

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV071520\
 Data File : VV071506.D
 Acq On : 15 Jul 2020 12:55
 Operator : SY/MD
 Sample : L3281-20
 Misc : 25.0mL/MSVOA V/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 GAXY0

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.1

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

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

Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR071420WMA.M

Title : TRACE VOA SOM01.0

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.116	3	8	22	rVB2	60732	68895	4.79%	0.664%
2	1.226	36	42	56	rVB	36336	36052	2.51%	0.347%
3	1.316	61	70	80	rBV2	90022	106544	7.41%	1.027%
4	1.579	144	152	164	rBV	54944	66272	4.61%	0.639%
5	1.647	165	173	184	rBV	55671	70780	4.92%	0.682%
6	1.824	220	228	239	rVB4	15737	23987	1.67%	0.231%
7	2.126	311	322	337	rBV	132980	198521	13.80%	1.913%
8	2.315	373	381	392	rVB	10311	16996	1.18%	0.164%
9	2.598	462	469	479	rVB4	11930	21732	1.51%	0.209%
10	2.782	513	526	539	rBV4	14234	29369	2.04%	0.283%
11	3.061	601	613	629	rBV2	15276	32099	2.23%	0.309%
12	3.788	827	839	853	rVB5	17153	42015	2.92%	0.405%
13	3.939	866	886	919	rBV2	568988	1438702	100.00%	13.866%
14	4.383	1008	1024	1049	rBV	84580	216704	15.06%	2.089%
15	4.701	1104	1123	1140	rBV6	26123	71464	4.97%	0.689%
16	4.936	1182	1196	1220	rBV6	12343	38600	2.68%	0.372%
17	5.074	1222	1239	1255	rBV3	183545	454617	31.60%	4.381%
18	5.206	1267	1280	1289	rBV9	8725	19839	1.38%	0.191%
19	5.277	1294	1302	1313	rVV8	8149	16973	1.18%	0.164%
20	5.351	1313	1325	1337	rVB8	14192	33265	2.31%	0.321%
21	5.518	1360	1377	1390	rBV3	23063	49372	3.43%	0.476%
22	5.643	1403	1416	1451	rBV	174774	373572	25.97%	3.600%
23	6.097	1544	1557	1567	rBV2	105406	218220	15.17%	2.103%
24	6.155	1568	1575	1591	rVB4	50969	108088	7.51%	1.042%
25	6.486	1665	1678	1691	rVB7	8715	19593	1.36%	0.189%
26	6.653	1723	1730	1742	rVB7	7216	15644	1.09%	0.151%
27	6.939	1810	1819	1827	rBV3	17709	29741	2.07%	0.287%
28	7.010	1828	1841	1856	rVB	55021	105020	7.30%	1.012%
29	7.103	1862	1870	1888	rVV5	13656	28653	1.99%	0.276%
30	7.187	1888	1896	1912	rVB7	9460	22910	1.59%	0.221%
31	7.286	1918	1927	1933	rBV5	23160	43172	3.00%	0.416%
32	7.338	1933	1943	1958	rVV	230243	405975	28.22%	3.913%
33	7.415	1958	1967	1976	rVV2	20707	38685	2.69%	0.373%
34	7.589	2012	2021	2030	rBV3	36414	60199	4.18%	0.580%

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 ALS Vial : 9 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.1

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Filtering: 5

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Max Peaks: 100

Peak Location: TOP

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

Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR071420WMA.M
 Title : TRACE VOA SOM01.0

35	7.646	2031	2039	2045	rBV	31673	51442	3.58%	0.496%
36	7.807	2080	2089	2099	rBV5	9576	16903	1.17%	0.163%
37	8.109	2171	2183	2209	rBV	223366	410375	28.52%	3.955%
38	8.312	2238	2246	2255	rVB3	18102	29050	2.02%	0.280%
39	8.373	2255	2265	2273	rBV3	20283	35159	2.44%	0.339%
40	8.531	2301	2314	2324	rBV3	10483	22866	1.59%	0.220%
41	8.611	2336	2339	2353	rVV10	9501	22698	1.58%	0.219%
42	8.688	2355	2363	2373	rVB4	19335	33550	2.33%	0.323%
43	8.810	2388	2401	2410	rBV2	21332	34580	2.40%	0.333%
44	8.875	2410	2421	2435	rBV	261651	428199	29.76%	4.127%
45	9.032	2457	2470	2484	rBV	112586	182602	12.69%	1.760%
46	9.158	2490	2509	2522	rBV	370558	620796	43.15%	5.983%
47	9.225	2522	2530	2545	rVB3	56124	91308	6.35%	0.880%
48	9.495	2598	2614	2622	rBV8	14948	30481	2.12%	0.294%
49	9.566	2628	2636	2653	rVB2	21719	38739	2.69%	0.373%
50	9.730	2680	2687	2697	rVB6	19659	32996	2.29%	0.318%
51	9.820	2703	2715	2724	rBV8	17496	34206	2.38%	0.330%
52	9.952	2745	2756	2765	rBV2	18605	33860	2.35%	0.326%
53	10.029	2774	2780	2790	rVB7	10683	16993	1.18%	0.164%
54	10.142	2809	2815	2827	rVB7	10711	19911	1.38%	0.192%
55	10.238	2832	2845	2864	rBV	86309	159620	11.09%	1.538%
56	10.376	2876	2888	2906	rBV2	36613	66637	4.63%	0.642%
57	10.470	2906	2917	2933	rVV	159646	345784	24.03%	3.333%
58	10.560	2936	2945	2953	rVV	97146	160360	11.15%	1.545%
59	10.601	2953	2958	2970	rVB2	30895	45979	3.20%	0.443%
60	10.759	2996	3007	3025	rVB	79472	133953	9.31%	1.291%
61	10.936	3042	3062	3085	rBV	363001	560425	38.95%	5.401%
62	11.177	3128	3137	3143	rVV9	9734	16450	1.14%	0.159%
63	11.215	3143	3149	3155	rVV2	15879	21125	1.47%	0.204%
64	11.270	3155	3166	3183	rVV	308543	479125	33.30%	4.618%
65	11.357	3183	3193	3211	rVB	115623	181043	12.58%	1.745%
66	11.553	3242	3254	3262	rBV2	81039	150569	10.47%	1.451%
67	11.595	3263	3267	3274	rVV2	38654	56739	3.94%	0.547%
68	11.646	3274	3283	3307	rVB2	289043	541496	37.64%	5.219%
69	11.839	3338	3343	3355	rVB2	16569	26004	1.81%	0.251%
70	11.932	3363	3372	3377	rBV	31802	47468	3.30%	0.457%
71	11.961	3378	3381	3392	rVB2	32215	40177	2.79%	0.387%

