

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU060922\
 Data File : VU048984.D
 Acq On : 09 Jun 2022 18:29
 Operator : SY/MD
 Sample : N3248-09
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 EXYY0

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\SFAMULM051622WMA.M
 Title : VOC Analysis

Signal : TIC: VU048984.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.601	72	81	98	rBV2	311634	385844	7.34%	1.400%
2	1.922	166	181	222	rVB	1772065	2637689	50.16%	9.569%
3	2.565	367	381	395	rBV2	256521	459035	8.73%	1.665%
4	2.948	494	500	501	rBV2	1160	1060	0.02%	0.004%
5	3.041	522	529	539	rVB3	2228	3486	0.07%	0.013%
6	3.182	561	573	589	rBV	3525	8047	0.15%	0.029%
7	3.350	614	625	644	rVB2	28102	56377	1.07%	0.205%
8	3.867	766	786	814	rBV	1317208	3105629	59.05%	11.267%
9	4.568	998	1004	1007	rBV2	597	555	0.01%	0.002%
10	4.665	1007	1034	1061	rBV2	1463246	3607779	68.60%	13.088%
11	5.063	1140	1158	1182	rVB	227784	496226	9.44%	1.800%
12	5.211	1199	1204	1205	rBV4	660	558	0.01%	0.002%
13	5.314	1215	1236	1269	rBV	2320136	5258889	100.00%	19.078%
14	5.723	1342	1363	1397	rBV2	429105	1207230	22.96%	4.380%
15	5.899	1408	1418	1427	rBV6	3490	7659	0.15%	0.028%
16	5.970	1433	1440	1452	rVB8	1656	3026	0.06%	0.011%
17	6.250	1512	1527	1544	rBV	389185	738191	14.04%	2.678%
18	6.311	1544	1546	1554	rVB4	1274	1258	0.02%	0.005%
19	6.485	1595	1600	1605	rVB4	627	581	0.01%	0.002%
20	6.542	1605	1618	1632	rBV3	53403	103113	1.96%	0.374%
21	6.690	1650	1664	1679	rBV	290318	560247	10.65%	2.032%
22	6.973	1743	1752	1759	rBV5	1692	2460	0.05%	0.009%
23	7.179	1808	1816	1827	rVV6	1695	3802	0.07%	0.014%
24	7.256	1837	1840	1844	rVV4	811	649	0.01%	0.002%
25	7.565	1921	1936	1955	rBV	164669	283710	5.39%	1.029%
26	7.790	1998	2006	2015	rVV4	2020	3082	0.06%	0.011%
27	7.851	2021	2025	2026	rBV2	756	529	0.01%	0.002%
28	7.899	2026	2040	2051	rVV	544605	927148	17.63%	3.363%
29	7.970	2051	2062	2077	rVB	321378	548583	10.43%	1.990%
30	8.179	2112	2127	2142	rBV2	118124	198028	3.77%	0.718%
31	8.366	2181	2185	2190	rBV3	651	481	0.01%	0.002%
32	8.472	2208	2218	2231	rBV6	2492	4942	0.09%	0.018%
33	8.552	2231	2243	2256	rBV	142976	243332	4.63%	0.883%
34	8.632	2256	2268	2302	rBV	445371	802436	15.26%	2.911%
35	8.806	2320	2322	2329	rVB4	1095	994	0.02%	0.004%
36	8.861	2332	2339	2347	rVB5	2109	3174	0.06%	0.012%

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Integrator: RTE

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Stop Thrs : 0

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

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Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM051622WMA.M
 Title : VOC Analysis

37	8.925	2352	2359	2366	rBV5	1470	1582	0.03%	0.006%
38	9.124	2415	2421	2425	rVV4	837	896	0.02%	0.003%
39	9.269	2460	2466	2467	rBV2	641	508	0.01%	0.002%
40	9.356	2487	2493	2496	rVB2	592	561	0.01%	0.002%
41	9.417	2499	2512	2535	rBV	540643	881523	16.76%	3.198%
42	9.571	2549	2560	2574	rVB	100386	157811	3.00%	0.573%
43	9.690	2584	2597	2614	rBV2	184114	314549	5.98%	1.141%
44	9.883	2652	2657	2660	rBV3	1229	940	0.02%	0.003%
45	10.034	2699	2704	2710	rVV5	922	1341	0.03%	0.005%
46	10.102	2712	2725	2743	rVB	122681	196715	3.74%	0.714%
47	10.275	2770	2779	2785	rVB6	4607	6266	0.12%	0.023%
48	10.362	2794	2806	2809	rBV5	1401	1977	0.04%	0.007%
49	10.378	2809	2811	2814	rVB2	989	521	0.01%	0.002%
50	10.484	2834	2844	2856	rVB2	17662	25825	0.49%	0.094%
51	10.571	2864	2871	2878	rBV2	2589	3413	0.06%	0.012%
52	10.645	2886	2894	2902	rBV6	3588	5981	0.11%	0.022%
53	10.697	2904	2910	2916	rVB6	1900	2678	0.05%	0.010%
54	10.758	2916	2929	2944	rBV	412359	676890	12.87%	2.456%
55	10.906	2964	2975	2991	rVB	35990	58145	1.11%	0.211%
56	10.999	2993	3004	3022	rVV	90042	199691	3.80%	0.724%
57	11.089	3022	3032	3044	rVB	53605	83335	1.58%	0.302%
58	11.234	3066	3077	3084	rBV4	6358	9314	0.18%	0.034%
59	11.291	3085	3095	3113	rVB	96384	153606	2.92%	0.557%
60	11.420	3129	3135	3138	rBV	2165	2477	0.05%	0.009%
61	11.468	3139	3150	3167	rVB	292771	444136	8.45%	1.611%
62	11.607	3186	3193	3197	rVV2	5133	6808	0.13%	0.025%
63	11.645	3199	3205	3215	rVB2	13624	20170	0.38%	0.073%
64	11.745	3226	3236	3244	rBV2	19550	28800	0.55%	0.104%
65	11.812	3244	3257	3272	rVV	546799	869431	16.53%	3.154%
66	11.896	3272	3283	3297	rVB	141692	217547	4.14%	0.789%
67	12.008	3310	3318	3327	rVB2	3664	5092	0.10%	0.018%
68	12.095	3334	3345	3352	rBV2	66334	115244	2.19%	0.418%
69	12.195	3363	3376	3394	rVB	555808	945860	17.99%	3.431%
70	12.304	3399	3410	3421	rVB9	4696	7783	0.15%	0.028%
71	12.378	3425	3433	3443	rVB	16192	24404	0.46%	0.089%
72	12.491	3452	3468	3481	rBV4	49187	117480	2.23%	0.426%
73	12.571	3485	3493	3502	rVV2	26752	42615	0.81%	0.155%
74	12.610	3502	3505	3518	rVB8	5751	7806	0.15%	0.028%
75	12.687	3518	3529	3547	rBV5	16489	44293	0.84%	0.161%
76	12.774	3547	3556	3561	rVV5	1443	2571	0.05%	0.009%

