

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081821\
 Data File : VV021866.D
 Acq On : 18 Aug 2021 15:28
 Operator : SY/MD
 Sample : M3448-03
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 DBKD5

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\SFAMVLM081621WMA.M

Title : VOC Analysis

Signal : TIC: VV021866.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.304	76	82	95	rVB	180601	179913	12.67%	1.107%
2	1.561	155	162	178	rVB	148726	176020	12.40%	1.083%
3	2.105	322	331	347	rVB	294682	431782	30.42%	2.657%
4	2.613	486	489	499	rVV6	848	1123	0.08%	0.007%
5	2.822	552	554	558	rVB2	567	313	0.02%	0.002%
6	2.883	570	573	577	rVB2	440	248	0.02%	0.002%
7	3.028	614	618	621	rBV3	431	419	0.03%	0.003%
8	3.079	630	634	637	rBV3	255	217	0.02%	0.001%
9	3.343	713	716	719	rBV	330	249	0.02%	0.002%
10	3.539	770	777	778	rBV	232	203	0.01%	0.001%
11	3.703	825	828	832	rBV2	249	206	0.01%	0.001%
12	3.876	869	882	926	rBV2	124744	350772	24.71%	2.158%
13	4.352	1013	1030	1055	rBV	217386	543146	38.26%	3.342%
14	4.664	1124	1127	1128	rBV2	571	236	0.02%	0.001%
15	4.732	1146	1148	1151	rVB2	364	197	0.01%	0.001%
16	4.786	1162	1165	1169	rVB3	427	257	0.02%	0.002%
17	4.873	1188	1192	1194	rBV	354	257	0.02%	0.002%
18	4.970	1220	1222	1226	rBV	252	198	0.01%	0.001%
19	5.047	1229	1246	1278	rBV2	523319	1346826	94.88%	8.287%
20	5.494	1383	1385	1389	rVB2	543	295	0.02%	0.002%
21	5.516	1389	1392	1395	rBV	320	273	0.02%	0.002%
22	5.564	1405	1407	1411	rBV2	418	239	0.02%	0.001%
23	5.619	1411	1424	1452	rBV	456820	921185	64.90%	5.668%
24	5.912	1504	1515	1530	rBV4	10107	21575	1.52%	0.133%
25	6.072	1549	1565	1592	rBV	303349	621214	43.76%	3.822%
26	6.413	1668	1671	1672	rBV3	496	287	0.02%	0.002%
27	6.458	1681	1685	1688	rBV2	477	394	0.03%	0.002%
28	6.625	1733	1737	1739	rBV3	672	435	0.03%	0.003%
29	6.642	1739	1742	1744	rBV2	719	427	0.03%	0.003%
30	6.728	1766	1769	1772	rBV3	524	366	0.03%	0.002%
31	6.918	1824	1828	1833	rBV2	330	312	0.02%	0.002%
32	6.944	1833	1836	1838	rBV	335	208	0.01%	0.001%
33	6.989	1838	1850	1870	rBV	161397	295571	20.82%	1.819%
34	7.195	1910	1914	1918	rBV4	761	664	0.05%	0.004%
35	7.320	1941	1953	1970	rBV	585417	1017745	71.70%	6.262%
36	7.625	2038	2048	2071	rBV	105780	189666	13.36%	1.167%

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Integrator: RTE

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Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
 Title : VOC Analysis

37	7.867	2122	2123	2126	rBV3	383	252	0.02%	0.002%
38	7.950	2138	2149	2155	rVB5	2104	3604	0.25%	0.022%
39	8.092	2181	2193	2220	rBV	344502	630690	44.43%	3.881%
40	8.667	2366	2372	2376	rBV3	331	354	0.02%	0.002%
41	8.741	2392	2395	2397	rBV2	395	220	0.02%	0.001%
42	8.854	2418	2430	2457	rBV	680463	1103526	77.74%	6.790%
43	9.018	2475	2481	2496	rVB	13003	22241	1.57%	0.137%
44	9.153	2516	2523	2532	rVB2	6857	10273	0.72%	0.063%
45	9.236	2548	2549	2553	rVB	592	271	0.02%	0.002%
46	9.278	2557	2562	2564	rBV2	393	324	0.02%	0.002%
47	9.294	2564	2567	2571	rBV2	246	194	0.01%	0.001%
48	9.313	2571	2573	2576	rBV2	352	186	0.01%	0.001%
49	9.413	2601	2604	2610	rVB3	447	374	0.03%	0.002%
50	9.445	2610	2614	2619	rBV2	476	414	0.03%	0.003%
51	9.551	2640	2647	2659	rVB2	5447	8483	0.60%	0.052%
52	9.709	2694	2696	2702	rBB3	277	264	0.02%	0.002%
53	9.899	2750	2755	2758	rVB2	307	245	0.02%	0.002%
54	9.940	2758	2768	2779	rBV4	3849	5856	0.41%	0.036%
55	10.098	2814	2817	2821	rVB3	489	343	0.02%	0.002%
56	10.130	2821	2827	2828	rBV2	585	585	0.04%	0.004%
57	10.217	2841	2854	2881	rBV	467811	741554	52.24%	4.563%
58	10.477	2929	2935	2947	rVB2	5003	8059	0.57%	0.050%
59	10.548	2951	2957	2964	rVB6	3081	4040	0.28%	0.025%
60	10.661	2990	2992	2995	rVB	528	247	0.02%	0.002%
61	10.683	2995	2999	3000	rBV	299	229	0.02%	0.001%
62	10.744	3007	3018	3033	rVV2	8413	13491	0.95%	0.083%
63	10.921	3063	3073	3087	rVB	14124	22191	1.56%	0.137%
64	10.992	3092	3095	3099	rBV3	368	245	0.02%	0.002%
65	11.040	3105	3110	3112	rBV3	408	299	0.02%	0.002%
66	11.085	3122	3124	3130	rVB4	1038	673	0.05%	0.004%
67	11.137	3138	3140	3142	rBV	397	192	0.01%	0.001%
68	11.149	3142	3144	3146	rBV2	582	320	0.02%	0.002%
69	11.191	3156	3157	3159	rBV2	564	271	0.02%	0.002%
70	11.249	3164	3175	3195	rBV	711742	1112990	78.41%	6.848%
71	11.529	3249	3262	3280	rBV	681280	1049086	73.91%	6.455%
72	11.625	3281	3292	3314	rVB	688312	1085984	76.51%	6.682%
73	11.918	3373	3383	3391	rBV4	3282	5243	0.37%	0.032%
74	12.021	3404	3415	3419	rBV2	11196	16985	1.20%	0.105%
75	12.117	3436	3445	3457	rBV	41889	65319	4.60%	0.402%
76	12.246	3479	3485	3487	rBV4	1039	935	0.07%	0.006%