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Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % of largest Peak
 Start Thrs: 0.1 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

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 Title : TRACE VOA SOM01.0

72	12.039	3393	3405	3412	rBV	49464	79535	5.53%	0.767%
73	12.138	3426	3436	3447	rBV	45856	84909	5.90%	0.818%
74	12.338	3489	3498	3504	rVV3	13050	18403	1.28%	0.177%
75	12.434	3520	3528	3537	rVB	25419	38787	2.70%	0.374%
76	12.489	3537	3545	3559	rVB	48750	76960	5.35%	0.742%
77	12.572	3563	3571	3576	rBV7	11916	17610	1.22%	0.170%
78	12.749	3617	3626	3640	rVB3	36630	61144	4.25%	0.589%
79	12.907	3665	3675	3686	rVB2	66765	105220	7.31%	1.014%
80	13.071	3719	3726	3735	rVB7	10583	17442	1.21%	0.168%
81	13.190	3753	3763	3770	rBV5	8914	15139	1.05%	0.146%
82	13.238	3770	3778	3785	rBV	15113	21916	1.52%	0.211%
83	13.293	3787	3795	3802	rBV4	19659	28976	2.01%	0.279%
84	13.527	3859	3868	3878	rVB	28855	44787	3.11%	0.432%
85	14.119	4041	4052	4064	rBV	9814	23201	1.61%	0.224%

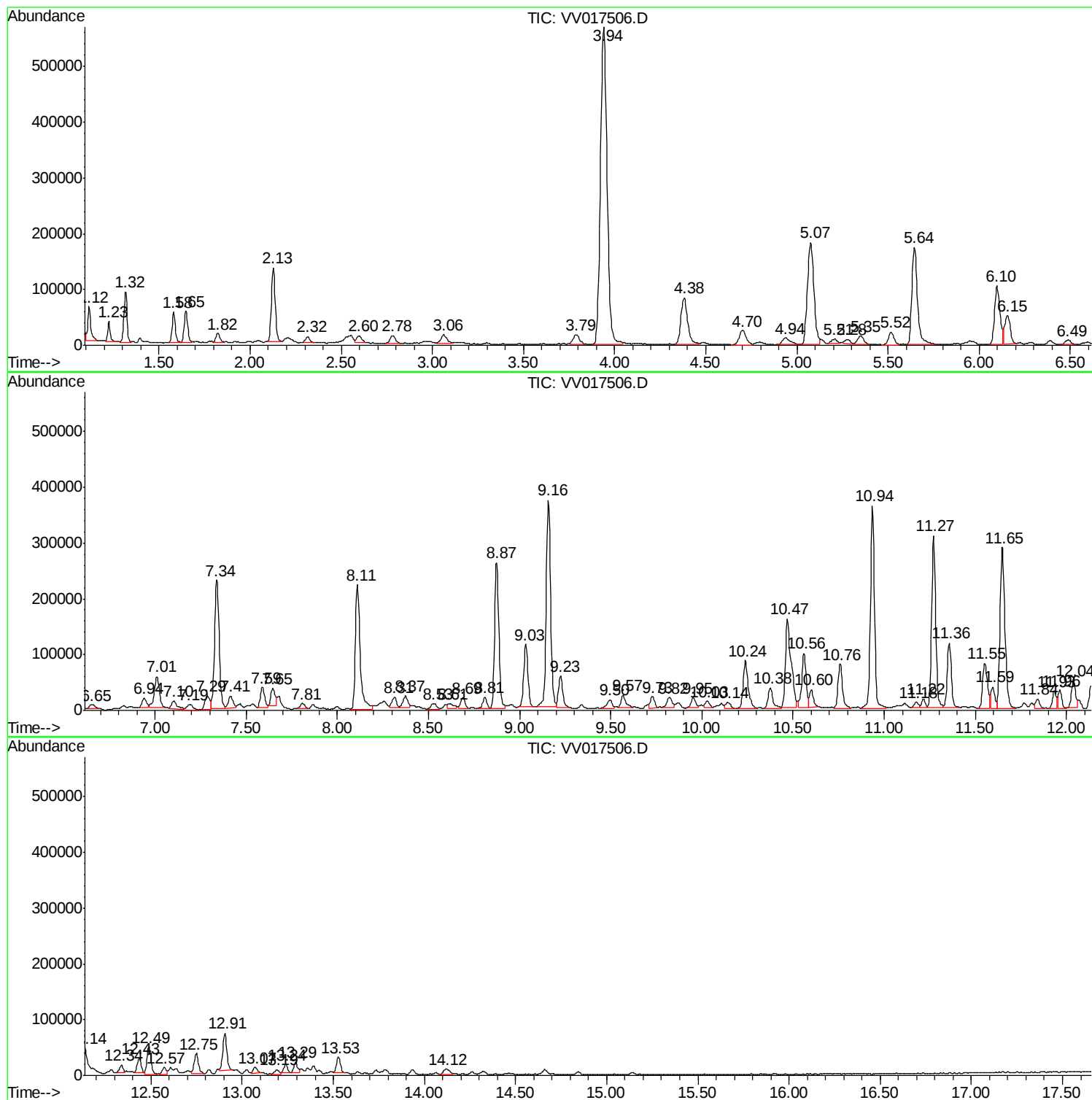
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ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
GAXY0

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR071420WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV071520\
Data File : VV017506.D
Acq On : 15 Jul 2020 12:55
Operator : SY/MD
Sample : L3281-20
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ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAXY0

