

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041122\
 Data File : VU047933.D
 Acq On : 12 Apr 2022 02:05
 Operator : SY/MD
 Sample : N2293-08ME
 Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 ETNN8ME

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\SFAMULM032422WMA.M
 Title : VOC Analysis

Signal : TIC: VU047933.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.356	3	5	24	rVB	17530	19354	0.17%	0.049%
2	1.597	73	80	117	rVB	126628	245267	2.15%	0.624%
3	1.903	170	175	203	rVB	80255	159231	1.39%	0.405%
4	2.546	363	375	404	rBV	328106	599655	5.24%	1.525%
5	2.777	439	447	461	rVB	5705	11050	0.10%	0.028%
6	2.922	486	492	494	rBV3	412	406	0.00%	0.001%
7	2.964	496	505	520	rBV2	6941	16249	0.14%	0.041%
8	3.215	579	583	584	rBV	482	353	0.00%	0.001%
9	3.346	619	624	628	rBV3	1144	1141	0.01%	0.003%
10	3.433	640	651	662	rVV5	1580	3191	0.03%	0.008%
11	3.681	723	728	740	rVB6	1590	2934	0.03%	0.007%
12	3.787	754	761	762	rBV2	459	540	0.00%	0.001%
13	3.829	768	774	777	rBV6	688	775	0.01%	0.002%
14	4.063	844	847	850	rVB3	452	363	0.00%	0.001%
15	4.237	895	901	907	rBV4	692	948	0.01%	0.002%
16	4.285	912	916	921	rBV2	452	515	0.00%	0.001%
17	4.517	980	988	990	rBV3	1046	1265	0.01%	0.003%
18	4.642	1012	1027	1082	rBV	139205	536817	4.69%	1.365%
19	4.980	1126	1132	1133	rBV	377	343	0.00%	0.001%
20	5.063	1140	1158	1183	rBV	330760	748478	6.55%	1.903%
21	5.163	1186	1189	1190	rBV2	773	380	0.00%	0.001%
22	5.243	1210	1214	1218	rBV5	837	1048	0.01%	0.003%
23	5.369	1241	1253	1256	rBV10	3135	5312	0.05%	0.014%
24	5.594	1318	1323	1328	rVB6	1088	1251	0.01%	0.003%
25	5.719	1340	1362	1373	rBV2	719820	1925140	16.84%	4.896%
26	5.893	1405	1416	1427	rVB4	7362	15133	0.13%	0.038%
27	5.993	1438	1447	1452	rBV6	1296	2290	0.02%	0.006%
28	6.038	1460	1461	1466	rVB3	1045	444	0.00%	0.001%
29	6.060	1466	1468	1476	rVB4	549	628	0.01%	0.002%
30	6.108	1477	1483	1484	rBV2	973	887	0.01%	0.002%
31	6.128	1484	1489	1497	rVB6	2139	3093	0.03%	0.008%
32	6.247	1512	1526	1549	rBV	691539	1355648	11.86%	3.447%
33	6.407	1574	1576	1584	rVB3	741	692	0.01%	0.002%
34	6.475	1588	1597	1606	rBV4	13659	26223	0.23%	0.067%
35	6.690	1649	1664	1682	rBV	422063	860890	7.53%	2.189%
36	6.899	1725	1729	1733	rBV2	561	378	0.00%	0.001%

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 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 ETNN8ME

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\SFAMULM032422WMA.M
 Title : VOC Analysis

37	6.944	1740	1743	1745	rBV2	440	363	0.00%	0.001%
38	7.057	1770	1778	1783	rBV3	978	1622	0.01%	0.004%
39	7.202	1811	1823	1825	rBV6	6777	11997	0.10%	0.031%
40	7.497	1901	1915	1925	rBV3	21683	47199	0.41%	0.120%
41	7.568	1926	1937	1956	rVB	213915	408517	3.57%	1.039%
42	7.694	1968	1976	1977	rBV6	2055	2435	0.02%	0.006%
43	7.899	2025	2040	2053	rBV	855657	1521187	13.30%	3.868%
44	7.973	2054	2063	2091	rVB2	96796	241657	2.11%	0.615%
45	8.182	2117	2128	2150	rBV	135990	277797	2.43%	0.706%
46	8.369	2174	2186	2206	rVB7	13917	37665	0.33%	0.096%
47	8.475	2216	2219	2224	rVB4	555	433	0.00%	0.001%
48	8.536	2235	2238	2240	rBV2	564	378	0.00%	0.001%
49	8.549	2240	2242	2244	rVV2	777	480	0.00%	0.001%
50	8.581	2244	2252	2260	rVV10	3365	6129	0.05%	0.016%
51	8.661	2260	2277	2307	rVV	580430	1228373	10.74%	3.124%
52	8.922	2353	2358	2359	rBV4	1192	964	0.01%	0.002%
53	9.034	2378	2393	2407	rVB6	10993	27515	0.24%	0.070%
54	9.118	2413	2419	2420	rBV3	541	481	0.00%	0.001%
55	9.192	2429	2442	2445	rBV5	1448	2740	0.02%	0.007%
56	9.240	2454	2457	2459	rBV2	593	435	0.00%	0.001%
57	9.365	2483	2496	2501	rBV4	9171	15154	0.13%	0.039%
58	9.452	2501	2523	2554	rVV2	5362170	11434277	100.00%	29.078%
59	9.581	2555	2563	2574	rVB4	18854	33144	0.29%	0.084%
60	9.697	2587	2599	2609	rBV2	18289	33766	0.30%	0.086%
61	9.944	2674	2676	2683	rVB5	797	843	0.01%	0.002%
62	10.015	2689	2698	2701	rBV6	2306	3250	0.03%	0.008%
63	10.108	2716	2727	2745	rVB2	14283	28718	0.25%	0.073%
64	10.240	2757	2768	2778	rBV9	5434	12248	0.11%	0.031%
65	10.333	2790	2797	2798	rBV5	1392	1222	0.01%	0.003%
66	10.353	2798	2803	2811	rVB8	2392	3785	0.03%	0.010%
67	10.475	2834	2841	2855	rVB8	4508	10178	0.09%	0.026%
68	10.552	2863	2865	2871	rVB4	936	724	0.01%	0.002%
69	10.658	2889	2898	2908	rVV8	14982	30736	0.27%	0.078%
70	10.716	2908	2916	2920	rVV4	12536	21853	0.19%	0.056%
71	10.764	2920	2931	2953	rVB	712320	1182584	10.34%	3.007%
72	10.983	2990	2999	3019	rVB7	17017	42838	0.37%	0.109%
73	11.066	3021	3025	3026	rVV3	1671	1193	0.01%	0.003%
74	11.108	3028	3038	3051	rVB5	14497	31454	0.28%	0.080%
75	11.250	3079	3082	3083	rBV3	1316	731	0.01%	0.002%
76	11.295	3092	3096	3109	rVB5	5017	8332	0.07%	0.021%

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 Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 ETNN8ME

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\SFAMULM032422WMA.M
 Title : VOC Analysis

77	11.472	3141	3151	3160	rBV4	18759	33326	0.29%	0.085%
78	11.584	3177	3186	3198	rBV7	7563	14763	0.13%	0.038%
79	11.645	3202	3205	3206	rBV3	1387	817	0.01%	0.002%
80	11.693	3215	3220	3221	rBV3	3439	2375	0.02%	0.006%
81	11.751	3231	3238	3245	rVB3	13281	18382	0.16%	0.047%
82	11.816	3245	3258	3273	rBV	1311602	2094183	18.31%	5.326%
83	11.954	3294	3301	3314	rVB5	21355	35989	0.31%	0.092%
84	12.053	3329	3332	3333	rBV2	924	546	0.00%	0.001%
85	12.198	3364	3377	3398	rBV2	1036824	2037823	17.82%	5.182%
86	12.327	3411	3417	3427	rVB	23350	32833	0.29%	0.083%
87	12.970	3606	3617	3630	rVB	795566	1241156	10.85%	3.156%
88	13.330	3718	3729	3737	rBV	648346	1007301	8.81%	2.562%
89	13.378	3738	3744	3754	rVB	209429	298439	2.61%	0.759%
90	14.044	3944	3951	3959	rBV2	43986	63566	0.56%	0.162%
91	14.359	4041	4049	4064	rVB	136351	201739	1.76%	0.513%
92	15.118	4277	4285	4291	rBV4	35786	58522	0.51%	0.149%
93	15.311	4339	4345	4355	rVB	91817	128045	1.12%	0.326%
94	15.545	4407	4418	4433	rVB	695089	1062551	9.29%	2.702%
95	15.796	4486	4496	4507	rBV	176911	270068	2.36%	0.687%
96	15.893	4511	4526	4540	rBV3	211693	578553	5.06%	1.471%
97	16.092	4581	4588	4600	rVB2	45317	68868	0.60%	0.175%
98	16.391	4669	4681	4706	rVB	4223704	6547659	57.26%	16.651%
99	17.031	4872	4880	4891	rVB2	91725	140436	1.23%	0.357%
100	17.259	4944	4951	4967	rVB2	76284	121722	1.06%	0.310%