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 Title : VOC Analysis

77	12.365	3514	3522	3531	rBV	19182	28760	2.03%	0.177%
78	12.416	3531	3538	3545	rBV	8720	11745	0.83%	0.072%
79	12.468	3546	3554	3567	rBV3	31095	59197	4.17%	0.364%
80	12.641	3602	3608	3610	rBV4	2356	2272	0.16%	0.014%
81	12.728	3625	3635	3653	rVB	64682	102882	7.25%	0.633%
82	12.886	3670	3684	3696	rBV	165713	259298	18.27%	1.595%
83	13.005	3717	3721	3726	rBV4	1742	1851	0.13%	0.011%
84	13.053	3726	3736	3747	rVV4	23319	39284	2.77%	0.242%
85	13.117	3754	3756	3757	rBV	454	184	0.01%	0.001%
86	13.275	3797	3805	3811	rBV5	10068	14476	1.02%	0.089%
87	13.432	3852	3854	3857	rBV3	781	515	0.04%	0.003%
88	13.468	3861	3865	3867	rBV2	882	615	0.04%	0.004%
89	13.506	3867	3877	3885	rBV6	11588	23372	1.65%	0.144%
90	13.677	3923	3930	3935	rBV9	3629	5926	0.42%	0.036%
91	13.911	3993	4003	4016	rBV2	14574	23389	1.65%	0.144%
92	14.095	4051	4060	4073	rBV4	16034	35446	2.50%	0.218%
93	14.580	4201	4211	4219	rBV	62975	101178	7.13%	0.623%
94	14.635	4220	4228	4240	rVB	139083	214705	15.13%	1.321%
95	14.754	4258	4265	4273	rBV3	9278	16404	1.16%	0.101%
96	14.818	4274	4285	4298	rVV	899216	1419491	100.00%	8.734%
97	15.667	4540	4549	4571	rVB2	82624	164658	11.60%	1.013%
98	15.812	4585	4594	4601	rBV	166733	252989	17.82%	1.557%
99	16.181	4702	4709	4723	rVB	36434	55328	3.90%	0.340%
100	16.487	4793	4804	4825	rVB	906098	1398032	98.49%	8.602%

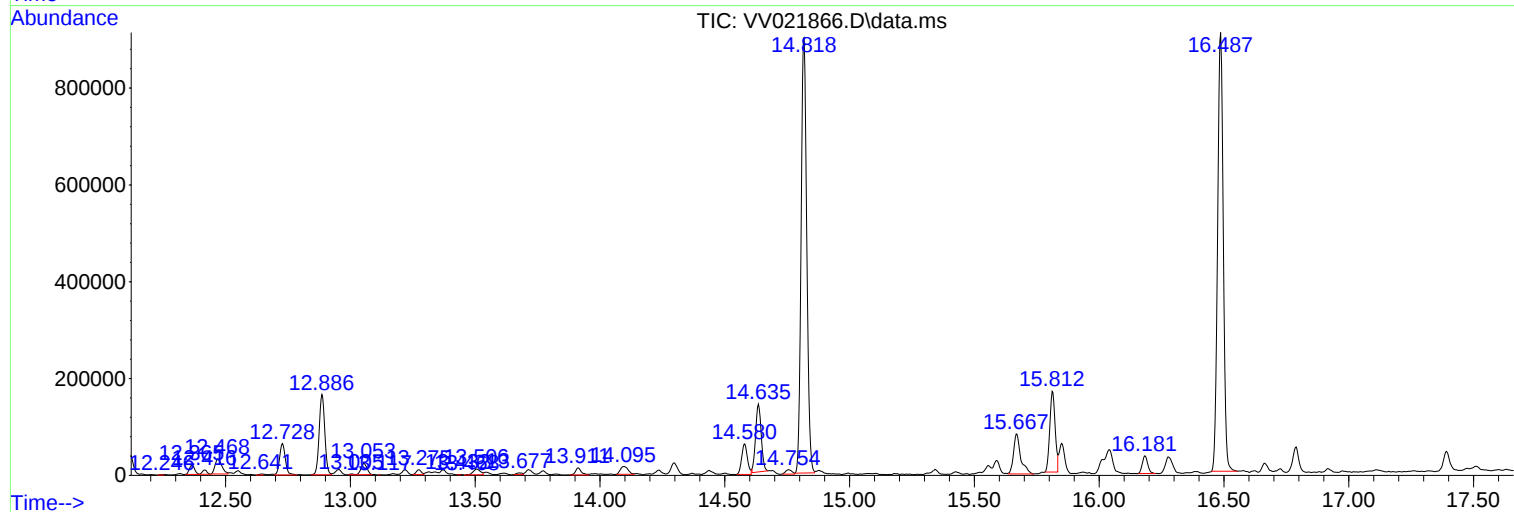
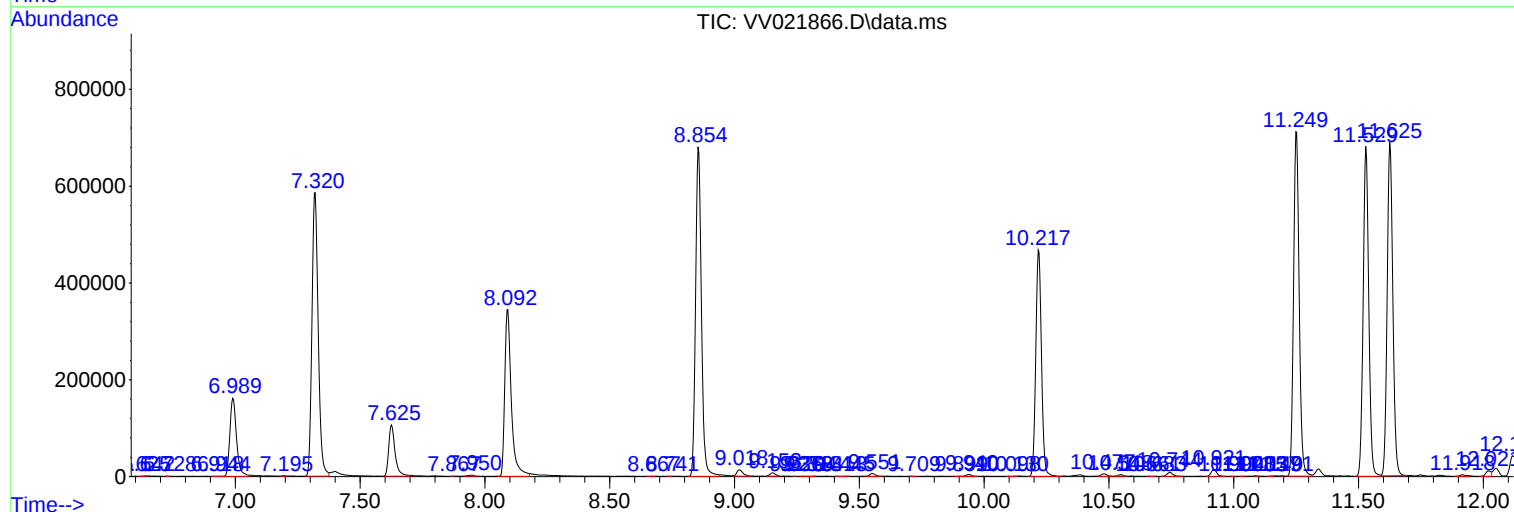
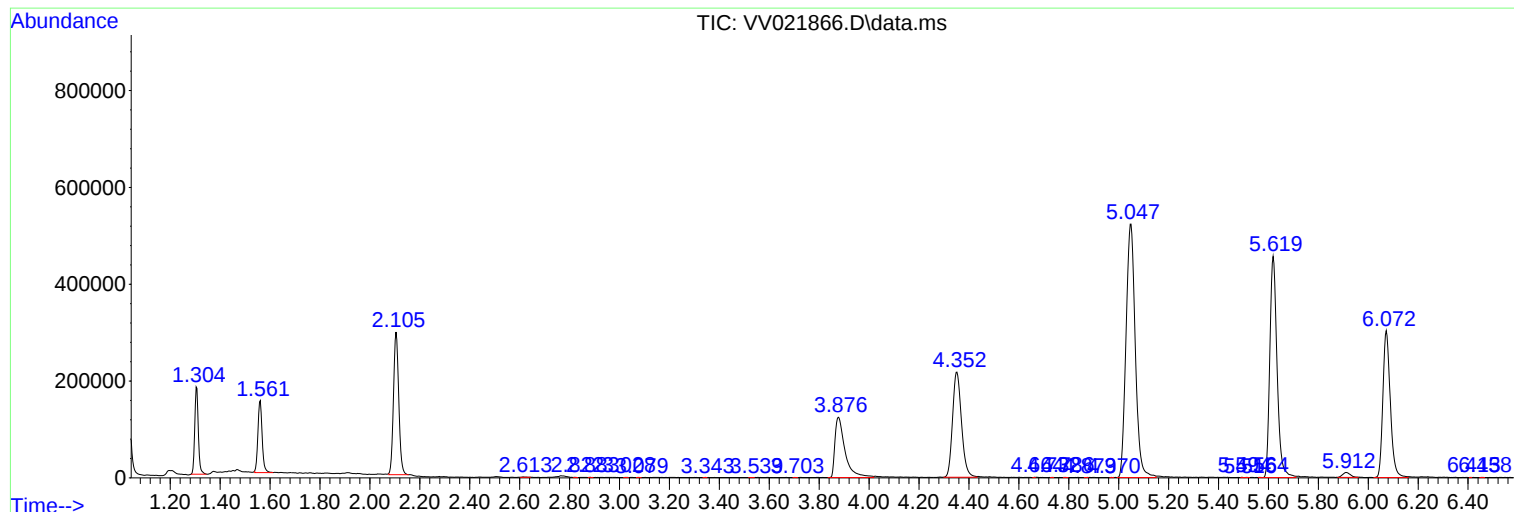
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ALS Vial : 15 Sample Multiplier: 1