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 Title : VOC Analysis

77	12.812	3564	3568	3574	rVB5	1707	2022	0.04%	0.007%
78	12.880	3579	3589	3597	rVB3	9413	15472	0.29%	0.056%
79	12.976	3603	3619	3627	rBV3	11954	21638	0.41%	0.078%
80	13.028	3627	3635	3646	rVB2	22878	34413	0.65%	0.125%
81	13.111	3651	3661	3667	rBV6	1891	2998	0.06%	0.011%
82	13.256	3699	3706	3710	rBV5	1564	2221	0.04%	0.008%
83	13.295	3713	3718	3724	rVB7	3892	5078	0.10%	0.018%
84	13.410	3747	3754	3757	rBV5	1667	2220	0.04%	0.008%
85	13.449	3758	3766	3781	rVB3	28481	47335	0.90%	0.172%
86	13.523	3783	3789	3797	rBV7	2114	3022	0.06%	0.011%
87	13.626	3814	3821	3829	rVB5	8054	10942	0.21%	0.040%
88	13.738	3852	3856	3859	rVB3	893	677	0.01%	0.002%
89	13.790	3865	3872	3879	rVB3	1141	1961	0.04%	0.007%
90	13.838	3879	3887	3894	rBV8	2159	3568	0.07%	0.013%
91	13.909	3905	3909	3912	rBV4	585	669	0.01%	0.002%
92	13.944	3916	3920	3923	rVV3	972	908	0.02%	0.003%
93	14.092	3952	3966	3979	rBV	26975	48778	0.93%	0.177%
94	14.295	4023	4029	4030	rBV2	692	581	0.01%	0.002%
95	14.336	4038	4042	4044	rBV4	920	550	0.01%	0.002%
96	14.687	4148	4151	4156	rVB4	852	764	0.01%	0.003%
97	14.957	4230	4235	4239	rBV3	469	569	0.01%	0.002%
98	15.423	4371	4380	4383	rBV5	2513	3863	0.07%	0.014%
99	16.278	4643	4646	4649	rBV3	1291	956	0.02%	0.003%
100	16.593	4740	4744	4746	rBV4	1473	1365	0.03%	0.005%

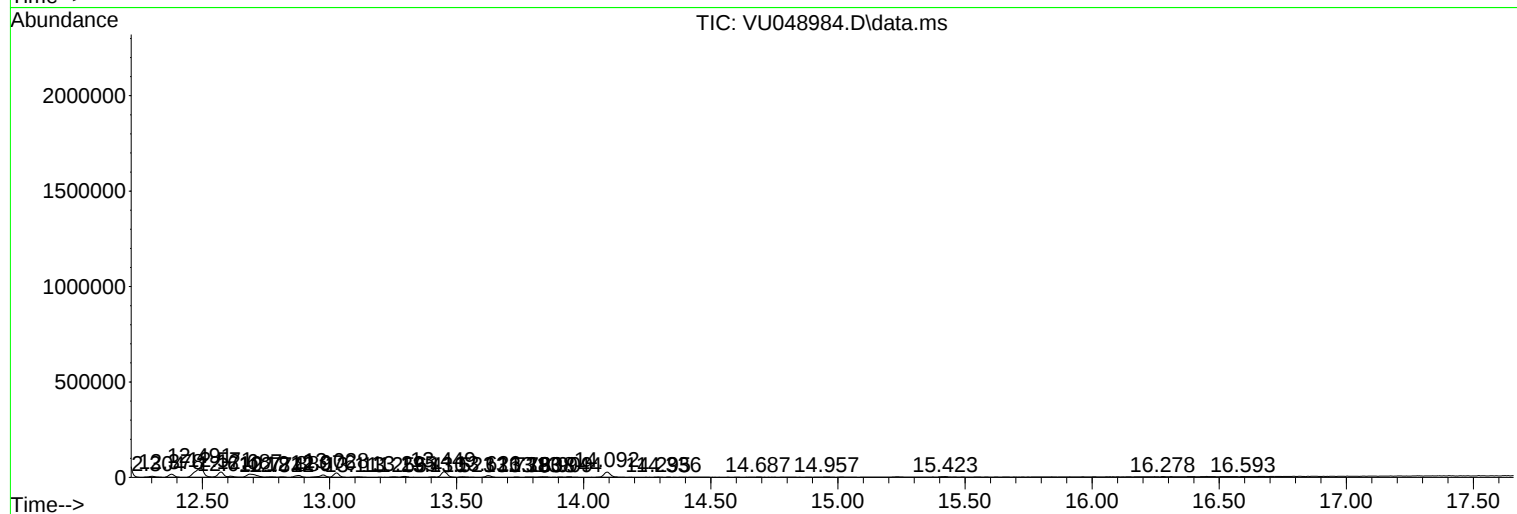
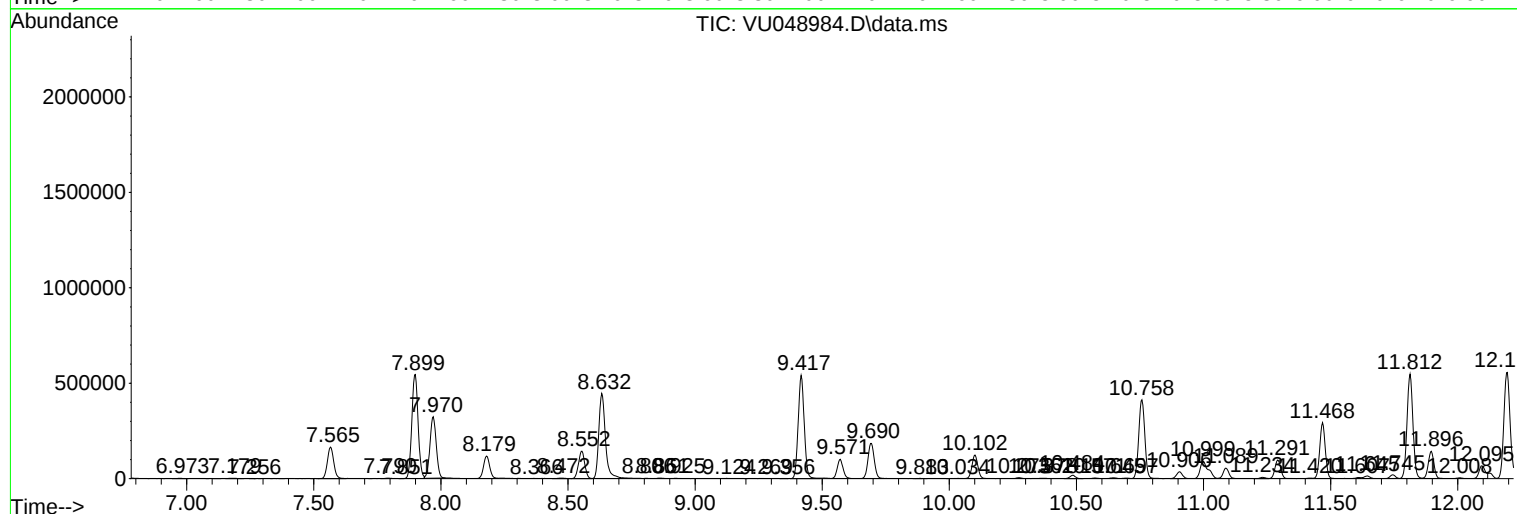
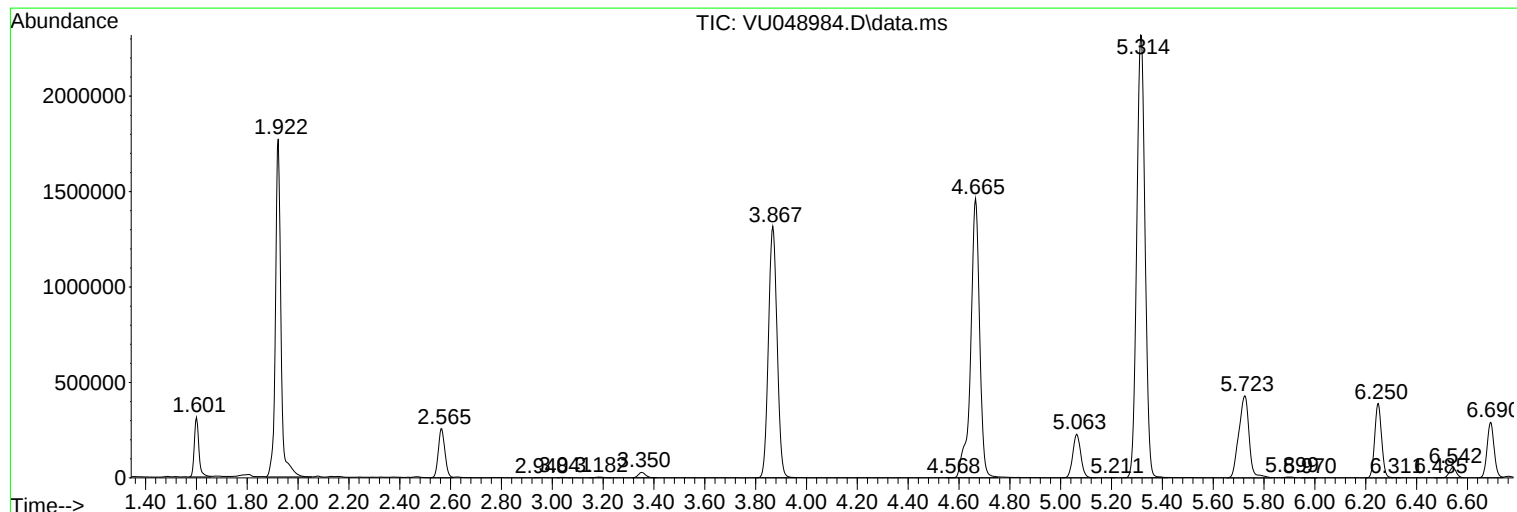
Sum of corrected areas: 27564984