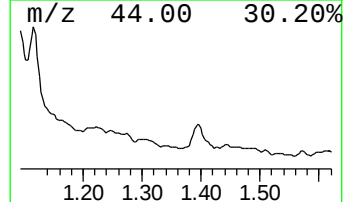
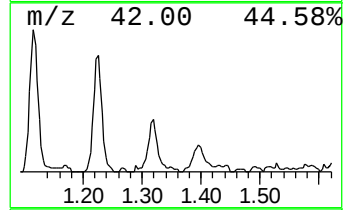
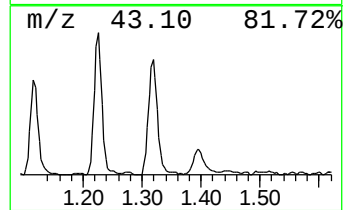
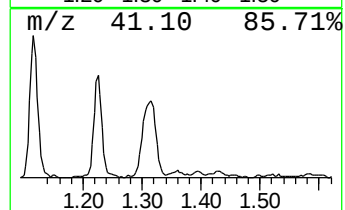
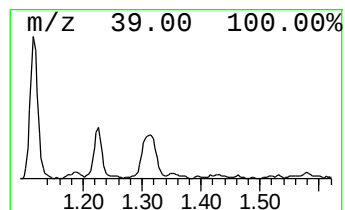
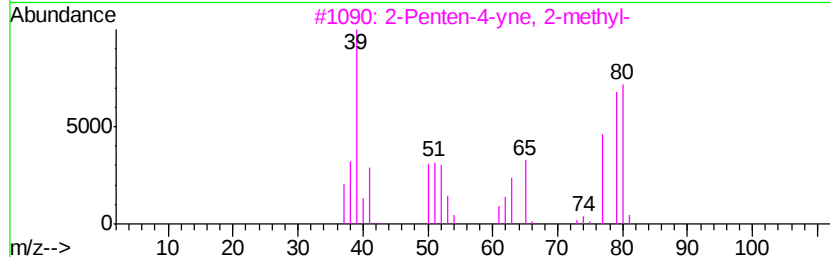
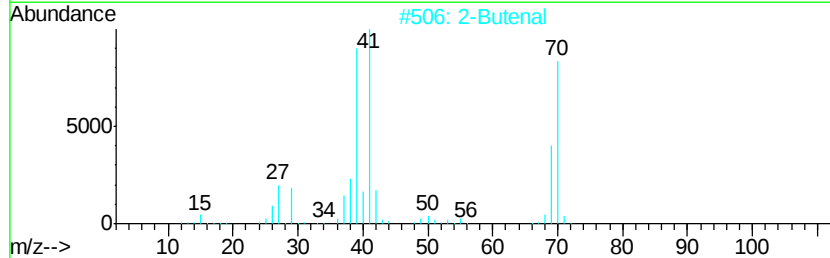
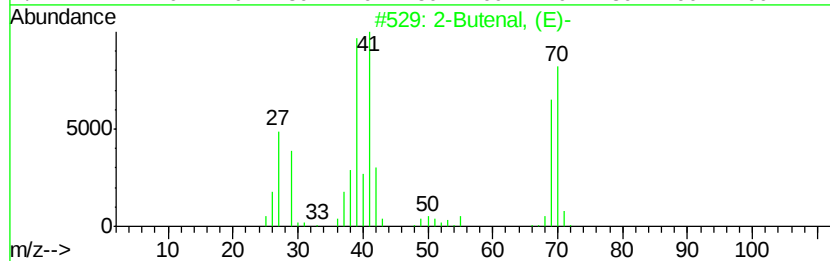
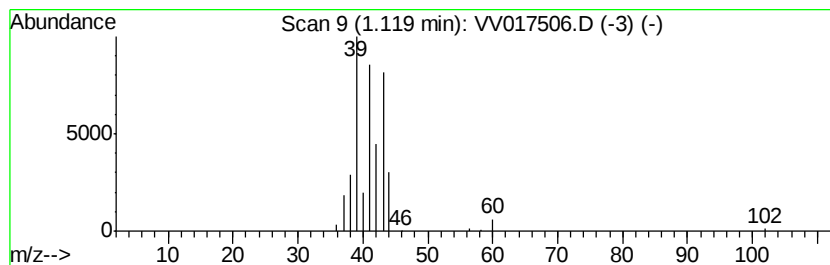
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Butenal, (E)- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.12	0.92 ug/L	68895	1,4-Difluorobenzene	5.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Butenal, (E)-	70	C4H6O	000123-73-9	72
2		2-Butenal	70	C4H6O	004170-30-3	39
3		2-Penten-4-yne, 2-methyl-	80	C6H8	001595-53-5	9
4		2-Butynal	68	C4H4O	001119-19-3	9
5		2-Ethyl-3-vinylloxirane	98	C6H10O	034485-78-4	9