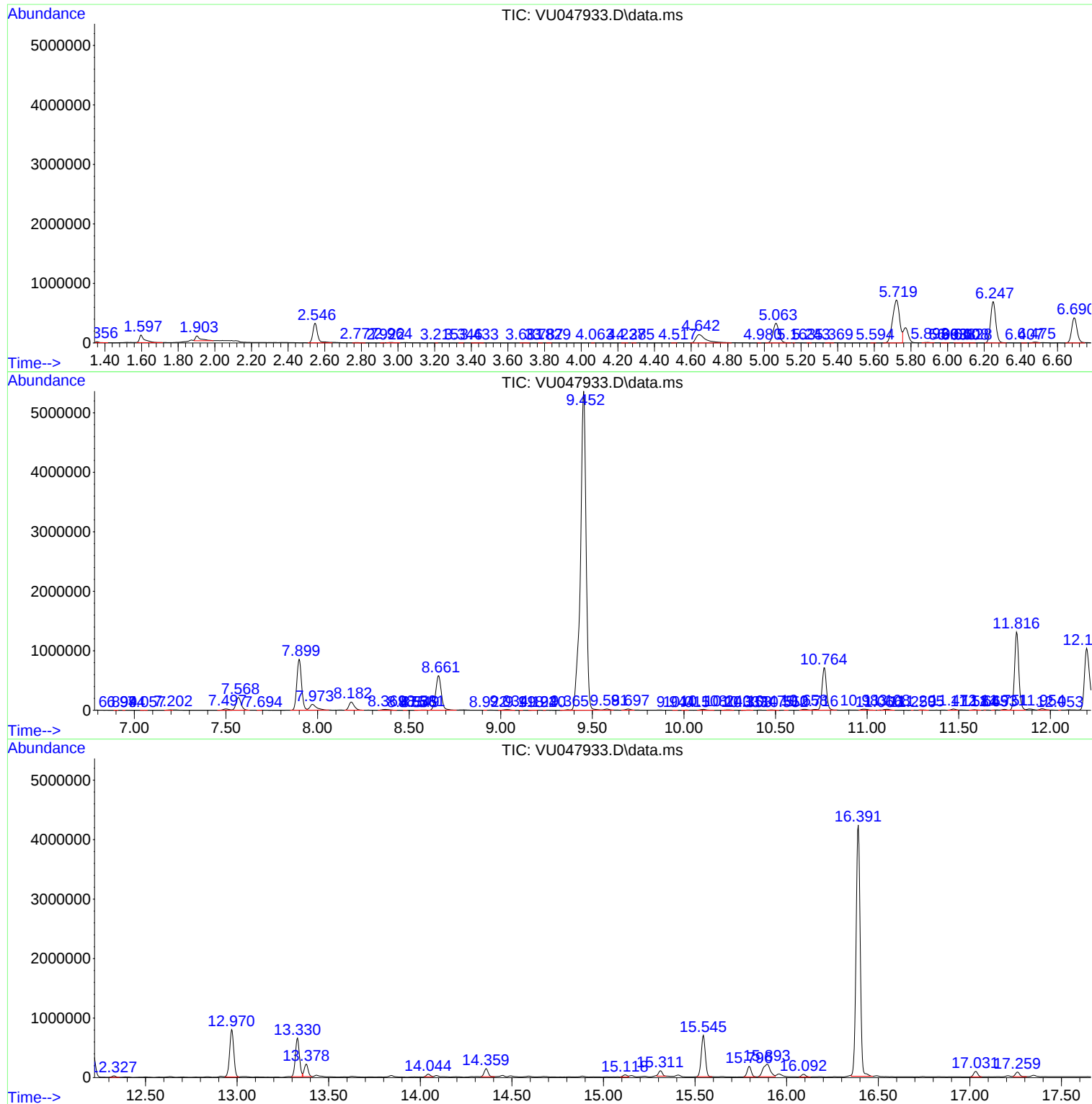
Sum of corrected areas: 39323371

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041122\
Data File : VU047933.D
Acq On : 12 Apr 2022 02:05
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Sample : N2293-08ME
Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 41 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
ETNN8ME

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM032422WMA.M
Quant Title : VOC Analysis

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041122\
Data File : VU047933.D
Acq On : 12 Apr 2022 02:05
Operator : SY/MD
Sample : N2293-08ME
Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 41 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
ETNN8ME

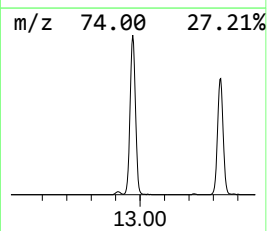
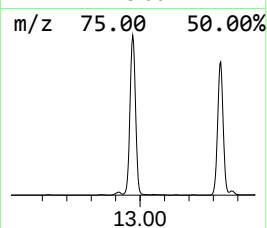
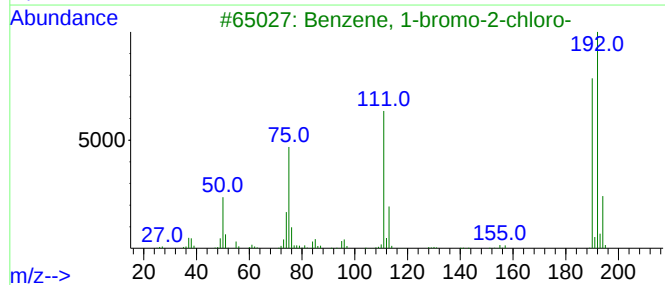
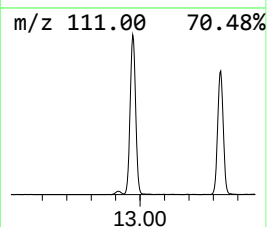
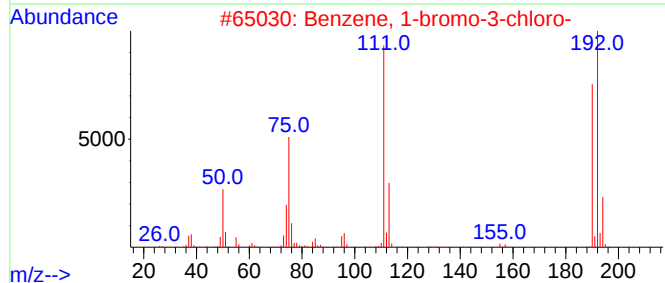
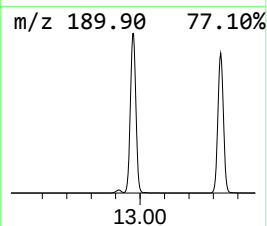
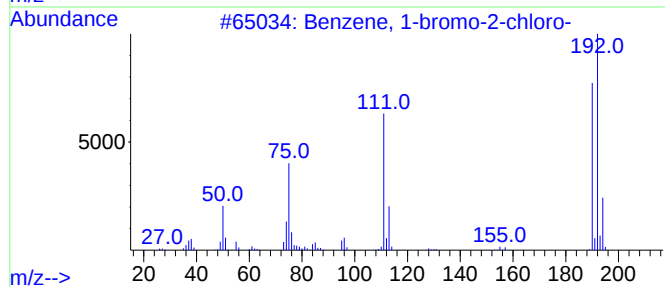
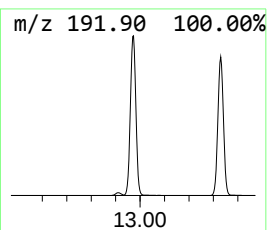
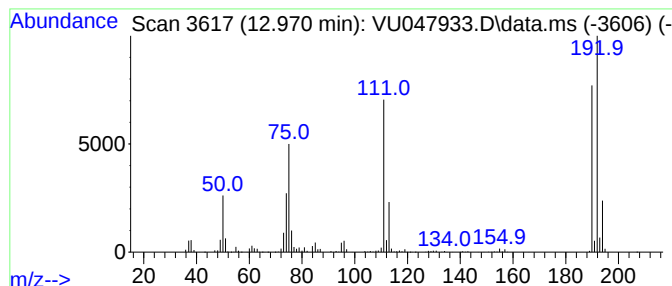
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM032422WMA.M
Quant Title : VOC Analysis