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Instrument :
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ClientSampleId :
EXYY0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM051622WMA.M
Quant Title : VOC Analysis

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



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Instrument :
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ClientSampleId :
EXYY0

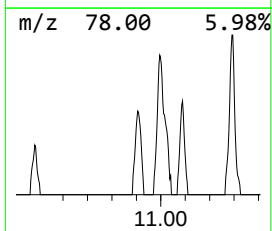
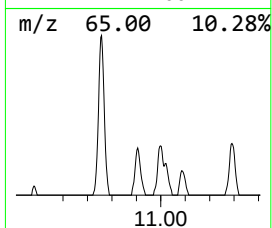
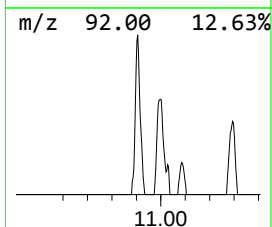
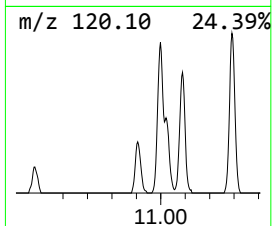
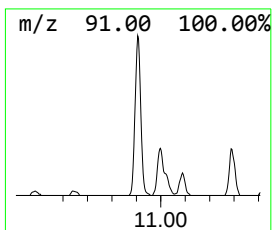
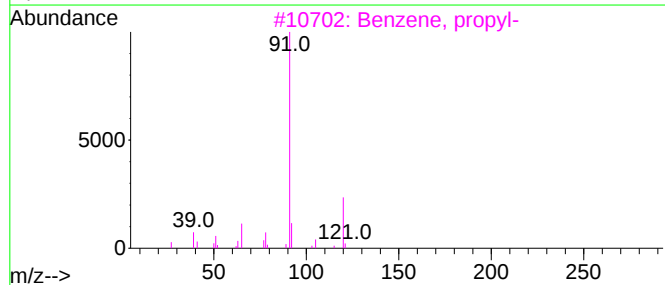
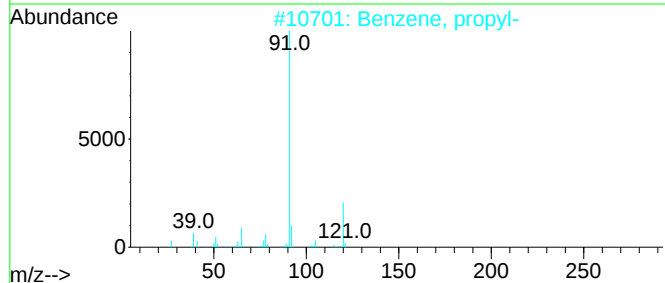
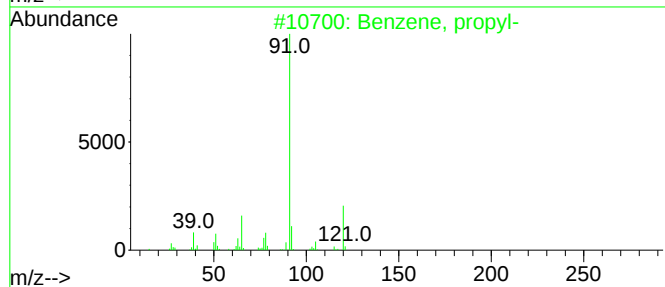
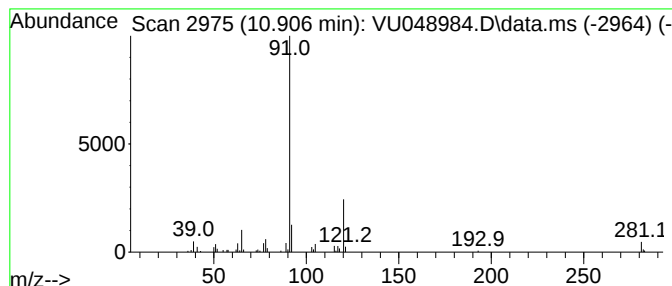
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM051622WMA.M
Quant Title : VOC Analysis