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Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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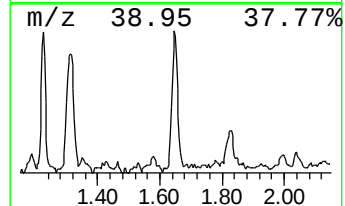
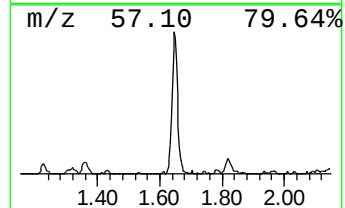
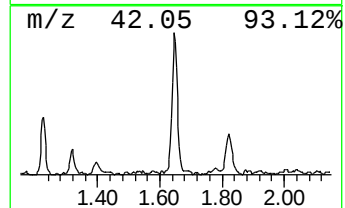
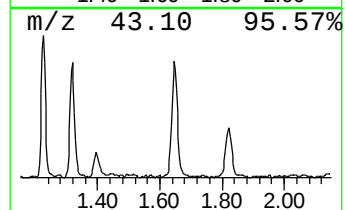
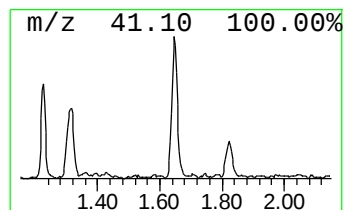
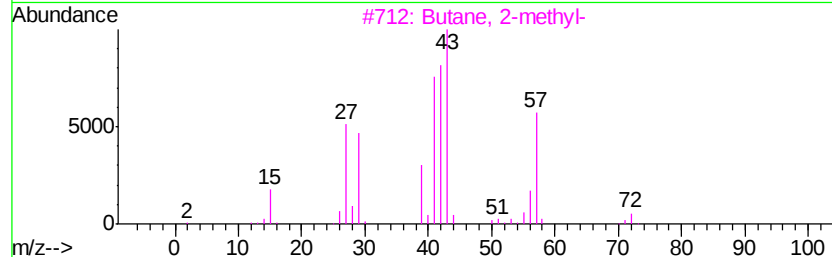
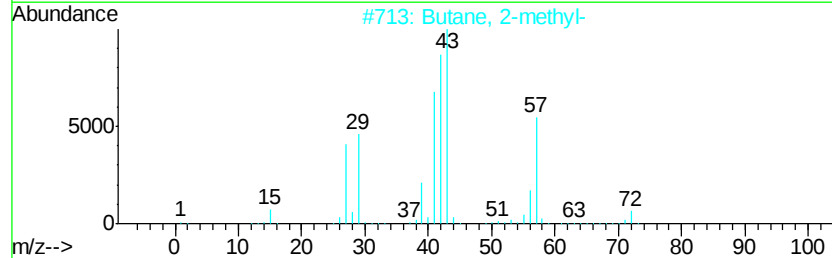
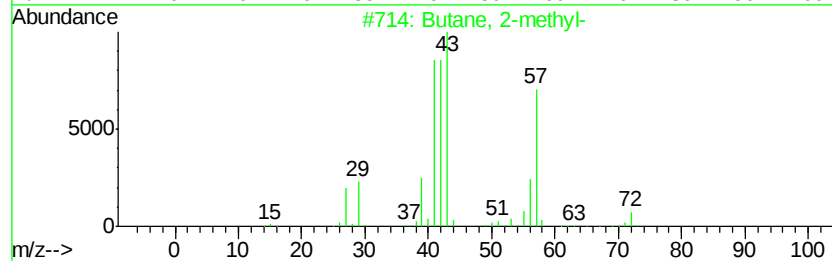
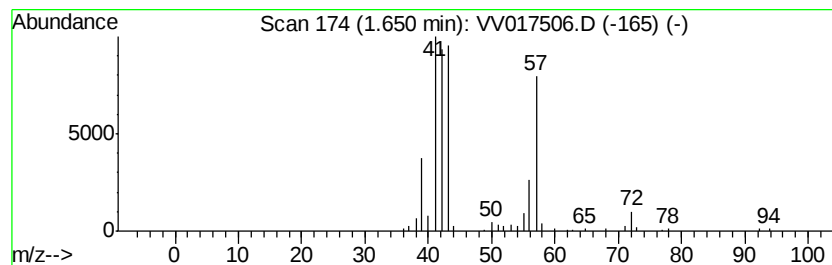
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR071420WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.65	0.95 ug/L	70780	1,4-Difluorobenzene	5.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methyl-	72	C5H12	000078-78-4	90
2		Butane, 2-methyl-	72	C5H12	000078-78-4	80
3		Butane, 2-methyl-	72	C5H12	000078-78-4	80
4		3-Buten-1-ol	72	C4H8O	000627-27-0	37
5		2(3H)-Furanone, dihydro-4-methyl-	100	C5H8O2	001679-49-8	25



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Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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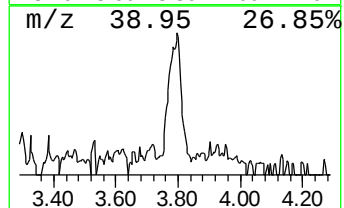
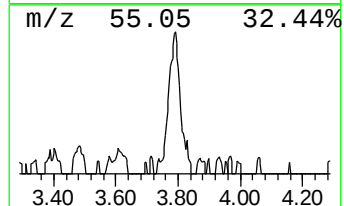
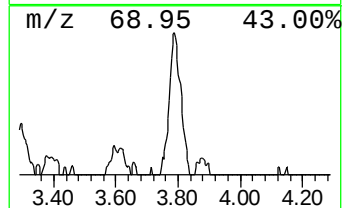
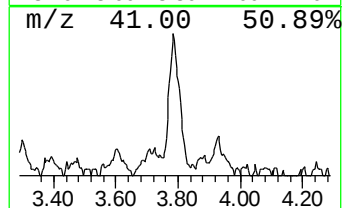
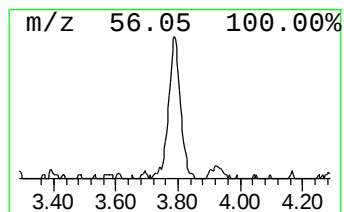
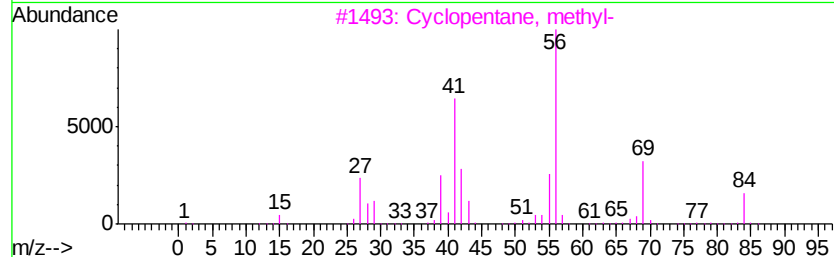
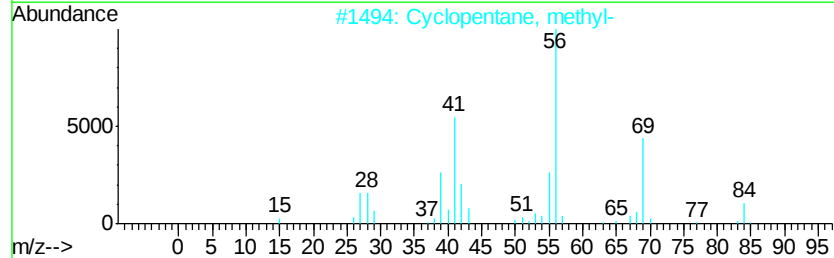
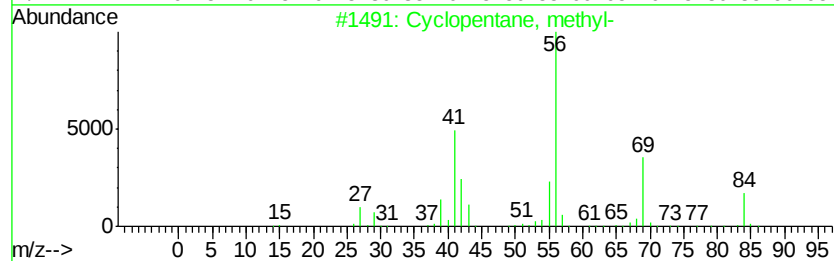
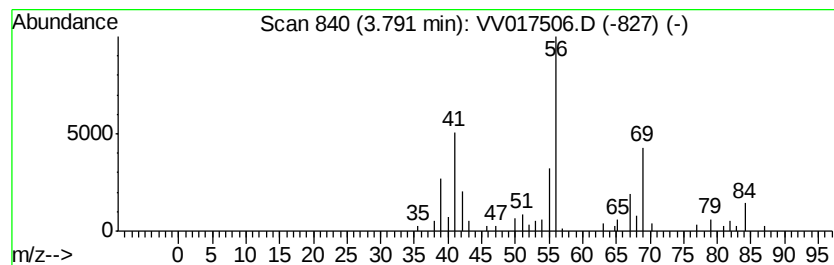
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 (DEL) Alkane: Cyclic3.79 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.79	0.56 ug/L	42015	1,4-Difluorobenzene	5.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	72
2		Cyclopentane, methyl-	84	C6H12	000096-37-7	72
3		Cyclopentane, methyl-	84	C6H12	000096-37-7	72
4		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	59
5		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	58