Instrument :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

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



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DBKD5

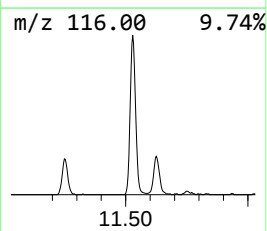
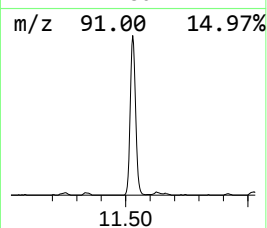
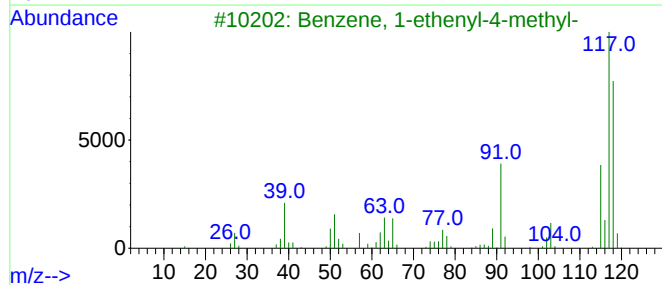
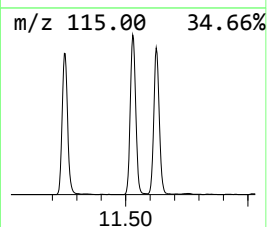
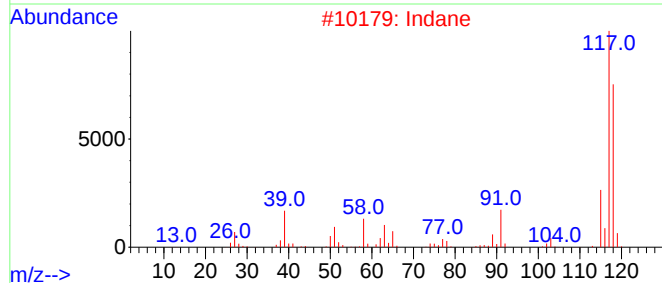
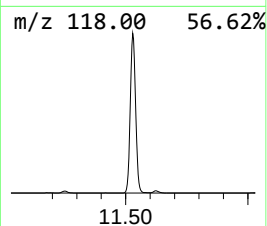
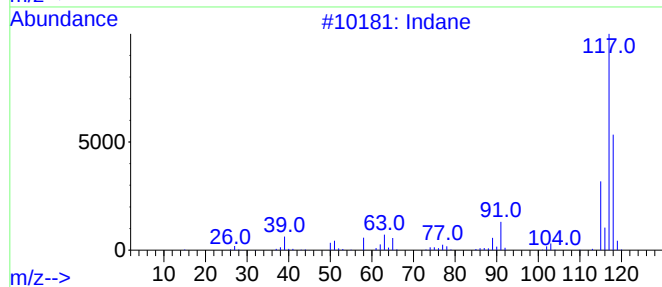
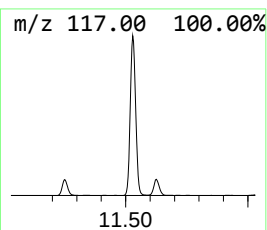
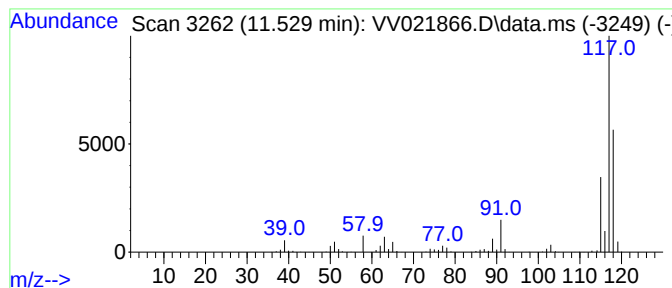
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

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

Peak Number 2 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.529	47.13 ug/L	1049090	1,4-Dichlorobenzene-d4	11.252

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	93
2			Indane	118	C9H10	000496-11-7	91
3			Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	83
4			Delta-cyclene	118	C9H10	007785-10-6	80
5			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	80