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

Peak Number 2 Benzene, propyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.906	3.34 ug/L	58145	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	87
2			Benzene, propyl-	120	C9H12	000103-65-1	81
3			Benzene, propyl-	120	C9H12	000103-65-1	74
4			Benzene, propyl-	120	C9H12	000103-65-1	74
5			N-Isobutyl(phenyl)methanesulfona...	227	C11H17NO2S	144615-94-1	72



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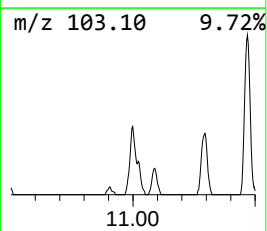
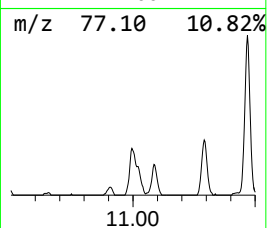
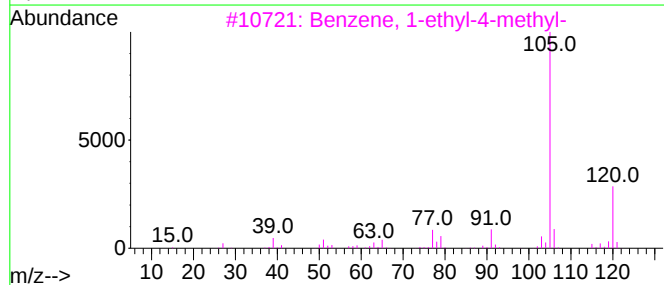
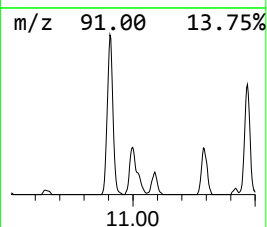
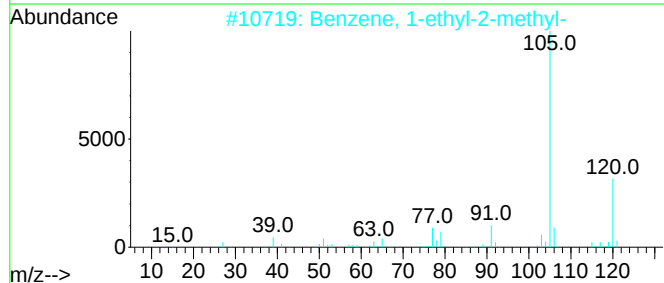
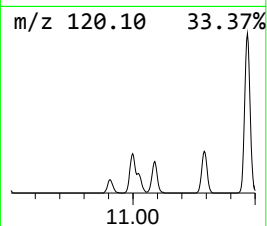
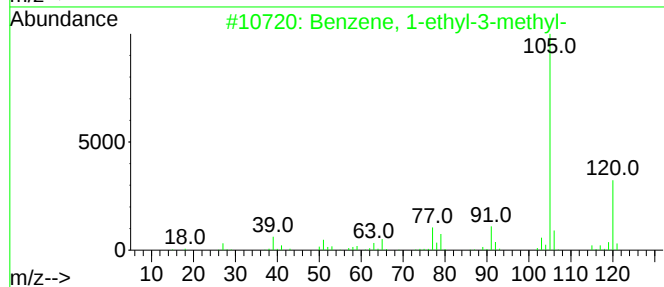
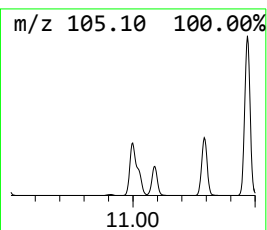
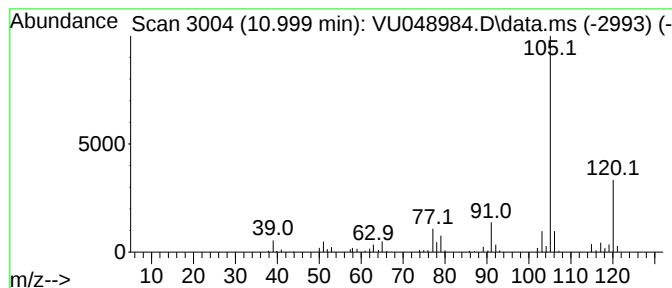
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM051622WMA.M
Quant Title : VOC Analysis

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.999	11.48 ug/L	199691	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	97
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94



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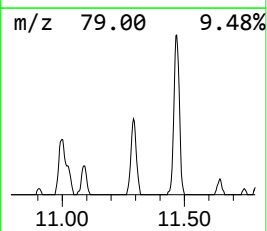
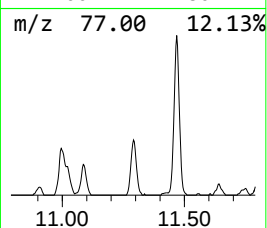
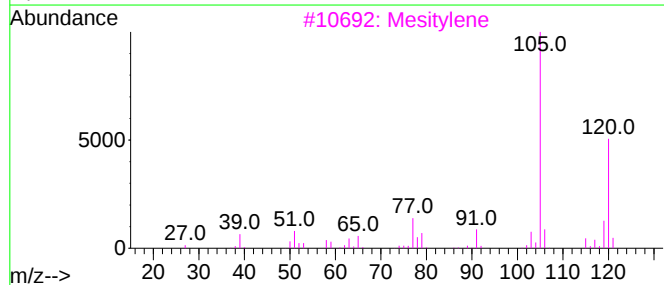
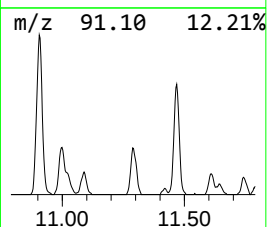
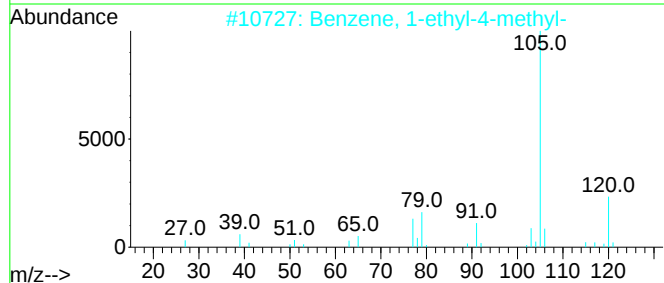
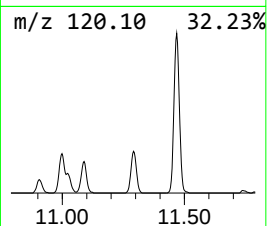
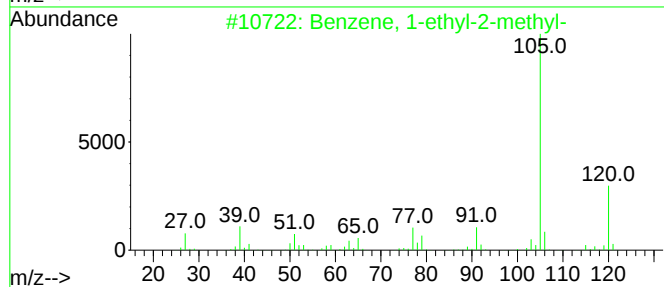
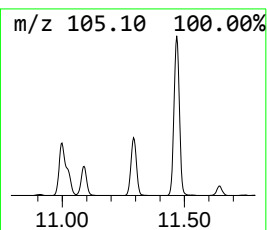
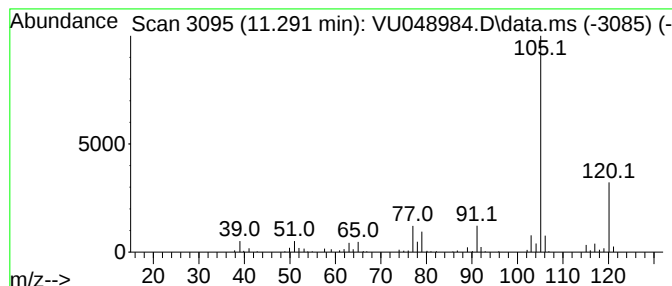
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Quant Title : VOC Analysis