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

Peak Number 2 Benzene, 1-bromo-2-chloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.970	29.63 ug/L	1241160	1,4-Dichlorobenzene-d4	11.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	99
2			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	98
3			Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	98
4			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	98
5			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98



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Sample : N2293-08ME
Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 41 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
ETNN8ME

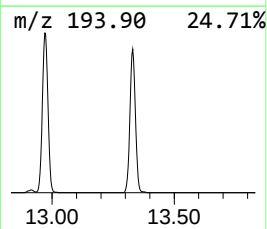
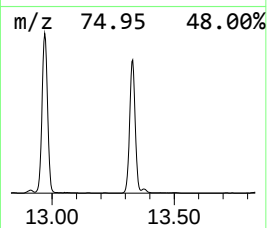
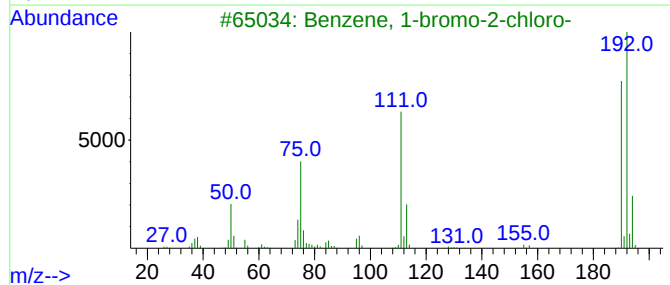
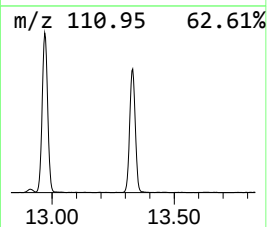
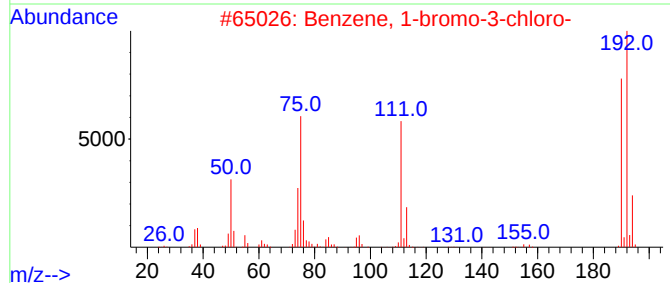
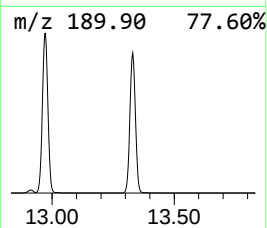
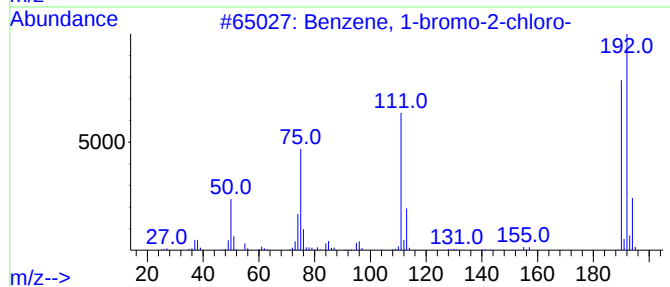
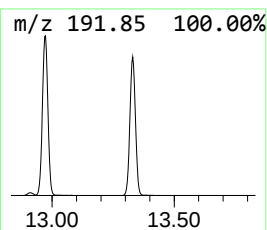
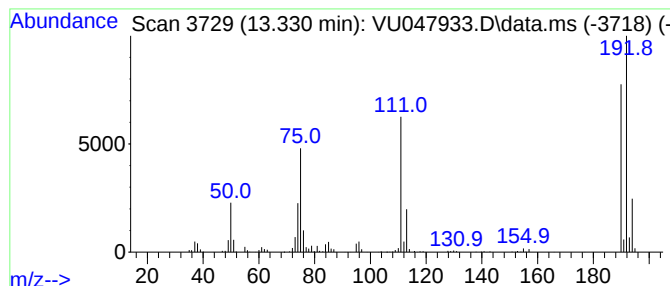
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM032422WMA.M
Quant Title : VOC Analysis

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

Peak Number 3 Benzene, 1-bromo-3-chloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.330	24.05 ug/L	1007300	1,4-Dichlorobenzene-d4	11.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	99
2			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	99
3			Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	99
4			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98
5			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98



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Operator : SY/MD
Sample : N2293-08ME
Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 41 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
ETNN8ME

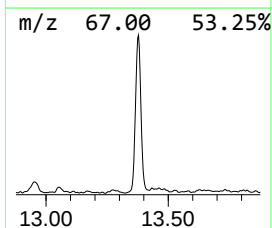
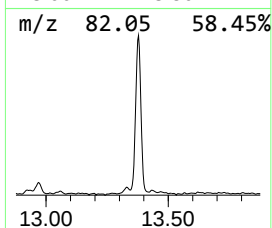
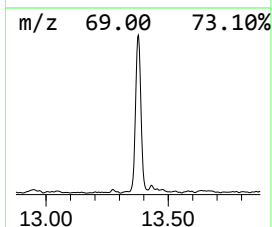
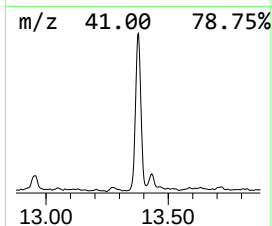
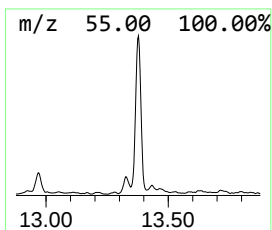
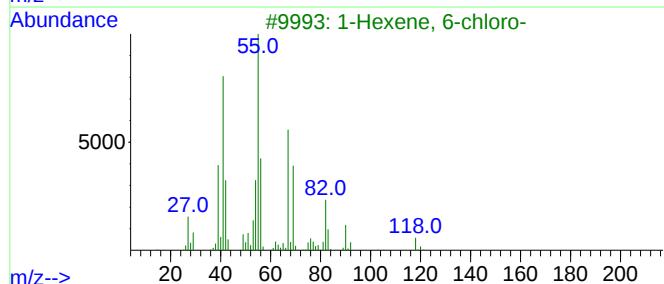
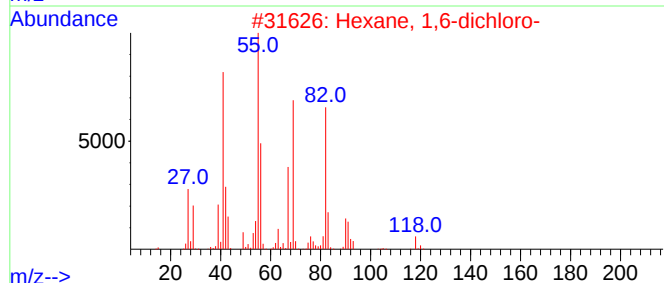
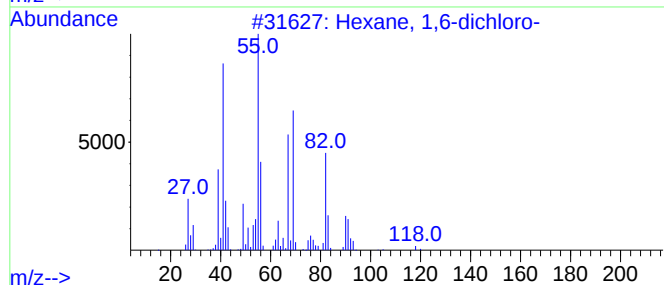
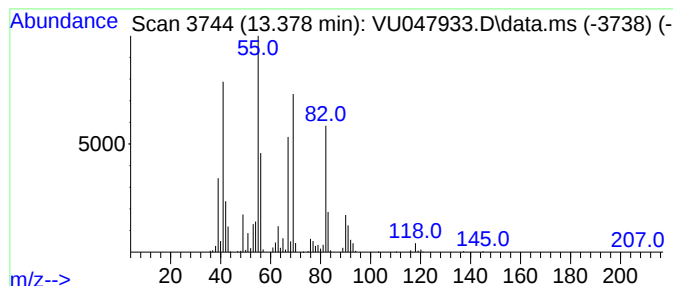
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Quant Title : VOC Analysis