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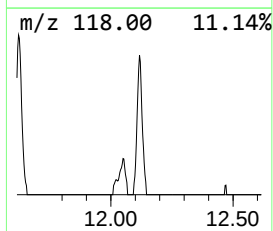
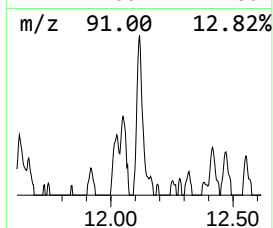
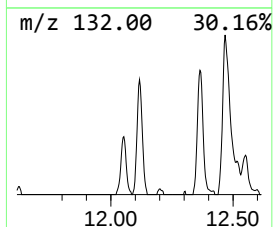
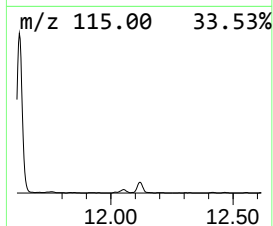
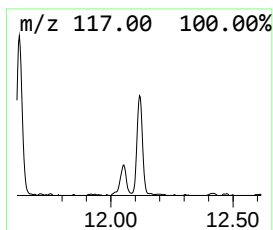
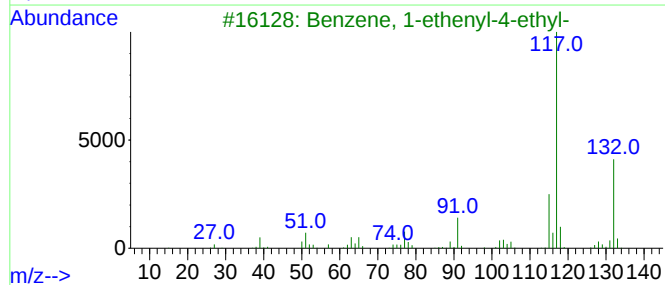
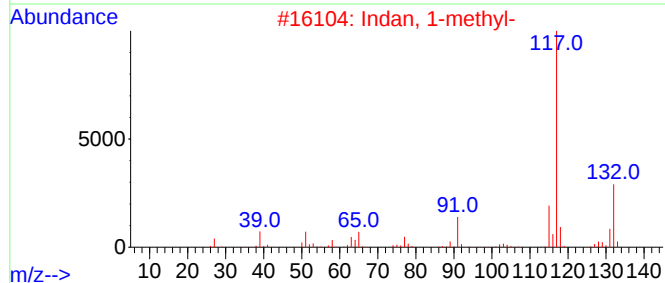
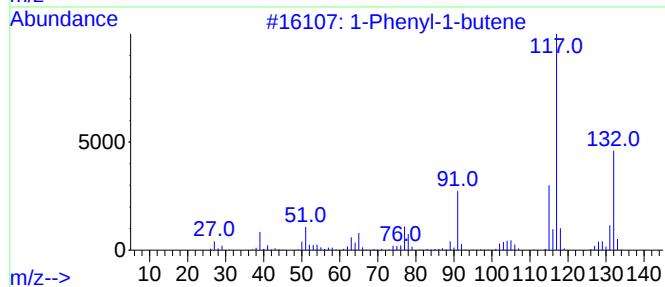
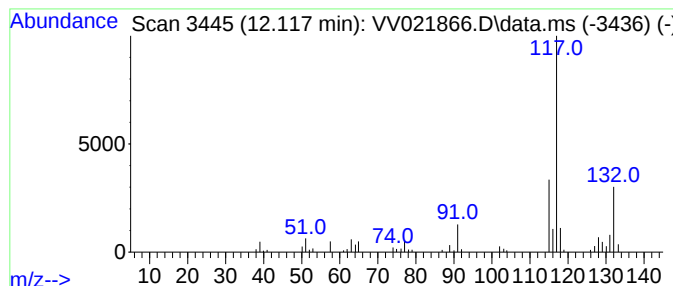
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

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

Peak Number 3 1-Phenyl-1-butene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.117	2.93 ug/L	65319	1,4-Dichlorobenzene-d4	11.252

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenyl-1-butene	132	C10H12	000824-90-8	90
2			Indan, 1-methyl-	132	C10H12	000767-58-8	90
3			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
4			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	86
5			Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	86



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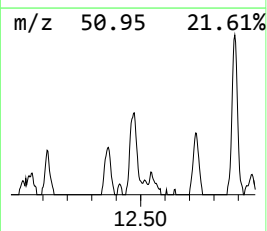
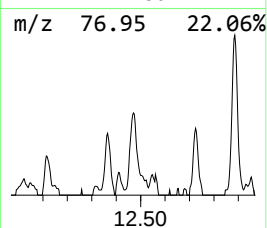
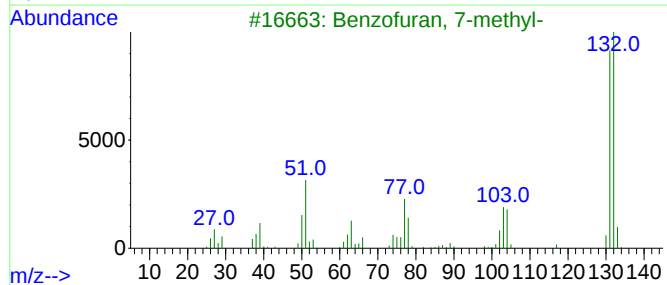
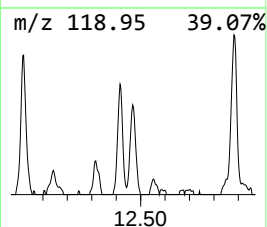
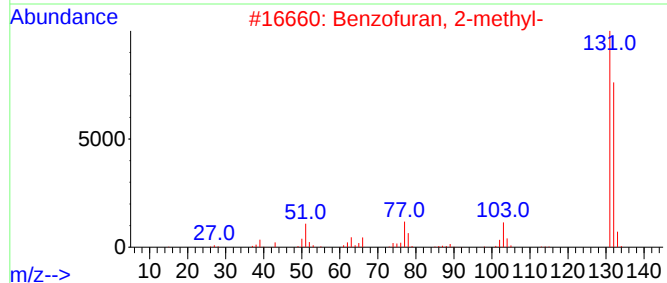
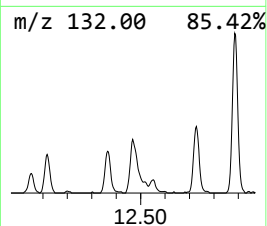
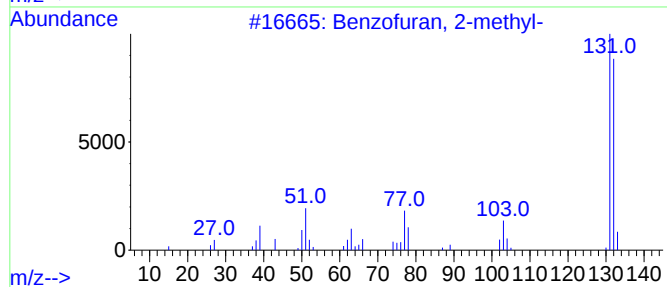
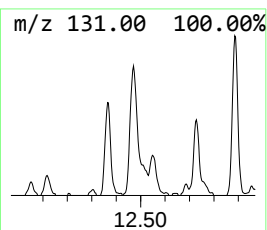
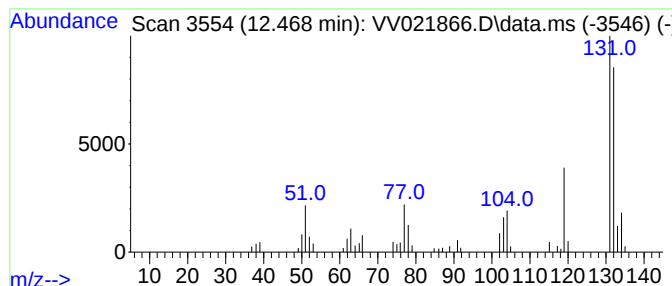
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.M
Quant Title : VOC Analysis