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

Peak Number 4 Benzene, 1-ethyl-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.291	8.83 ug/L	153606	1,4-Dichlorobenzene-d4	11.812

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU060922\
Data File : VU048984.D
Acq On : 09 Jun 2022 18:29
Operator : SY/MD
Sample : N3248-09
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXYY0

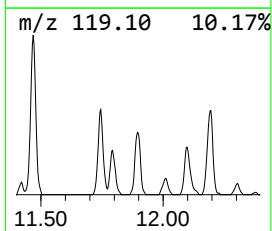
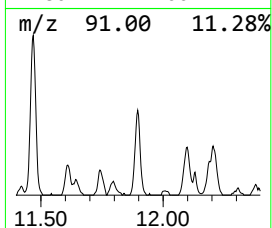
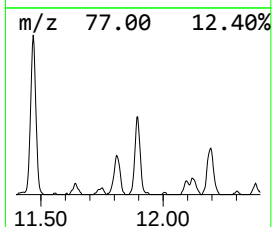
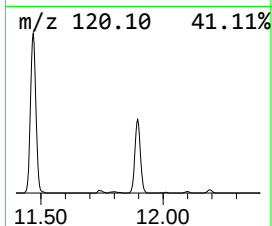
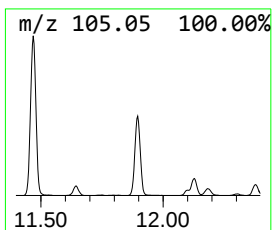
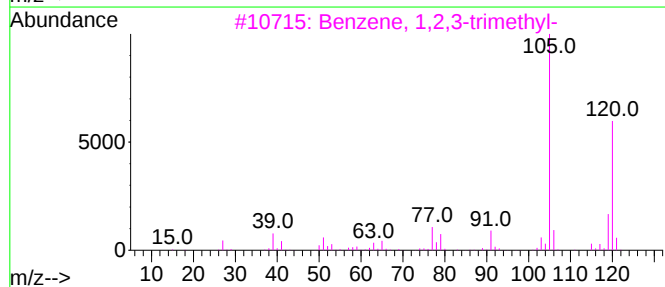
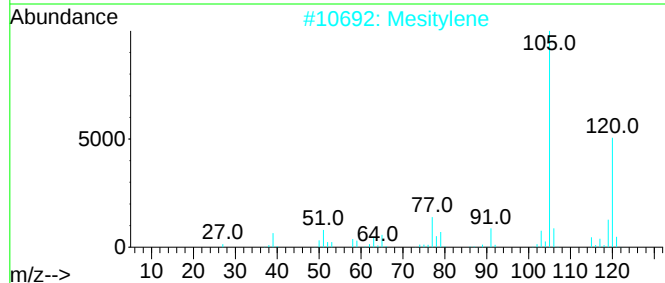
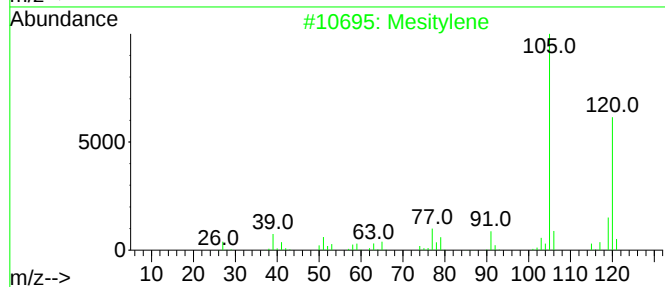
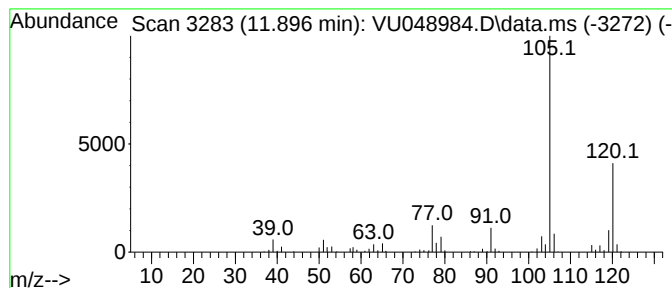
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Quant Title : VOC Analysis

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

Peak Number 5 Benzene, 1,2,3-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.896	12.51 ug/L	217547	1,4-Dichlorobenzene-d4	11.812

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			Mesitylene	120	C9H12	000108-67-8	94
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU060922\
Data File : VU048984.D
Acq On : 09 Jun 2022 18:29
Operator : SY/MD
Sample : N3248-09
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXYY0