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

Peak Number 4 Hexane, 1,6-dichloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.378	7.13 ug/L	298439	1,4-Dichlorobenzene-d4	11.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 1,6-dichloro-	154	C6H12Cl2	002163-00-0	80
2			Hexane, 1,6-dichloro-	154	C6H12Cl2	002163-00-0	72
3			1-Hexene, 6-chloro-	118	C6H11Cl	000928-89-2	58
4			4-Methyl-1,3-pentadiene	82	C6H10	000926-56-7	38
5			2H-Pyran, 2-(4-chlorobutoxy)tetr...	192	C9H17ClO2	041302-05-0	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041122\
Data File : VU047933.D
Acq On : 12 Apr 2022 02:05
Operator : SY/MD
Sample : N2293-08ME
Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 41 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
ETNN8ME

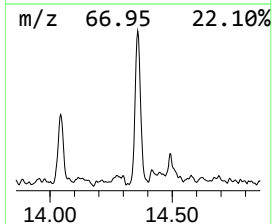
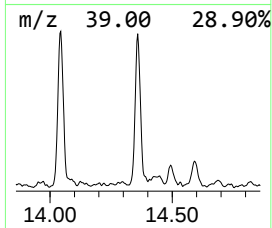
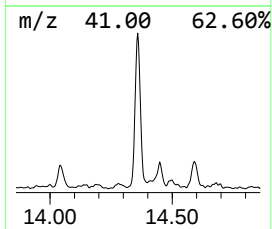
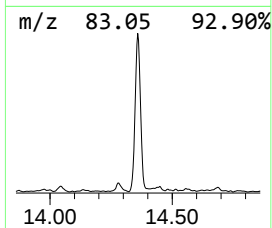
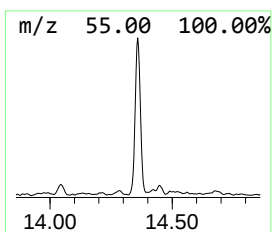
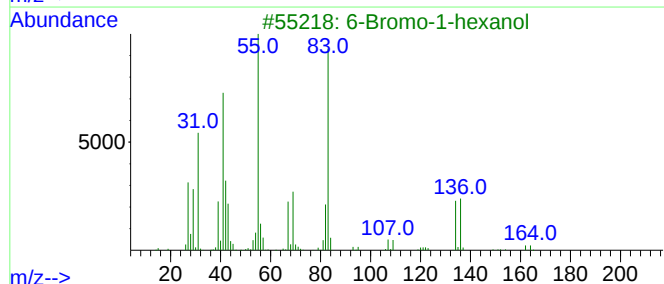
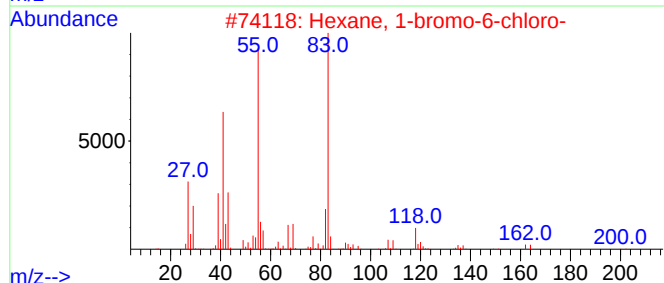
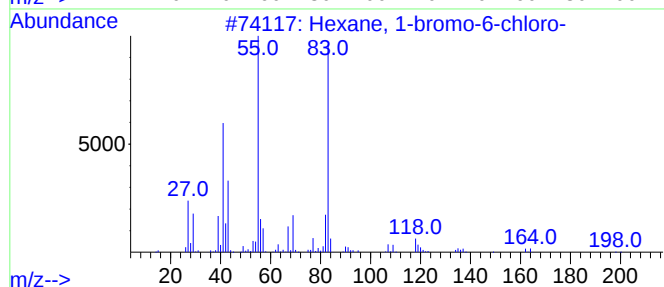
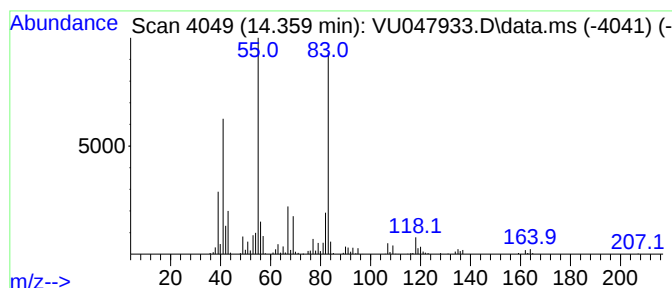
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Quant Title : VOC Analysis

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

Peak Number 5 Hexane, 1-bromo-6-chloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.359	4.82 ug/L	201739	1,4-Dichlorobenzene-d4	11.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 1-bromo-6-chloro-	198	C6H12BrCl	006294-17-3	80
2			Hexane, 1-bromo-6-chloro-	198	C6H12BrCl	006294-17-3	72
3			6-Bromo-1-hexanol	180	C6H13BrO	004286-55-9	64
4			Cyclopentane, 1-ethyl-1-methyl-	112	C8H16	016747-50-5	53
5			Oxalic acid, cyclohexyl isohexyl...	256	C14H24O4	1010309-30-7	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041122\
Data File : VU047933.D
Acq On : 12 Apr 2022 02:05
Operator : SY/MD
Sample : N2293-08ME
Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 41 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
ETNN8ME