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

Peak Number 4 Benzofuran, 2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.468	2.66 ug/L	59197	1,4-Dichlorobenzene-d4	11.252

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	95
2			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	95
3			Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	94
4			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
5			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	89



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081821\
Data File : VV021866.D
Acq On : 18 Aug 2021 15:28
Operator : SY/MD
Sample : M3448-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKD5

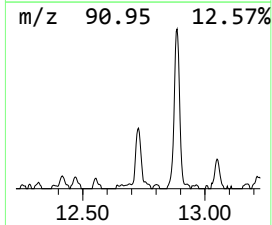
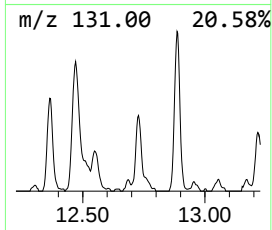
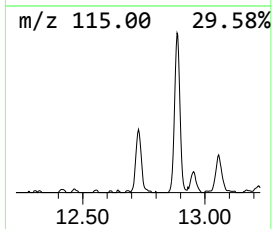
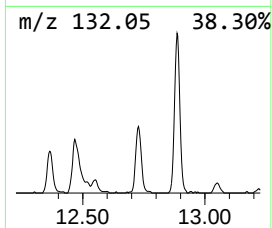
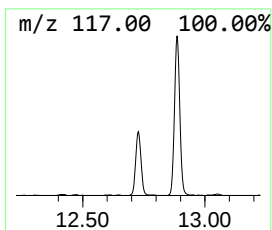
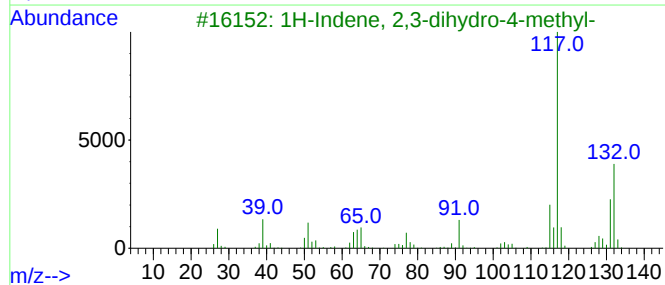
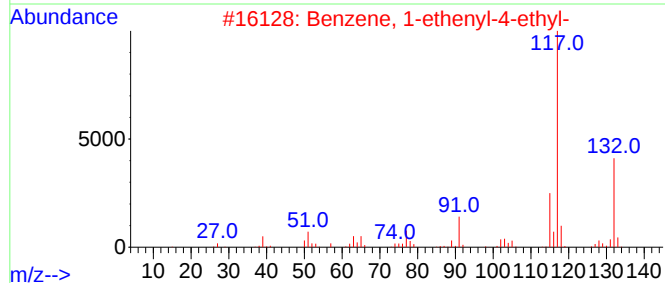
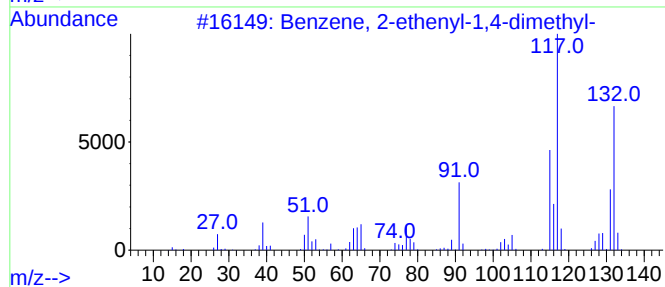
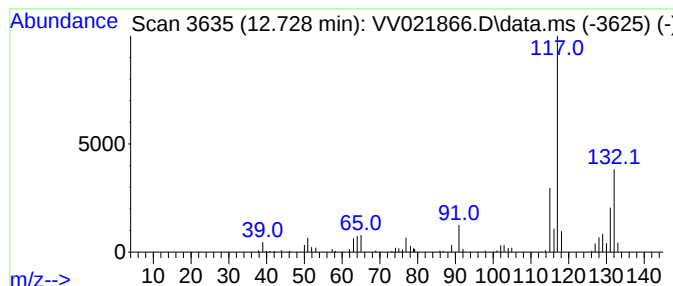
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Quant Title : VOC Analysis

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

Peak Number 5 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.728	4.62 ug/L	102882	1,4-Dichlorobenzene-d4	11.252

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
3			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	91
4			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
5			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081821\
Data File : VV021866.D
Acq On : 18 Aug 2021 15:28
Operator : SY/MD
Sample : M3448-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKD5