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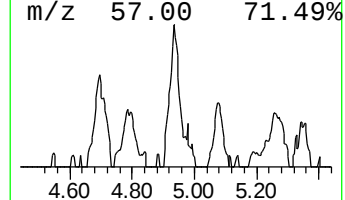
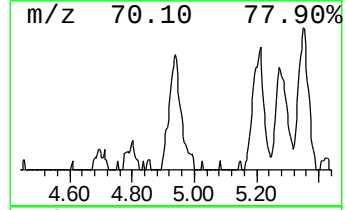
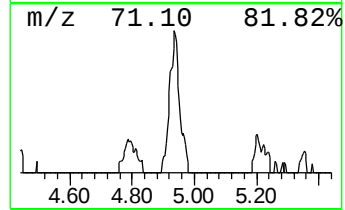
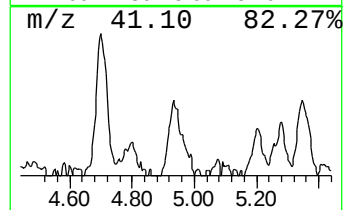
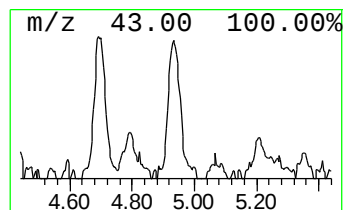
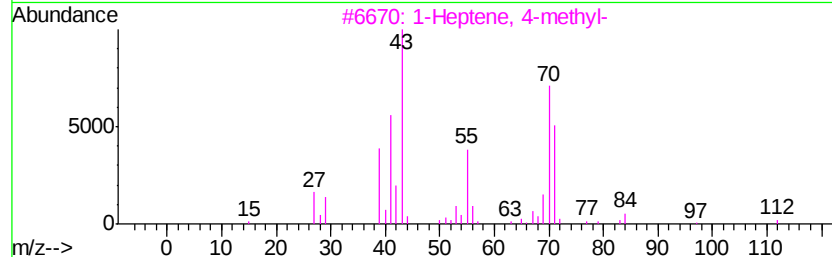
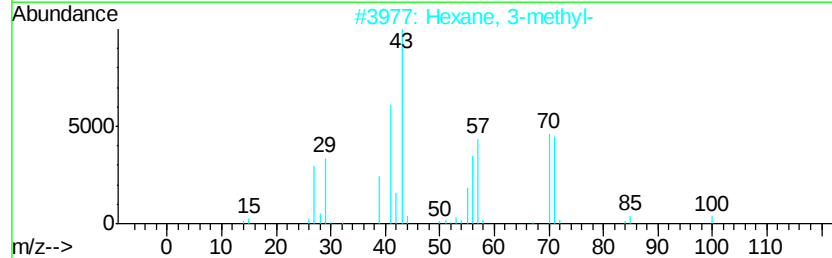
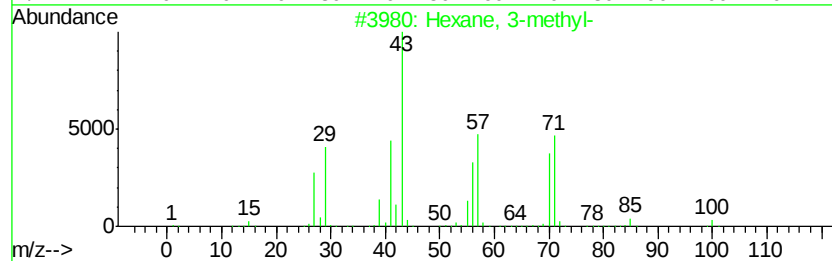
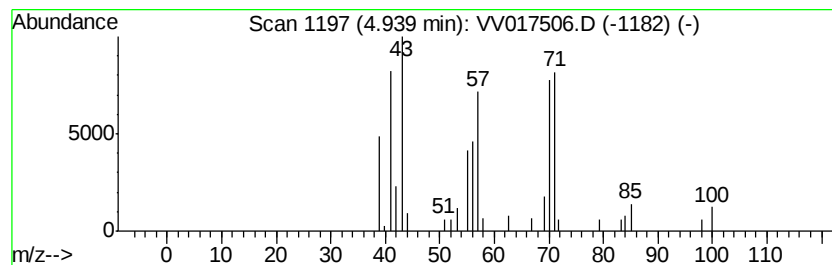
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.94	0.52 ug/L	38600	1,4-Difluorobenzene	5.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexane, 3-methyl-	100	C7H16	000589-34-4	53
2		Hexane, 3-methyl-	100	C7H16	000589-34-4	53
3		1-Heptene, 4-methyl-	112	C8H16	013151-05-8	47
4		Hexane, 3-methyl-	100	C7H16	000589-34-4	47
5		Trans-3-methylpent-3-ene-5-ol	100	C6H12O	030801-95-7	47



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV071520\
Data File : VV071506.D
Acq On : 15 Jul 2020 12:55
Operator : SY/MD
Sample : L3281-20
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAXY0