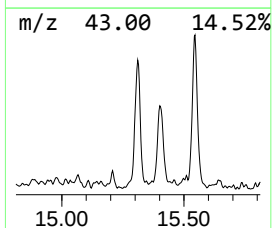
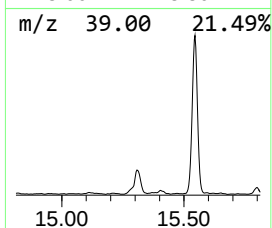
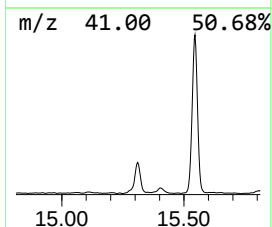
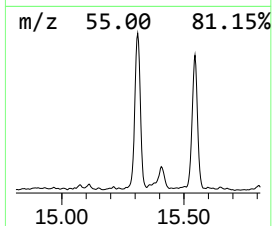
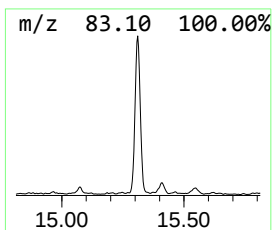
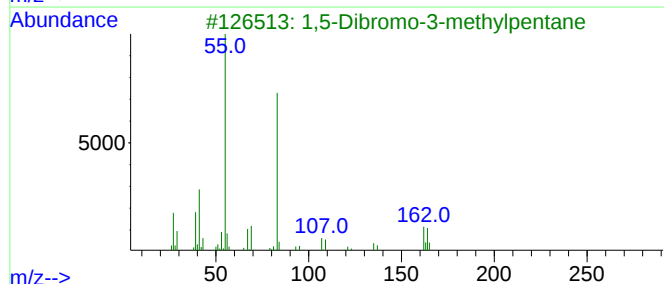
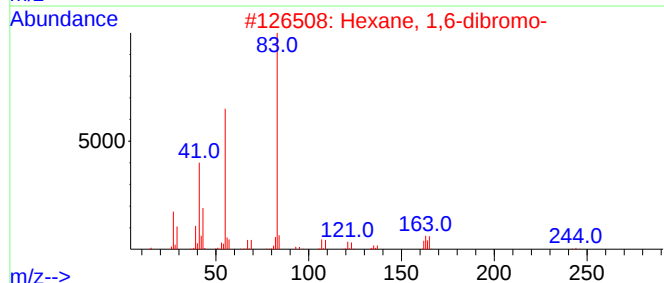
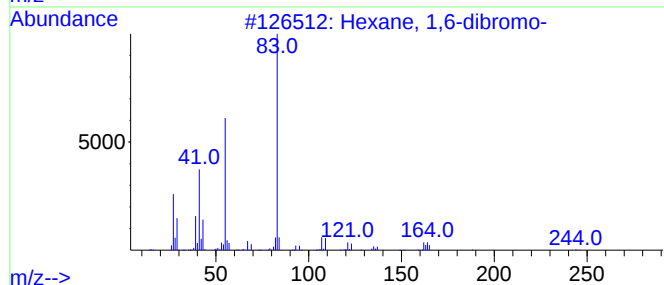
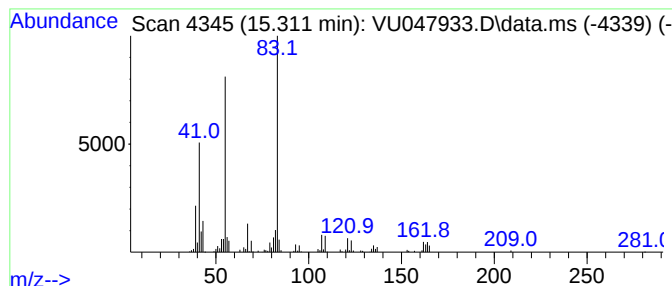
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Quant Title : VOC Analysis

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

Peak Number 6 Hexane, 1,6-dibromo- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.311	3.06 ug/L	128045	1,4-Dichlorobenzene-d4	11.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 1,6-dibromo-	242	C6H12Br2	000629-03-8	91
2			Hexane, 1,6-dibromo-	242	C6H12Br2	000629-03-8	83
3			1,5-Dibromo-3-methylpentane	242	C6H12Br2	004457-72-1	83
4			Hexane, 2,5-dibromo-	242	C6H12Br2	024774-58-1	78
5			Cyclohexane, bromo-	162	C6H11Br	000108-85-0	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041122\
Data File : VU047933.D
Acq On : 12 Apr 2022 02:05
Operator : SY/MD
Sample : N2293-08ME
Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 41 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
ETNN8ME

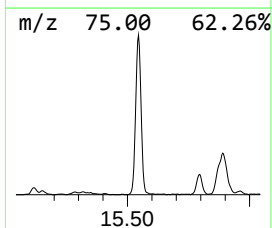
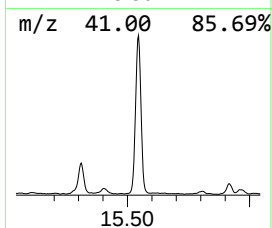
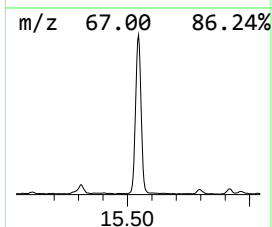
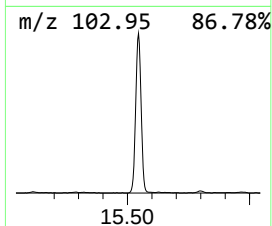
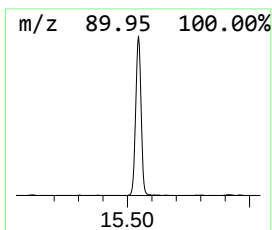
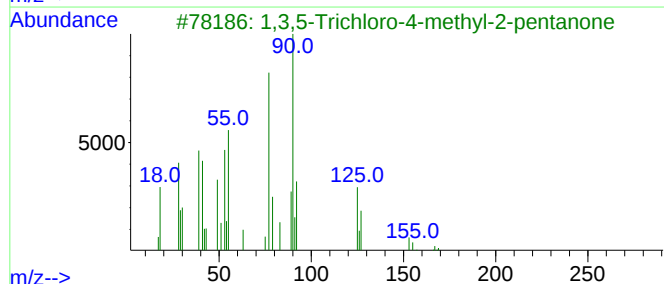
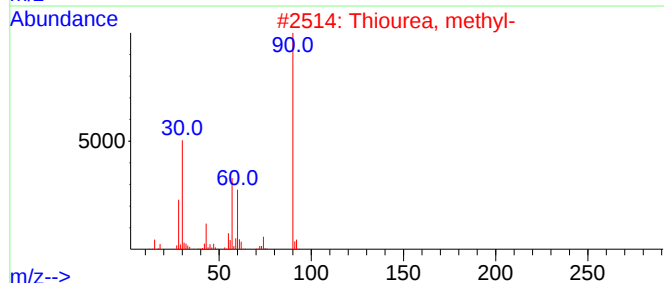
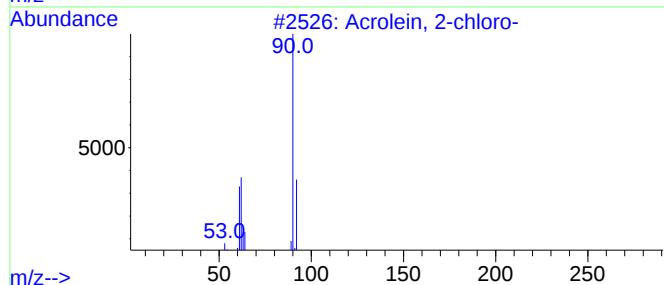
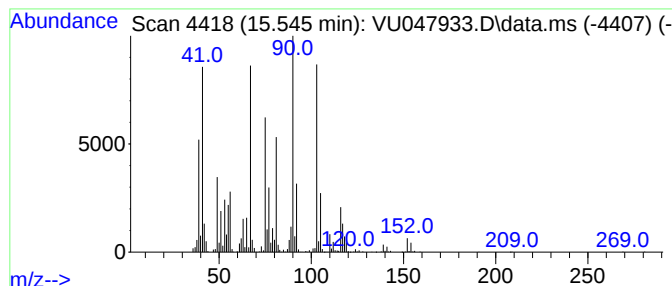
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Quant Title : VOC Analysis

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

Peak Number 7 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.545	25.37 ug/L	1062550	1,4-Dichlorobenzene-d4	11.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acrolein, 2-chloro-	90	C3H3ClO	000683-51-2	37
2			Thiourea, methyl-	90	C2H6N2S	000598-52-7	35
3			1,3,5-Trichloro-4-methyl-2-penta...	202	C6H9Cl3O	1000140-03-8	25
4			Propane, 2-(methylthio)-	90	C4H10S	001551-21-9	25
5			2,2-Bis(chloromethyl)-1-propanol	156	C5H10Cl2O	005355-54-4	23



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041122\
Data File : VU047933.D
Acq On : 12 Apr 2022 02:05
Operator : SY/MD
Sample : N2293-08ME
Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 41 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
ETNN8ME