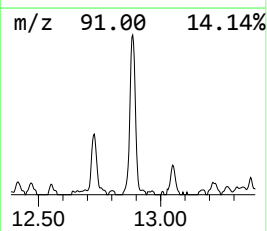
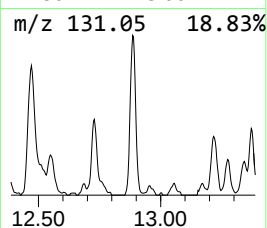
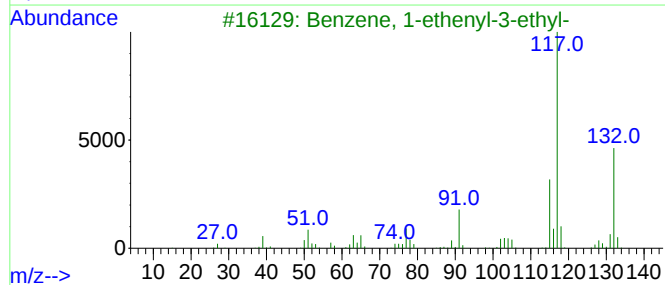
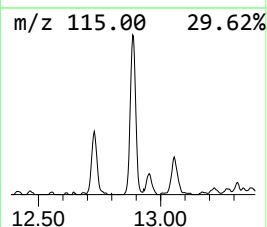
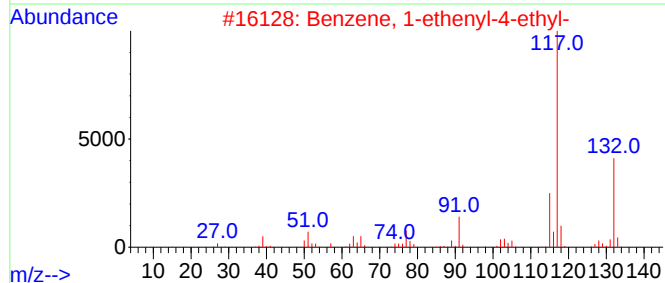
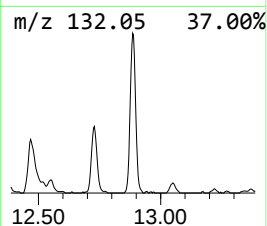
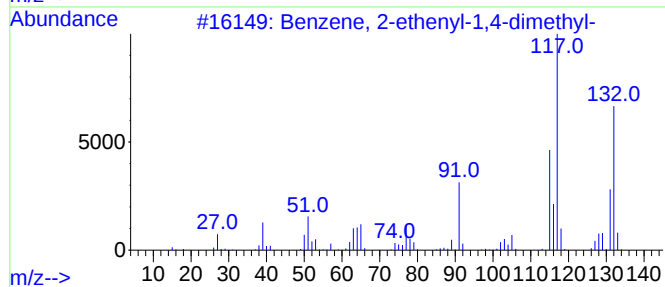
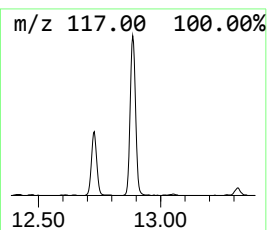
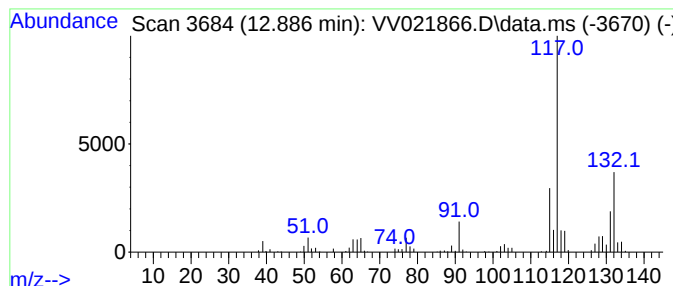
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Quant Title : VOC Analysis

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

Peak Number 6 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.886	11.65 ug/L	259298	1,4-Dichlorobenzene-d4	11.252

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
2			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
3			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
4			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
5			3-Phenylbut-1-ene	132	C10H12	000934-10-1	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081821\
Data File : VV021866.D
Acq On : 18 Aug 2021 15:28
Operator : SY/MD
Sample : M3448-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKD5

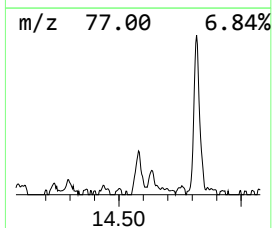
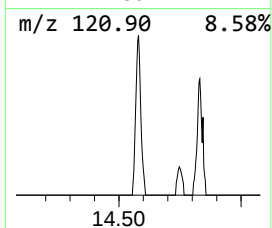
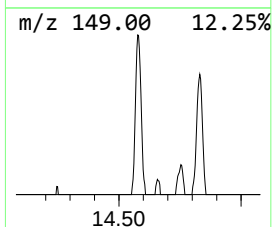
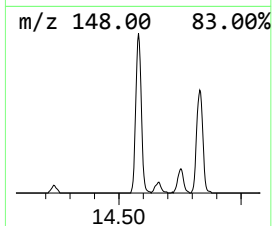
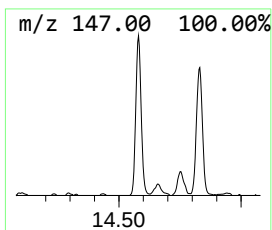
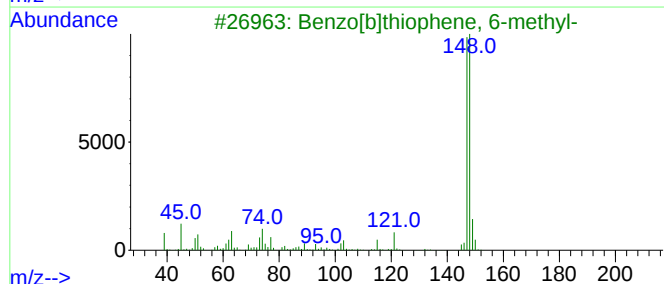
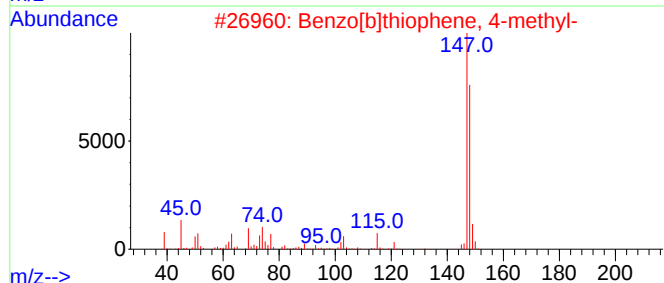
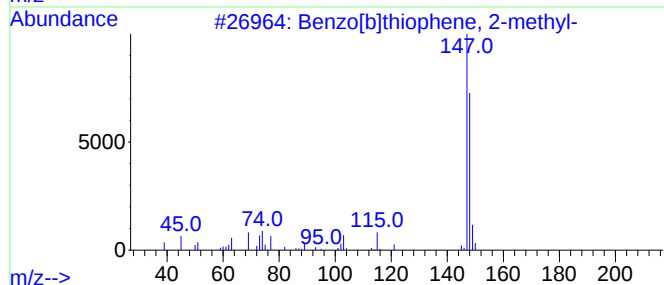
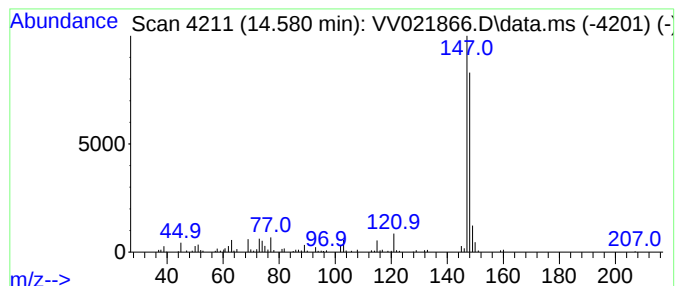
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Quant Title : VOC Analysis

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

Peak Number 7 Benzo[b]thiophene, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.580	4.55 ug/L	101178	1,4-Dichlorobenzene-d4	11.252