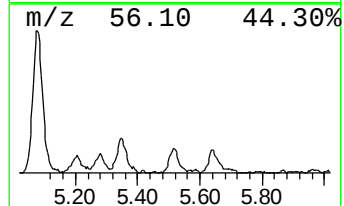
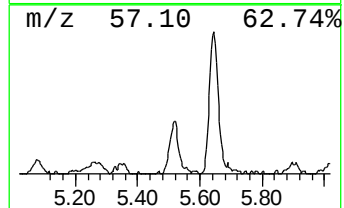
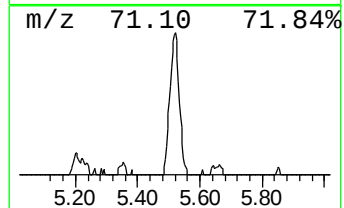
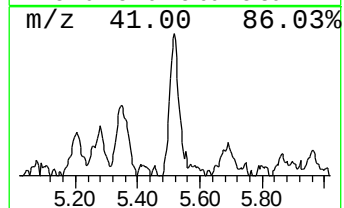
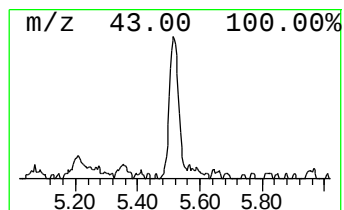
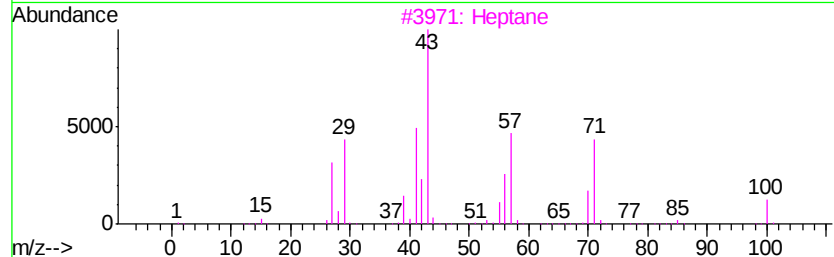
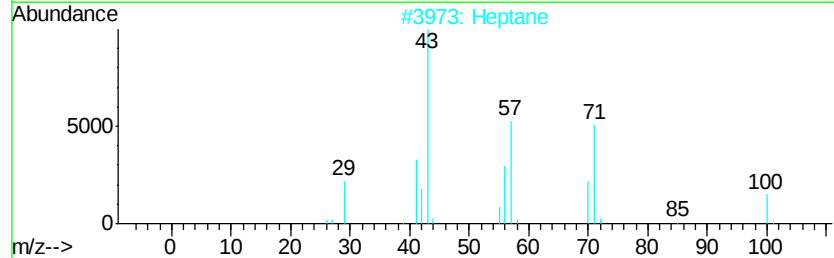
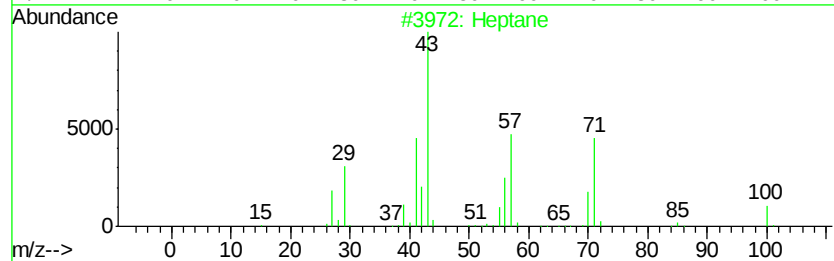
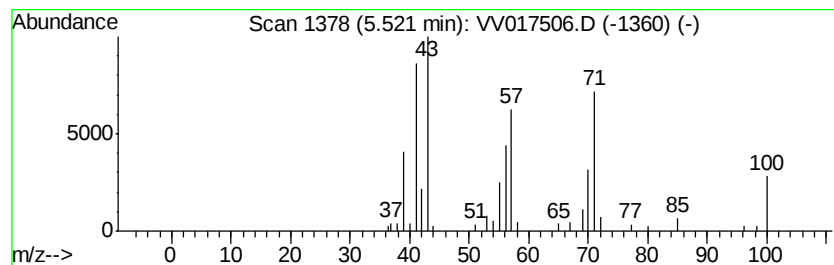
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.52	0.66 ug/L	49372	1,4-Difluorobenzene	5.64

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane		100	C7H16	000142-82-5	58
2	Heptane		100	C7H16	000142-82-5	53
3	Heptane		100	C7H16	000142-82-5	50
4	1-Hexene, 4-methyl-		98	C7H14	003769-23-1	50
5	Cyclopropane, butyl-		98	C7H14	000930-57-4	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV071520\
Data File : VV071506.D
Acq On : 15 Jul 2020 12:55
Operator : SY/MD
Sample : L3281-20
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 9 Sample Multiplier: 1

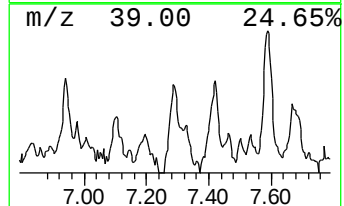
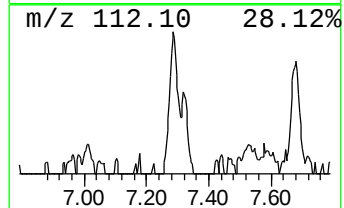
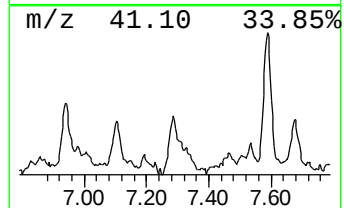
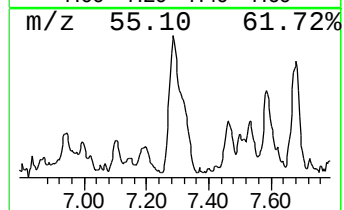
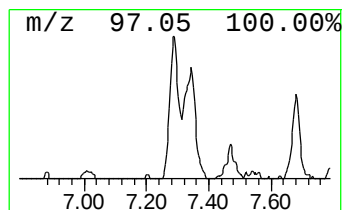
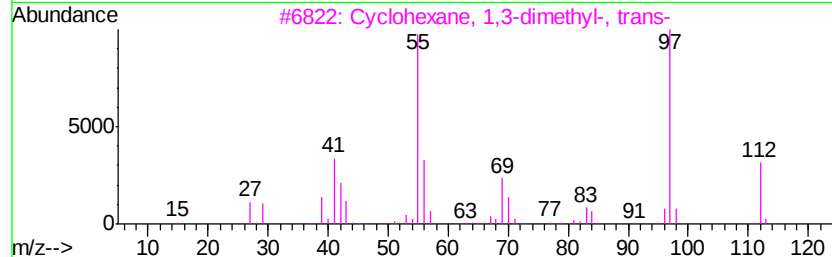
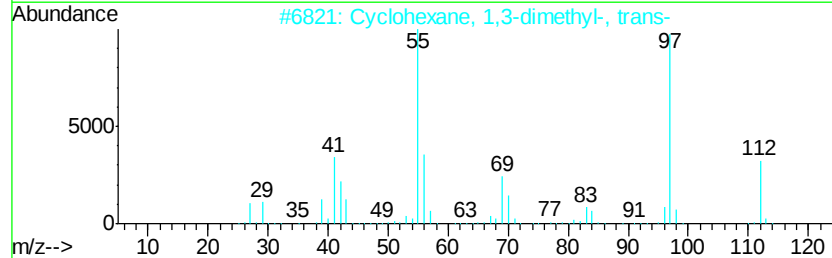
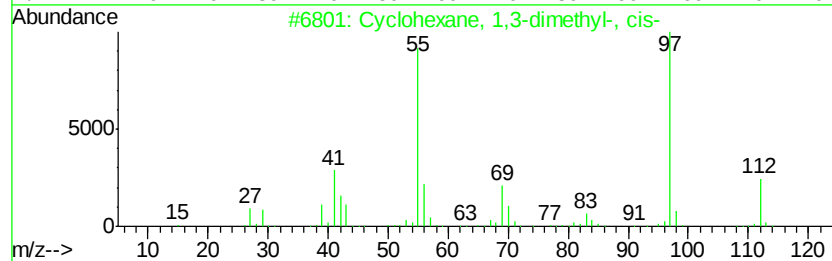
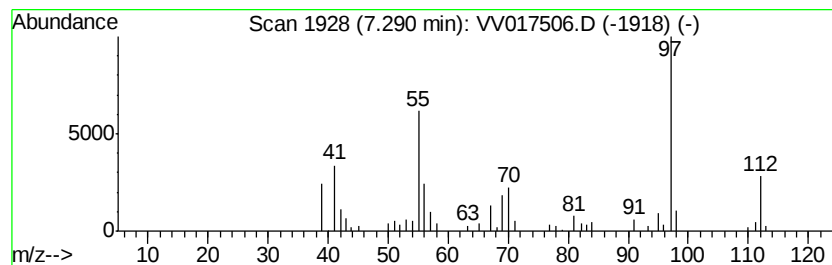
Instrument :
MSVOA_V
ClientSampleId :
GAXY0