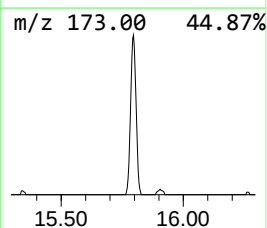
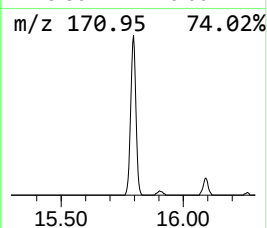
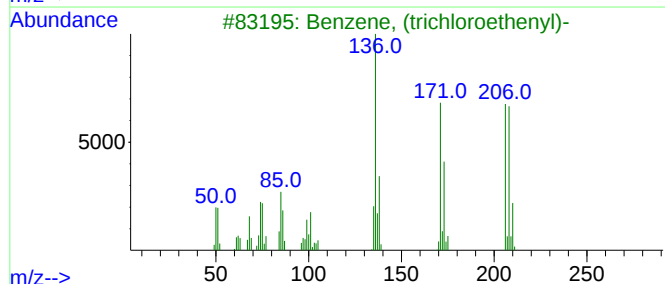
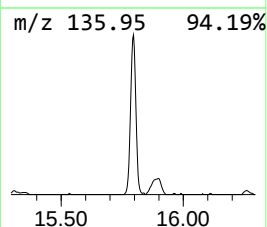
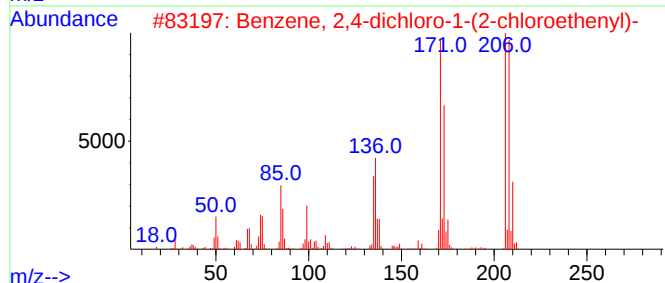
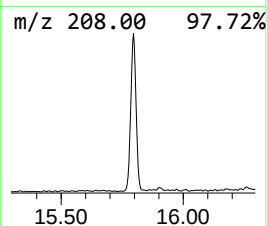
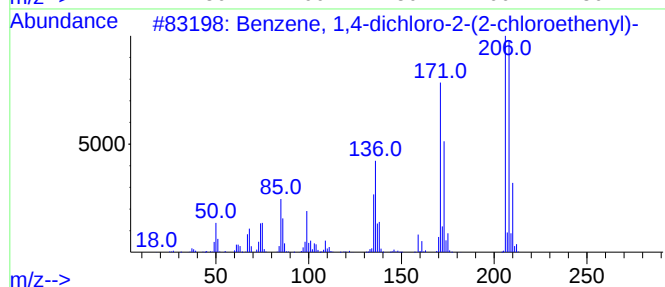
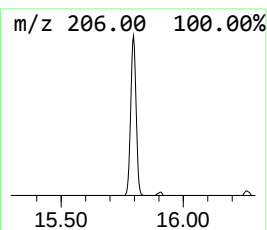
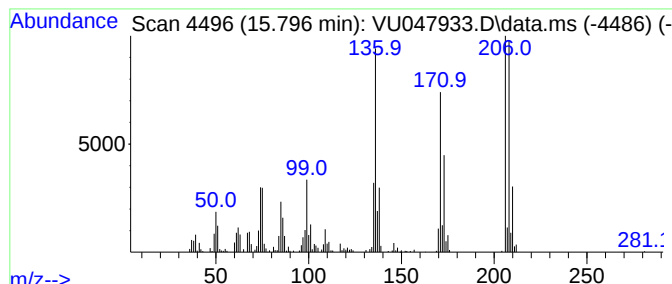
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Quant Title : VOC Analysis

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

Peak Number 8 Benzene, 1,4-dichloro-2-(2-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.796	6.45 ug/L	270068	1,4-Dichlorobenzene-d4	11.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-dichloro-2-(2-chlor...	206	C8H5Cl3	054935-00-1	98
2			Benzene, 2,4-dichloro-1-(2-chlor...	206	C8H5Cl3	045892-47-5	96
3			Benzene, (trichloroethenyl)-	206	C8H5Cl3	000700-60-7	93
4			1,1-Dichloro-2-(P-chlorophenyl)e...	206	C8H5Cl3	005263-17-2	53
5			Phenol, 4-bromo-2-chloro-	206	C6H4BrClO	003964-56-5	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041122\
Data File : VU047933.D
Acq On : 12 Apr 2022 02:05
Operator : SY/MD
Sample : N2293-08ME
Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 41 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
ETNN8ME

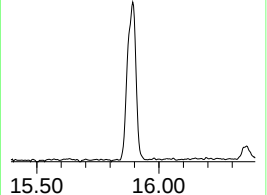
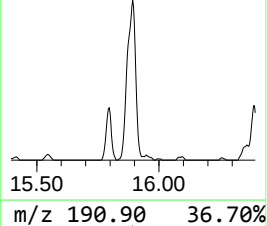
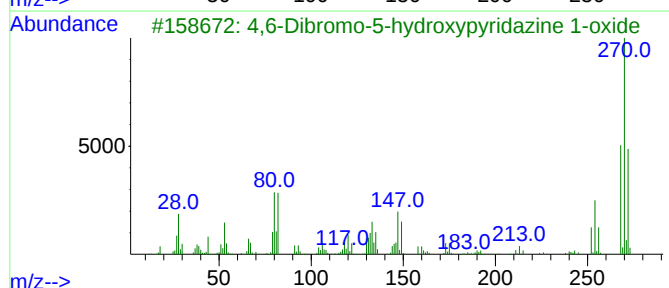
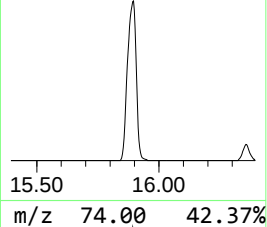
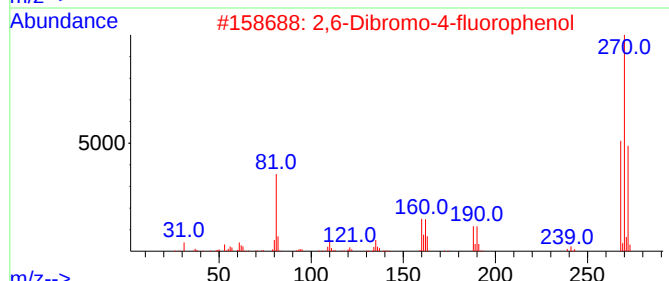
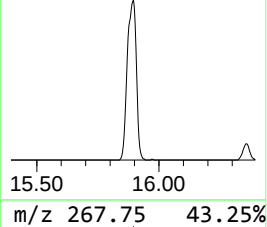
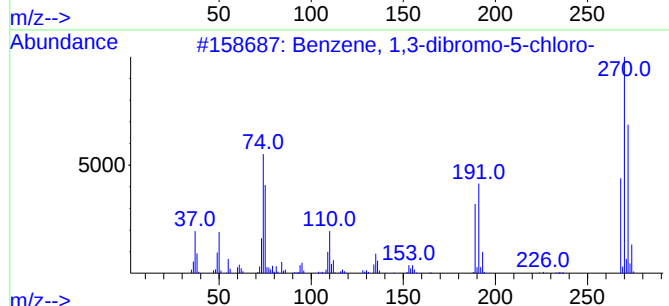
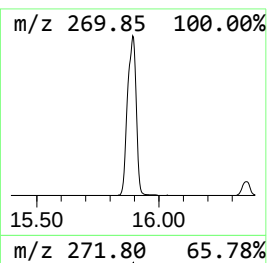
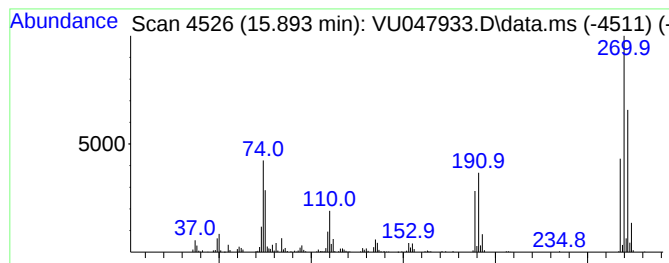
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Quant Title : VOC Analysis