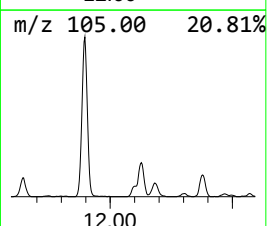
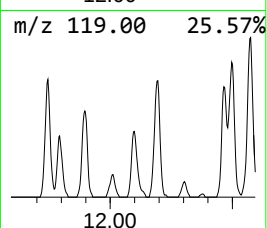
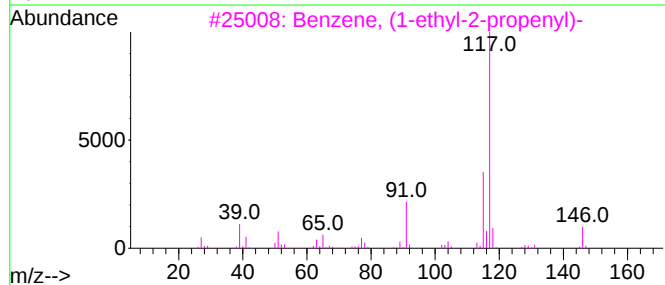
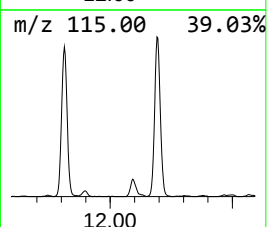
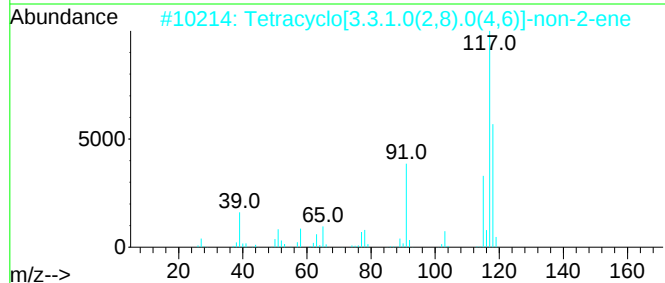
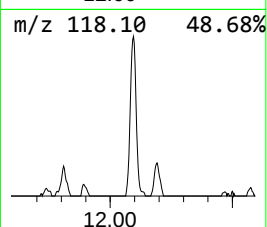
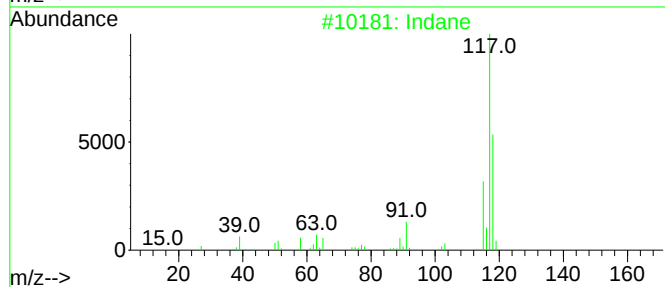
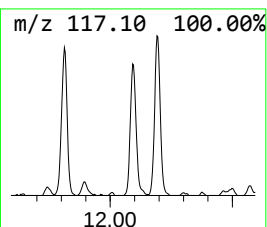
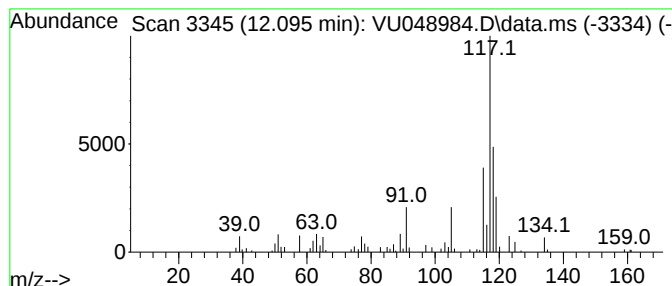
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Quant Title : VOC Analysis

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

Peak Number 6 Indane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.095	6.63 ug/L	115244	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	60
2			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	49
3			Benzene, (1-ethyl-2-propenyl)-	146	C11H14	019947-22-9	47
4			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	47
5			trans-Cinnamyl bromide	196	C9H9Br	026146-77-0	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU060922\
Data File : VU048984.D
Acq On : 09 Jun 2022 18:29
Operator : SY/MD
Sample : N3248-09
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXYY0

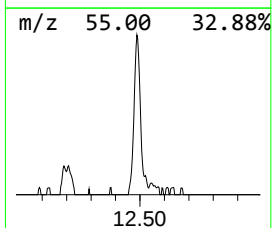
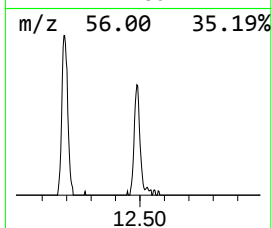
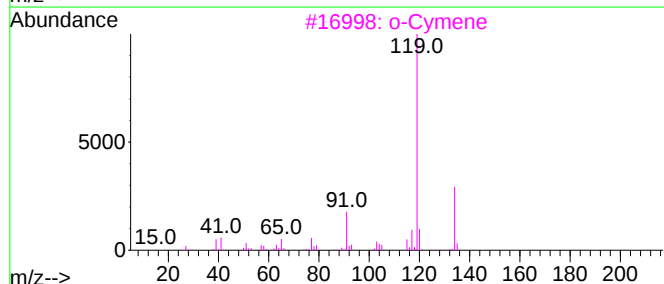
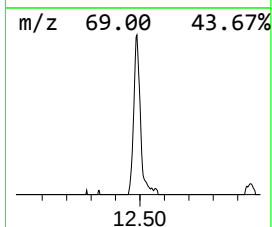
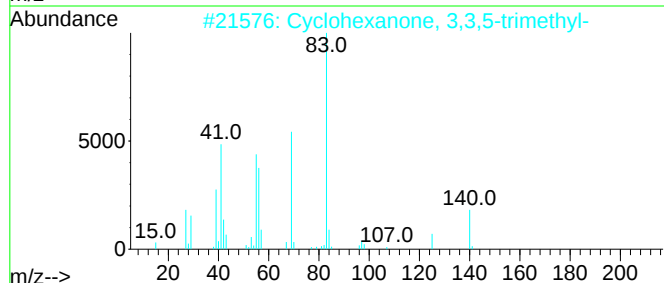
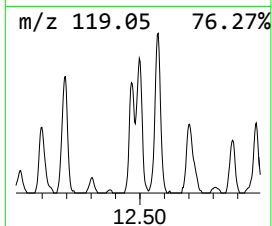
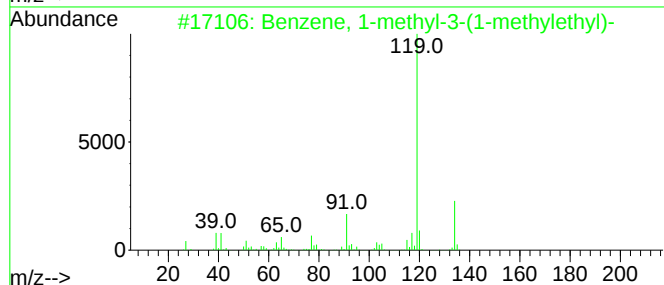
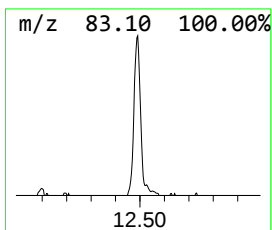
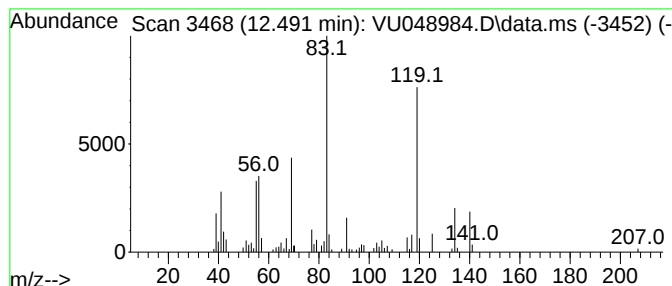
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Quant Title : VOC Analysis