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	91
2			Benzo[b]thiophene, 4-methyl-	148	C9H8S	014315-11-8	91
3			Benzo[b]thiophene, 6-methyl-	148	C9H8S	016587-47-6	91
4			Benzo[b]thiophene, 5-methyl-	148	C9H8S	014315-14-1	91
5			Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081821\
Data File : VV021866.D
Acq On : 18 Aug 2021 15:28
Operator : SY/MD
Sample : M3448-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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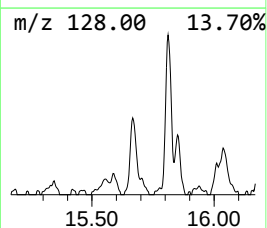
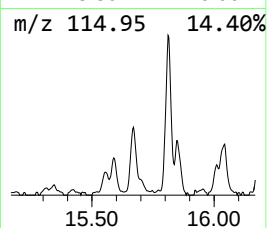
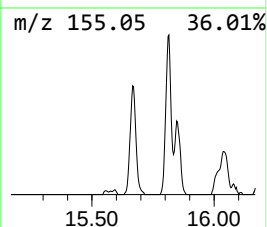
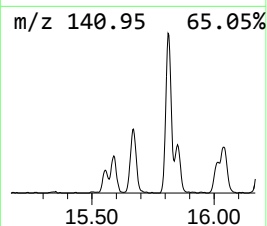
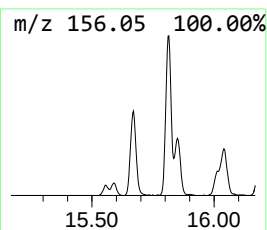
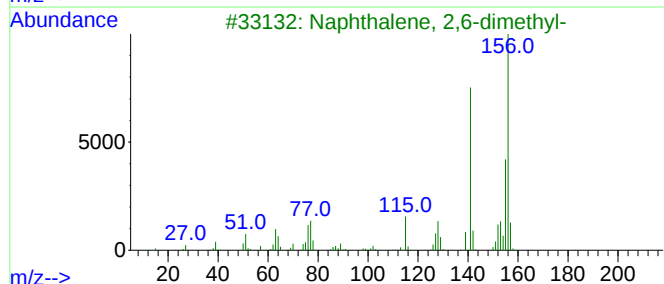
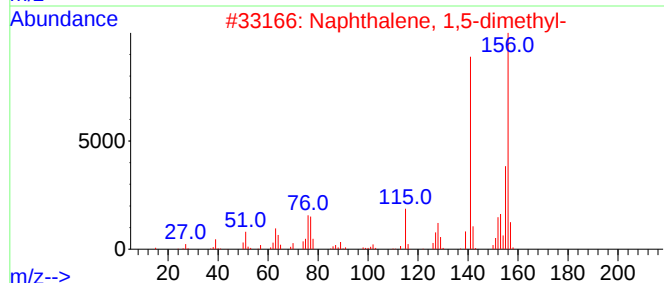
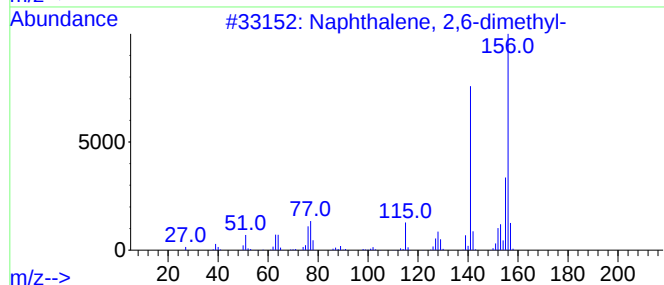
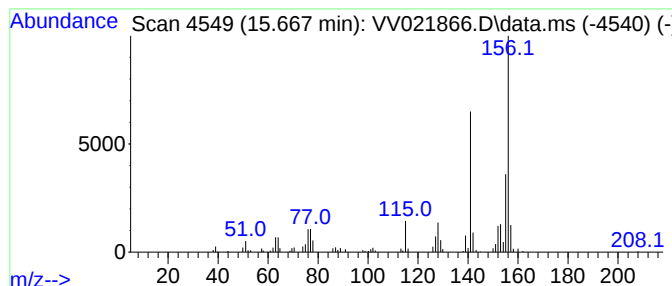
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Quant Title : VOC Analysis

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

Peak Number 10 Naphthalene, 2,6-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.667	7.40 ug/L	164658	1,4-Dichlorobenzene-d4	11.252

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	98
2			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	98
3			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	98
4			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
5			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081821\
Data File : VV021866.D
Acq On : 18 Aug 2021 15:28
Operator : SY/MD
Sample : M3448-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKD5

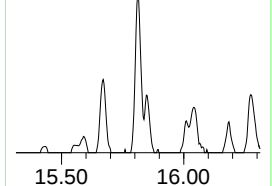
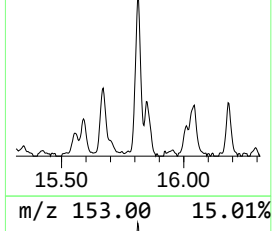
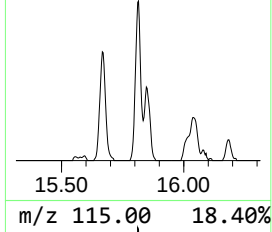
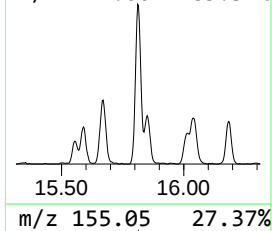
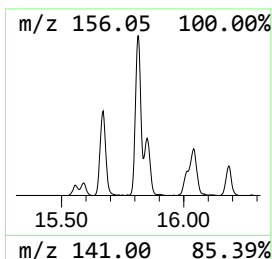
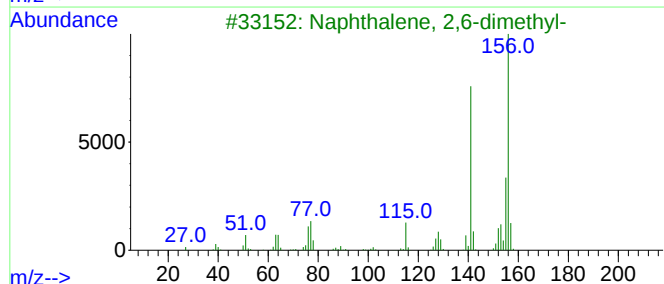
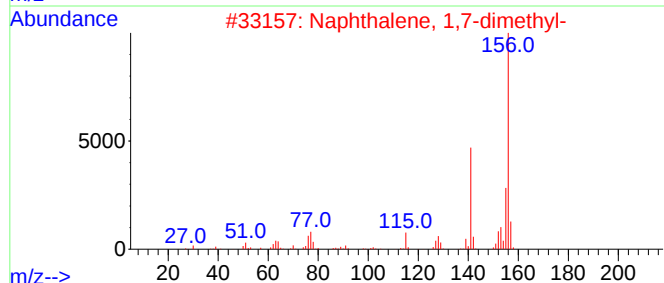
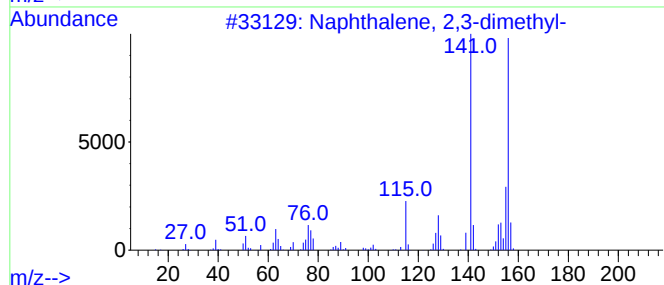
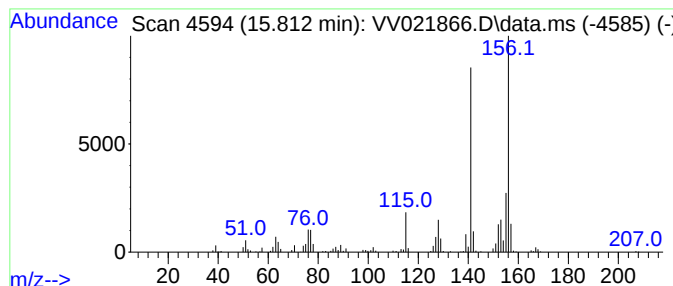
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Quant Title : VOC Analysis