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

Peak Number 9 Benzene, 1,3-dibromo-5-chloro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.893	13.81 ug/L	578553	1,4-Dichlorobenzene-d4	11.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dibromo-5-chloro-	268	C6H3Br2Cl	014862-52-3	94
2			2,6-Dibromo-4-fluorophenol	268	C6H3Br2FO	000344-20-7	38
3			4,6-Dibromo-5-hydroxypyridazine ...	268	C4H2Br2N2O2	019971-61-0	37
4			Pyrimidine, 4,6-dichloro-2-(meth...	270	C11H8Cl2N2S	064415-11-8	23
5			5,7-Dichloro-8-[(trimethylsilyl)...	285	C12H13Cl2NOSi	1000408-12-3	13



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041122\
Data File : VU047933.D
Acq On : 12 Apr 2022 02:05
Operator : SY/MD
Sample : N2293-08ME
Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 41 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
ETNN8ME

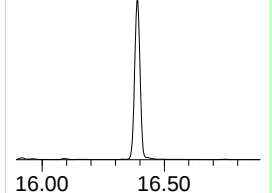
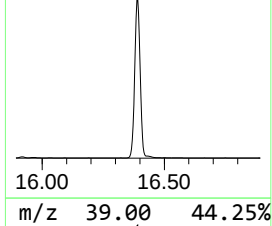
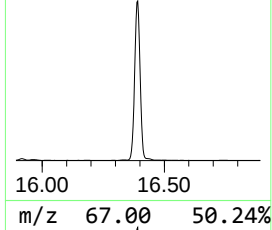
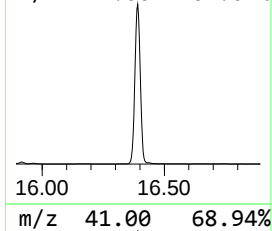
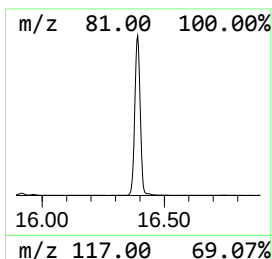
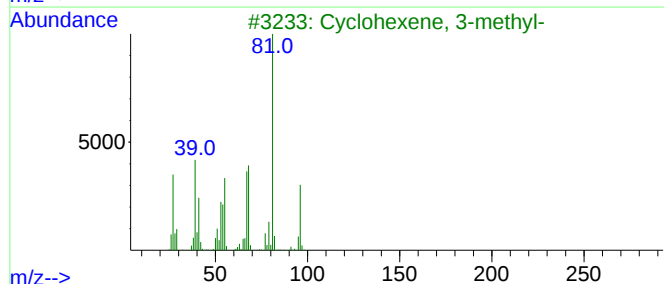
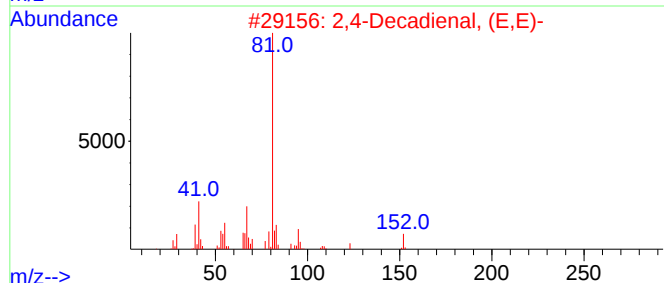
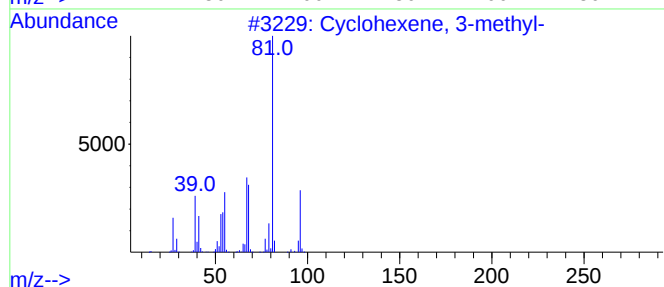
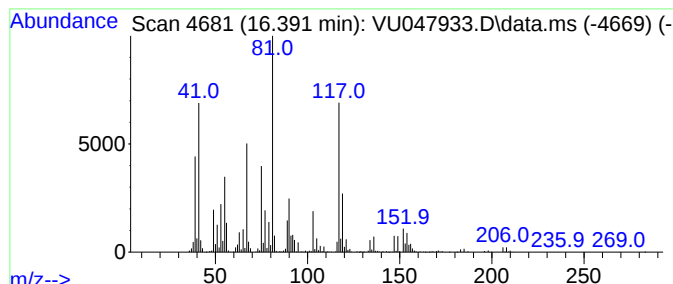
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Quant Title : VOC Analysis

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

Peak Number 10 unknown-02 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.391	156.33 ug/L	6547660	1,4-Dichlorobenzene-d4	11.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexene, 3-methyl-	96	C7H12	000591-48-0	22
2			2,4-Decadienal, (E,E)-	152	C10H16O	025152-84-5	22
3			Cyclohexene, 3-methyl-	96	C7H12	000591-48-0	22
4			(9-Oxabicyclo[3.3.1]non-6-en-3-y...	154	C9H14O2	1000191-02-8	12
5			2-Hexyne, 4-methyl-	96	C7H12	020198-49-6	12



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041122\
Data File : VU047933.D
Acq On : 12 Apr 2022 02:05
Operator : SY/MD
Sample : N2293-08ME
Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 41 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
ETNN8ME

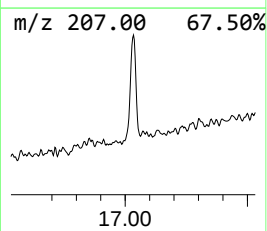
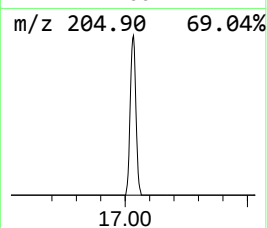
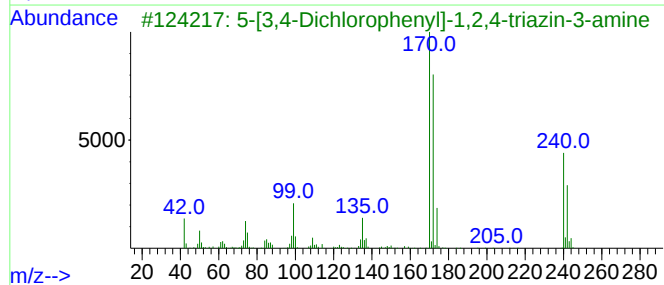
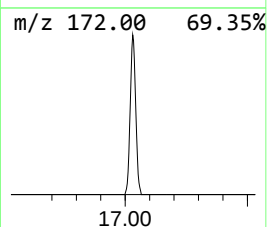
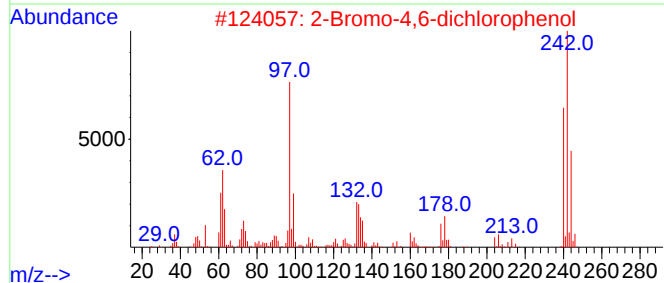
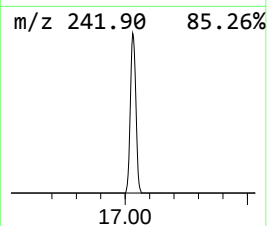
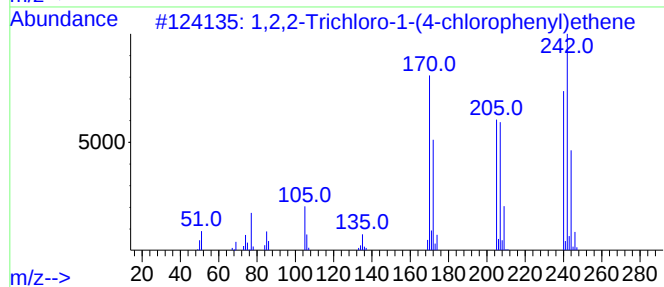
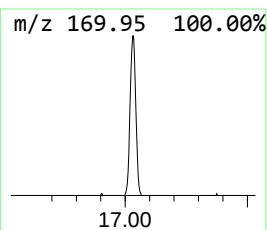
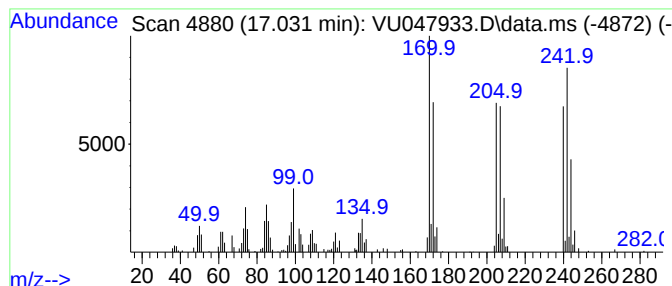
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Quant Title : VOC Analysis