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

Peak Number 7 Benzene, 1-methyl-3-(1-meth... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.491	6.76 ug/L	117480	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	86
2			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	83
3			o-Cymene	134	C10H14	000527-84-4	83
4			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	50
5			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU060922\
Data File : VU048984.D
Acq On : 09 Jun 2022 18:29
Operator : SY/MD
Sample : N3248-09
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXYY0

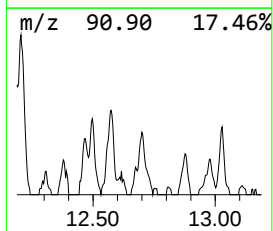
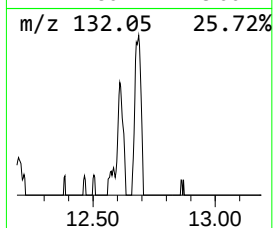
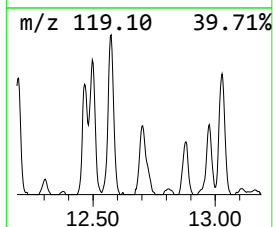
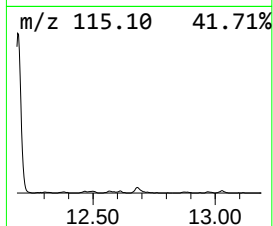
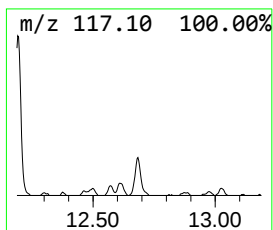
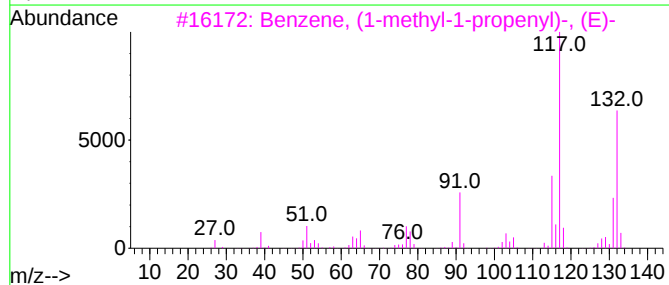
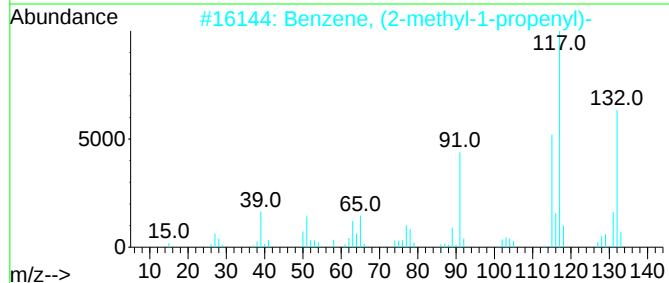
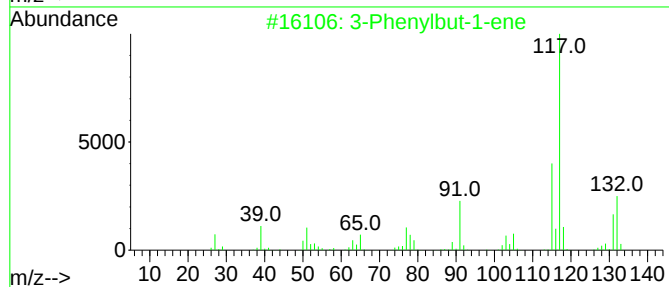
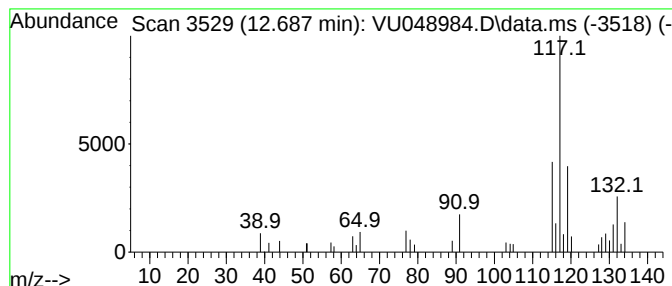
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Quant Title : VOC Analysis

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

Peak Number 8 3-Phenylbut-1-ene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.687	2.55 ug/L	44293	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Phenylbut-1-ene	132	C10H12	000934-10-1	76
2			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	76
3			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	64
4			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	64
5			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	62



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU060922\
Data File : VU048984.D
Acq On : 09 Jun 2022 18:29
Operator : SY/MD
Sample : N3248-09
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXYY0

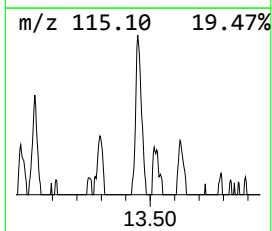
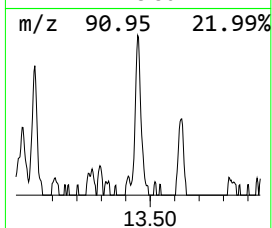
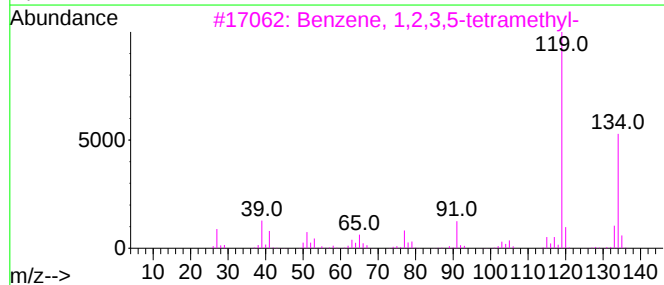
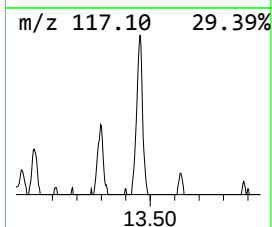
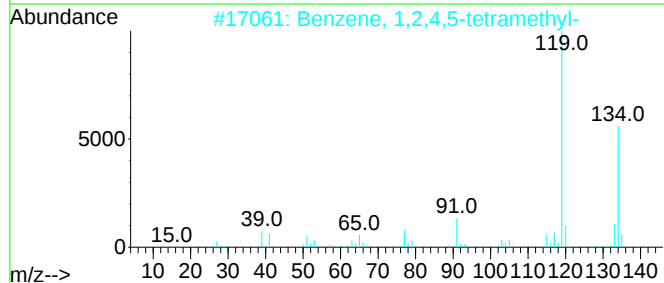
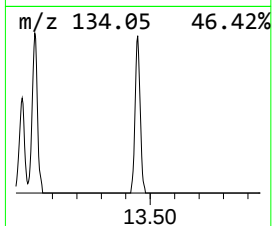
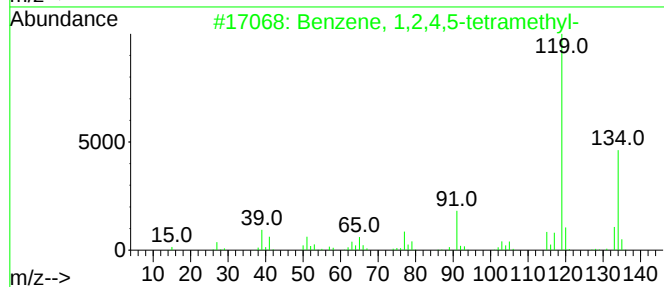
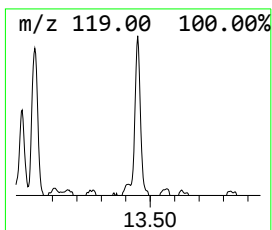
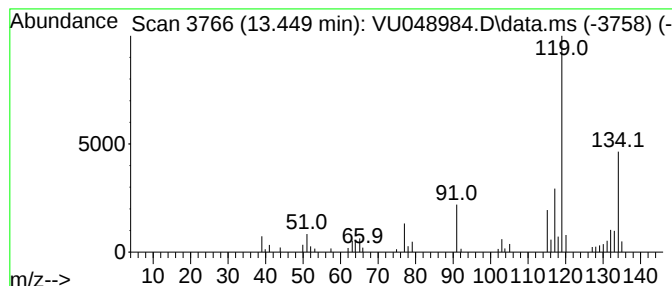
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Quant Title : VOC Analysis