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

Peak Number 11 Naphthalene, 2,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.812	11.37 ug/L	252989	1,4-Dichlorobenzene-d4	11.252

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
3			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
4			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
5			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081821\
Data File : VV021866.D
Acq On : 18 Aug 2021 15:28
Operator : SY/MD
Sample : M3448-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
DBKD5

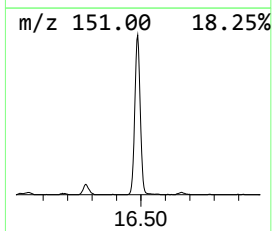
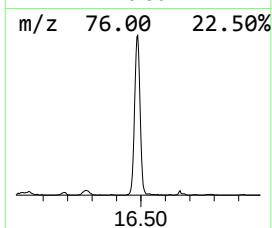
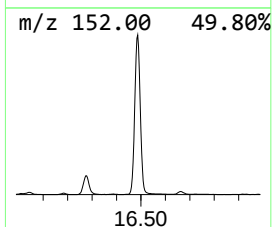
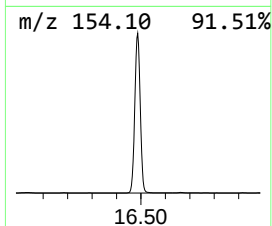
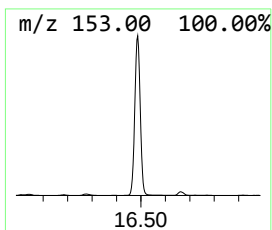
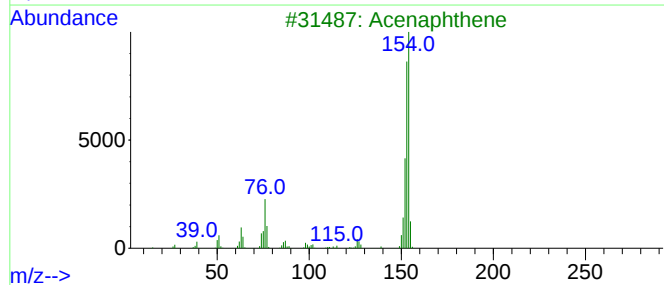
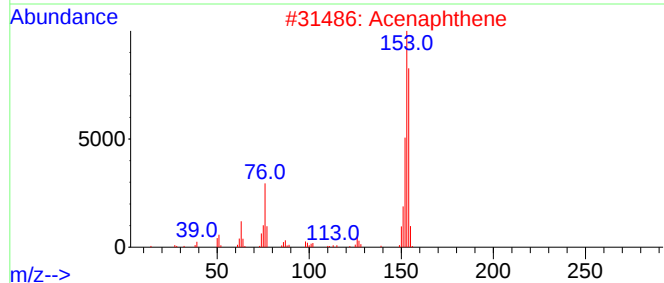
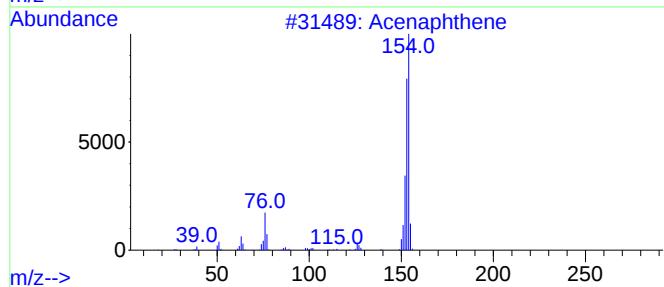
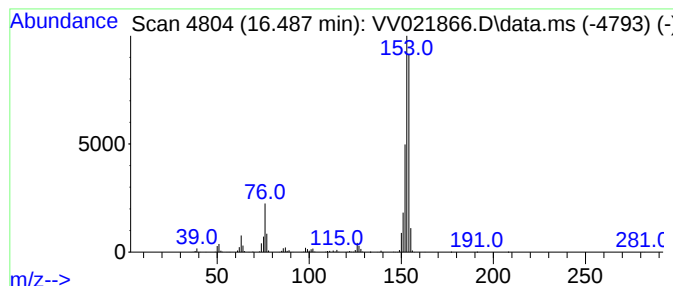
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Quant Title : VOC Analysis

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

Peak Number 12 Naphthalene, 2-ethenyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.487	62.81 ug/L	1398030	1,4-Dichlorobenzene-d4	11.252

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acenaphthene	154	C12H10	000083-32-9	94
2			Acenaphthene	154	C12H10	000083-32-9	89
3			Acenaphthene	154	C12H10	000083-32-9	72
4			Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	64
5			Hexa-2,4-diyn-1-ylbenzene	154	C12H10	000520-74-1	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV081821\
Data File : VV021866.D
Acq On : 18 Aug 2021 15:28
Operator : SY/MD
Sample : M3448-03
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
DBKD5

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM081621WMA.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
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1-Phenyl-1-butene	12.117	2.9	ug/L	65319	3	11.252	1112990	50.0
Benzofuran, 2-m...	12.468	2.7	ug/L	59197	3	11.252	1112990	50.0
Benzene, 2-ethe...	12.728	4.6	ug/L	102882	3	11.252	1112990	50.0
Benzene, 1-ethe...	12.886	11.7	ug/L	259298	3	11.252	1112990	50.0
Benzo[b]thiophe...	14.580	4.5	ug/L	101178	3	11.252	1112990	50.0
Naphthalene, 2,...	15.667	7.4	ug/L	164658	3	11.252	1112990	50.0
Naphthalene, 2,...	15.812	11.4	ug/L	252989	3	11.252	1112990	50.0
Naphthalene, 2-...	16.487	62.8	ug/L	1398030	3	11.252	1112990	50.0