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

Peak Number 11 1,2,2-Trichloro-1-(4-chloro... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.031	3.35 ug/L	140436	1,4-Dichlorobenzene-d4	11.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2,2-Trichloro-1-(4-chloropheny...	240	C8H4Cl4	004714-35-6	99
2			2-Bromo-4,6-dichlorophenol	240	C6H3BrCl2O	004524-77-0	50
3			5-[3,4-Dichlorophenyl]-1,2,4-tri...	240	C9H6Cl2N4	1000213-66-2	46
4			2-Chloro-4-fluoro-5-methylbenzen...	242	C7H5Cl2F02S	1000443-97-7	46
5			3-Amino-6-bromo-2,4-dichloropyri...	240	C5H3BrCl2N2	237435-16-4	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041122\
Data File : VU047933.D
Acq On : 12 Apr 2022 02:05
Operator : SY/MD
Sample : N2293-08ME
Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 41 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
ETNN8ME

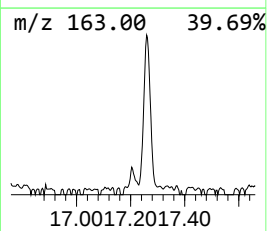
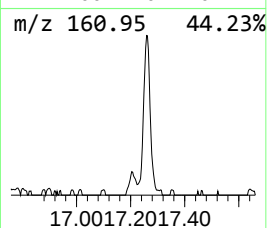
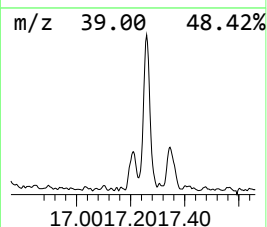
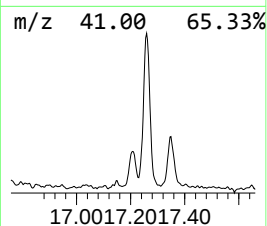
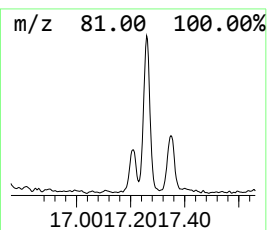
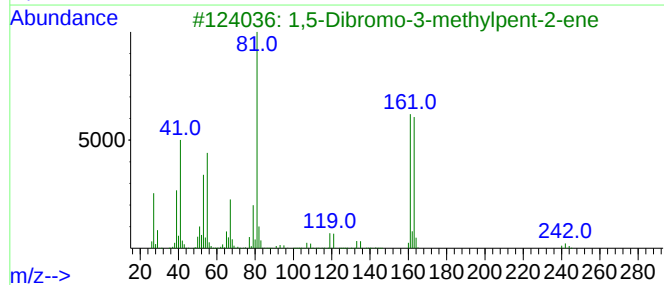
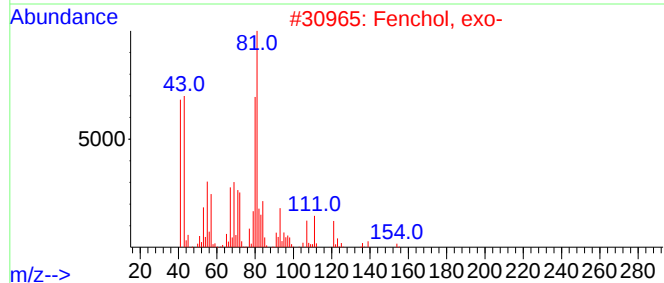
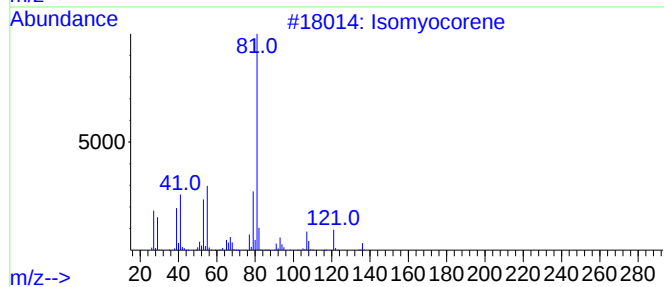
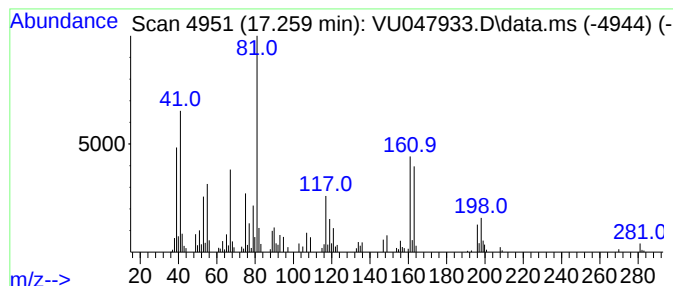
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Quant Title : VOC Analysis

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

Peak Number 12 unknown-03 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.259	2.91 ug/L	121722	1,4-Dichlorobenzene-d4	11.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isomycorene	136	C10H16	006876-07-9	38
2			Fenchol, exo-	154	C10H18O	022627-95-8	32
3			1,5-Dibromo-3-methylpent-2-ene	240	C6H10Br2	133545-67-2	27
4			2,3-Diazabicyclo[2.2.1]hept-2-en...	178	C11H18N2	150667-99-5	22
5			Cyclopropane, 1-methyl-1-isoprop...	96	C7H12	003422-07-9	22



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041122\
 Data File : VU047933.D
 Acq On : 12 Apr 2022 02:05
 Operator : SY/MD
 Sample : N2293-08ME
 Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 ETNN8ME

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM032422WMA.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
Benzene, 1-brom...	12.970	29.6	ug/L	1241160	3	11.816	2094180	50.0
Benzene, 1-brom...	13.330	24.1	ug/L	1007300	3	11.816	2094180	50.0
Hexane, 1,6-dic...	13.378	7.1	ug/L	298439	3	11.816	2094180	50.0
Hexane, 1-bromo...	14.359	4.8	ug/L	201739	3	11.816	2094180	50.0
Hexane, 1,6-dib...	15.311	3.1	ug/L	128045	3	11.816	2094180	50.0
unknown-01	15.545	25.4	ug/L	1062550	3	11.816	2094180	50.0
Benzene, 1,4-di...	15.796	6.5	ug/L	270068	3	11.816	2094180	50.0
Benzene, 1,3-di...	15.893	13.8	ug/L	578553	3	11.816	2094180	50.0
unknown-02	16.391	156.3	ug/L	6547660	3	11.816	2094180	50.0
1,2,2-Trichloro...	17.031	3.4	ug/L	140436	3	11.816	2094180	50.0
unknown-03	17.259	2.9	ug/L	121722	3	11.816	2094180	50.0