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

Peak Number 9 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.449	2.72 ug/L	47335	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	76
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	68
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	68
4			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	68
5			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU060922\
Data File : VU048984.D
Acq On : 09 Jun 2022 18:29
Operator : SY/MD
Sample : N3248-09
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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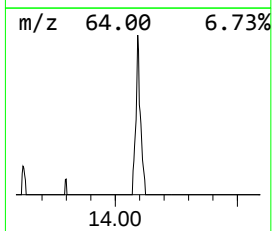
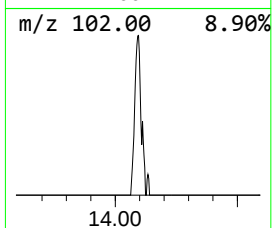
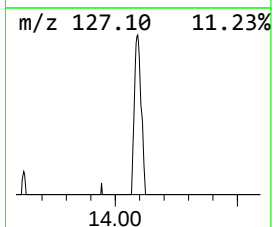
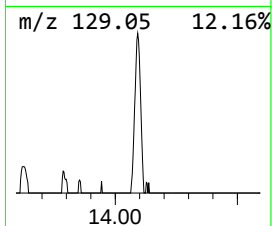
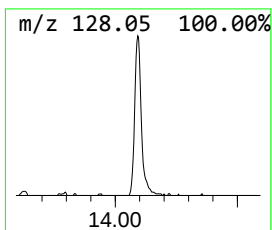
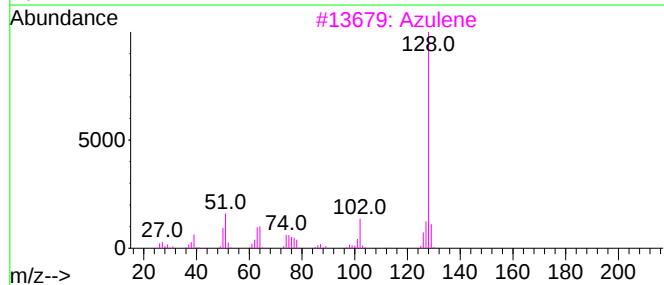
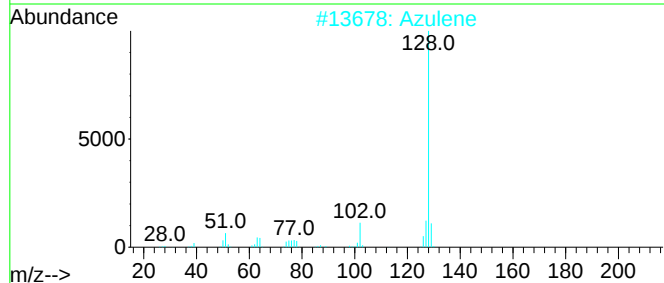
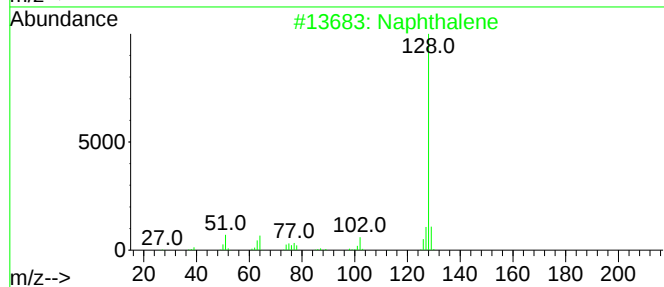
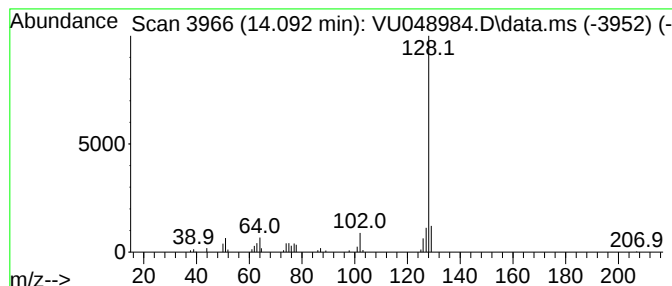
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Quant Title : VOC Analysis

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

Peak Number 10 Azulene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.092	2.81 ug/L	48778	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene	128	C10H8	000091-20-3	94
2			Azulene	128	C10H8	000275-51-4	91
3			Azulene	128	C10H8	000275-51-4	91
4			Naphthalene	128	C10H8	000091-20-3	91
5			Naphthalene	128	C10H8	000091-20-3	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU060922\
Data File : VU048984.D
Acq On : 09 Jun 2022 18:29
Operator : SY/MD
Sample : N3248-09
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXYY0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM051622WMA.M
Quant Title : VOC Analysis

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, propyl-	10.906	3.3	ug/L	58145	3	11.812	869431	50.0
Benzene, 1-ethy...	10.999	11.5	ug/L	199691	3	11.812	869431	50.0
Benzene, 1-ethy...	11.291	8.8	ug/L	153606	3	11.812	869431	50.0
Benzene, 1,2,3-...	11.896	12.5	ug/L	217547	3	11.812	869431	50.0
Indane	12.095	6.6	ug/L	115244	3	11.812	869431	50.0
Benzene, 1-meth...	12.491	6.8	ug/L	117480	3	11.812	869431	50.0
3-Phenylbut-1-ene	12.687	2.5	ug/L	44293	3	11.812	869431	50.0
Benzene, 1,2,4,...	13.449	2.7	ug/L	47335	3	11.812	869431	50.0
Azulene	14.092	2.8	ug/L	48778	3	11.812	869431	50.0