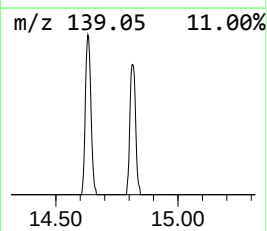
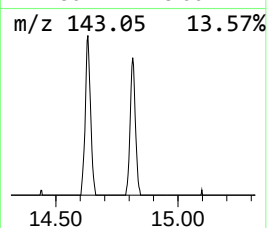
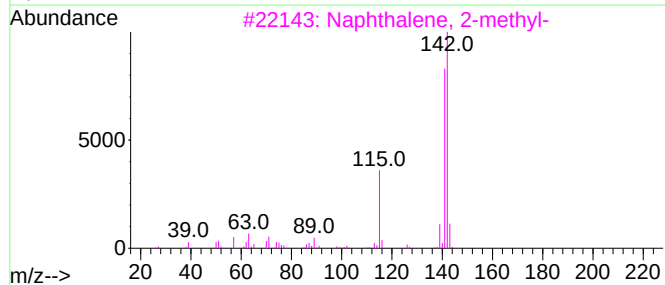
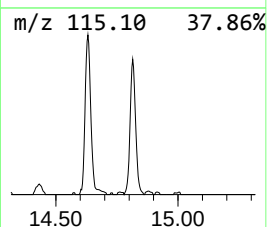
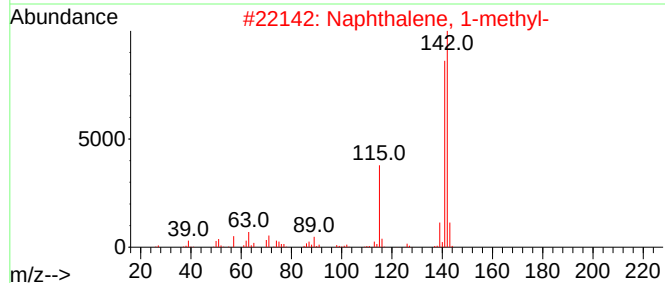
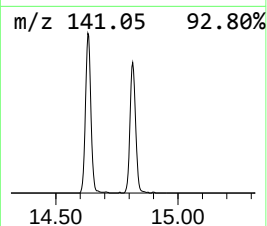
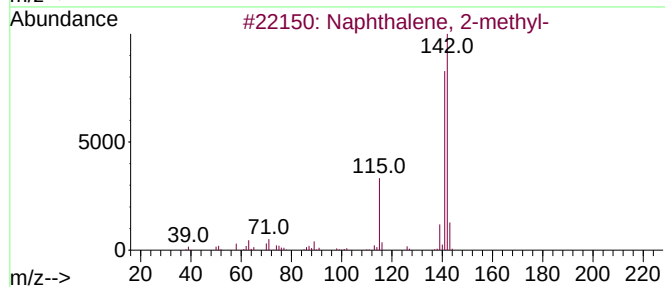
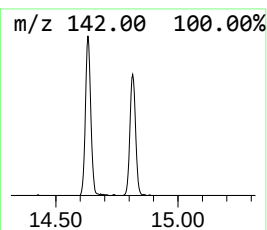
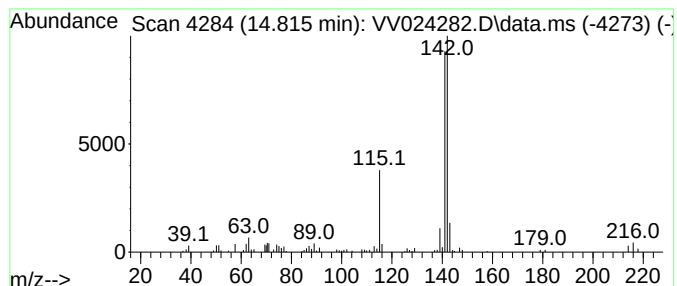
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM123121WMA.M
Quant Title : VOC Analysis

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

Peak Number 26 Naphthalene, 1-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.815	7.89 ug/L	132317	1,4-Dichlorobenzene-d4	11.246

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
4			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
5			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV010522\
 Data File : VV024282.D
 Acq On : 05 Jan 2022 16:56
 Operator : SY/MD
 Sample : N1025-10
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 BGLH4

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM123121WMA.M
 Quant Title : VOC Analysis

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethyl ether	1.950	119.5	ug/L	1404230	1	5.612	587700	50.0
2-Butanone, 3-m...	5.522	5.3	ug/L	61697	1	5.612	587700	50.0
2-Hexanone, 5-m...	9.236	18.0	ug/L	326357	2	8.850	907271	50.0
2-Hexanol, 5-me...	9.378	3.8	ug/L	68148	2	8.850	907271	50.0
1,5-Cyclooctadiene	10.001	15.2	ug/L	276172	2	8.850	907271	50.0
Benzene, propyl-	10.352	25.2	ug/L	423107	3	11.246	838373	50.0
Benzene, 1-ethy...	10.445	39.2	ug/L	656789	3	11.246	838373	50.0
Benzene, 1-ethy...	10.734	38.9	ug/L	652057	3	11.246	838373	50.0
Hydratropaldehyde	11.085	2.7	ug/L	44619	3	11.246	838373	50.0
Benzene, 1,2,3-...	11.333	44.6	ug/L	747158	3	11.246	838373	50.0
Indane	11.525	28.2	ug/L	472143	3	11.246	838373	50.0
Indene	11.744	9.9	ug/L	166557	3	11.246	838373	50.0
Benzene, 1-meth...	11.818	3.4	ug/L	57697	3	11.246	838373	50.0
Benzene, 2-ethy...	11.908	6.9	ug/L	116355	3	11.246	838373	50.0
p-Cymene	12.014	10.4	ug/L	174735	3	11.246	838373	50.0
1-Methyl-2-phen...	12.114	4.6	ug/L	77531	3	11.246	838373	50.0
Benzene, 1,2,4,...	12.413	4.3	ug/L	71794	3	11.246	838373	50.0
Benzene, 1,2,3,...	12.464	9.5	ug/L	159766	3	11.246	838373	50.0
Benzene, 1-ethe...	12.725	3.4	ug/L	57704	3	11.246	838373	50.0
Benzene, 2-ethe...	12.879	14.2	ug/L	237946	3	11.246	838373	50.0
Naphthalene, 1,...	13.046	4.6	ug/L	77040	3	11.246	838373	50.0
Azulene	13.500	202.7	ug/L	3399120	3	11.246	838373	50.0
Benzo[c]thiophene	13.619	5.6	ug/L	93793	3	11.246	838373	50.0
Naphthalene, 2-...	14.631	8.9	ug/L	149782	3	11.246	838373	50.0
Naphthalene, 1-...	14.815	7.9	ug/L	132317	3	11.246	838373	50.0