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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Cyclic7.29 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.	
7.29	0.50 ug/L	43172	Chlorobenzene-d5	8.87	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cvclohexane, 1,3-dimethyl-, cis-	112	C8H16	000638-04-0	64
2	Cvclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	64
3	Cvclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	64
4	Cvclohexane, 1,3-dimethyl-, cis-	112	C8H16	000638-04-0	64
5	Cvclohexane, 1,3-dimethyl-, cis-	112	C8H16	000638-04-0	59



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV071520\
Data File : VV017506.D
Acq On : 15 Jul 2020 12:55
Operator : SY/MD
Sample : L3281-20
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAXY0

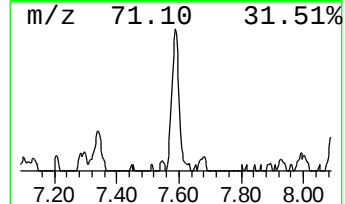
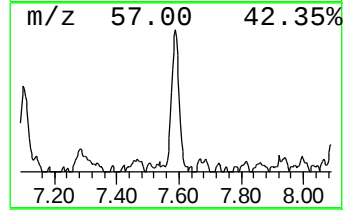
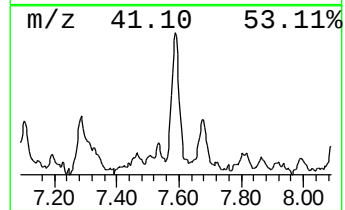
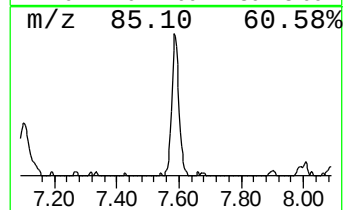
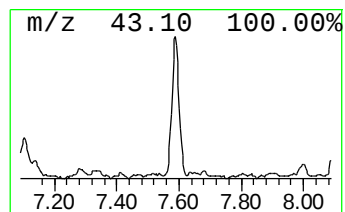
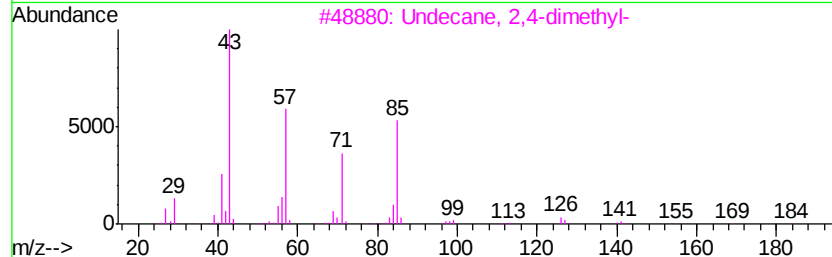
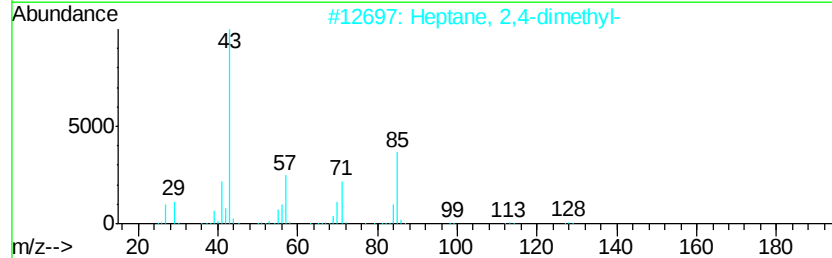
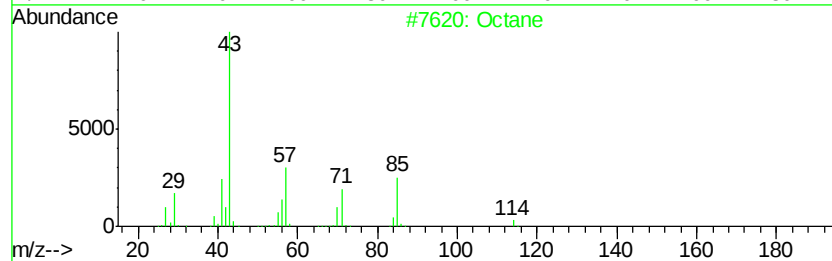
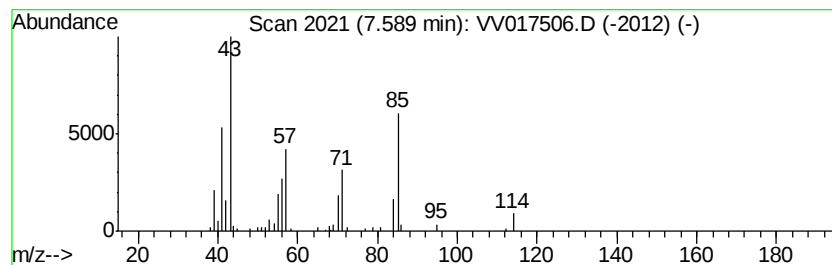
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	0.70 ug/L	60199	Chlorobenzene-d5	8.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octane	114	C8H18	000111-65-9	80
2		Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	64
3		Undecane, 2,4-dimethyl-	184	C13H28	017312-80-0	59
4		Octane	114	C8H18	000111-65-9	58
5		Heptane, 2,3-dimethyl-	128	C9H20	003074-71-3	50



Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV071520\
Data File : VV017506.D
Acq On : 15 Jul 2020 12:55
Operator : SY/MD
Sample : L3281-20
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAXY0

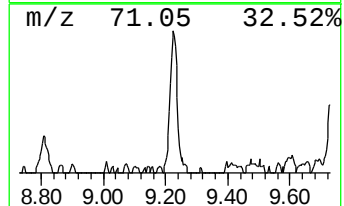
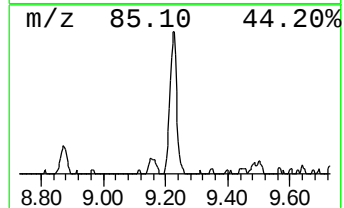
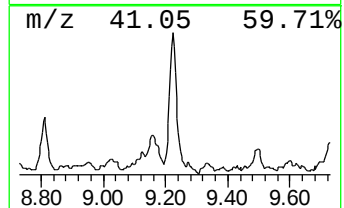
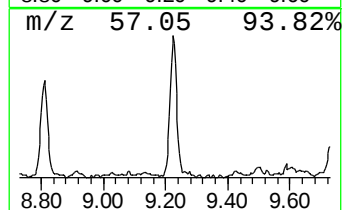
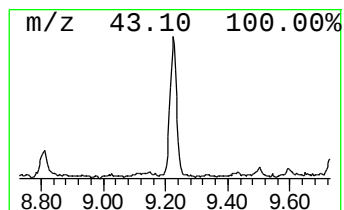
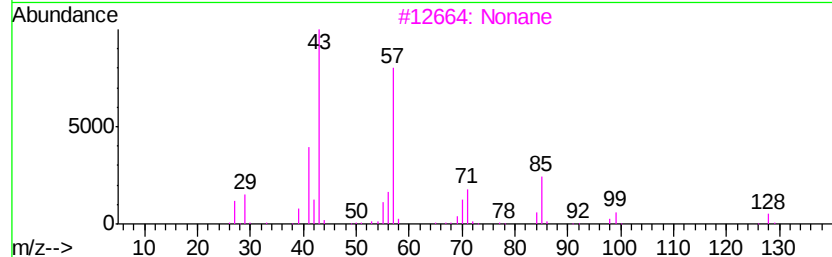
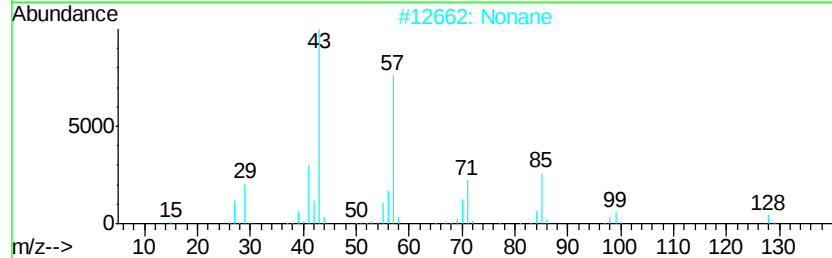
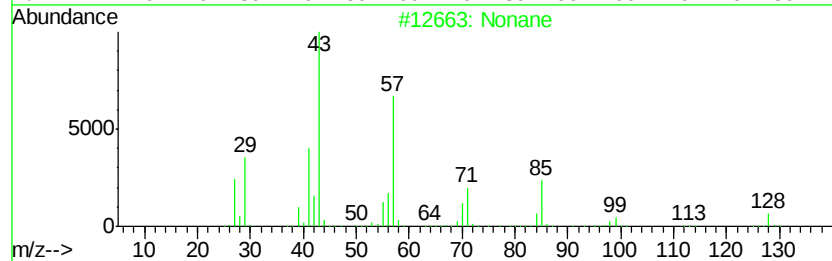
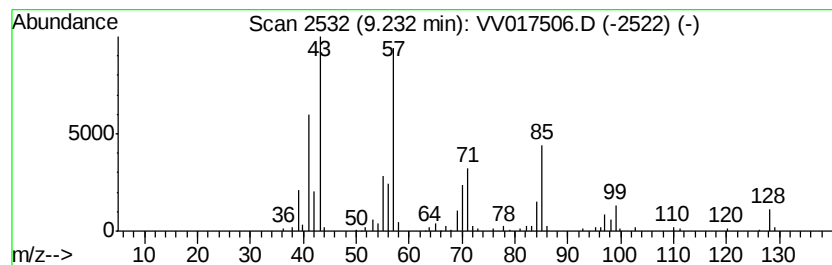
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.23	1.07 ug/L	91308	Chlorobenzene-d5	8.87

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane			128	C9H20	000111-84-2	93
2	Nonane			128	C9H20	000111-84-2	93
3	Nonane			128	C9H20	000111-84-2	81
4	Hexane, 2,4-dimethyl-			114	C8H18	000589-43-5	50
5	Octane			114	C8H18	000111-65-9	50



Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV071520\
Data File : VV017506.D
Acq On : 15 Jul 2020 12:55
Operator : SY/MD
Sample : L3281-20
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAXY0

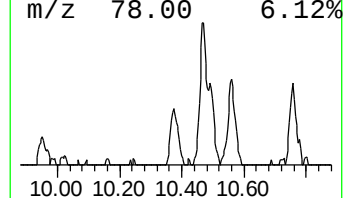
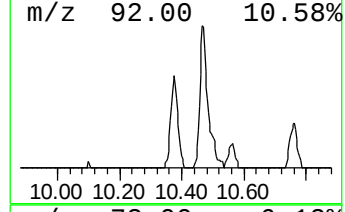
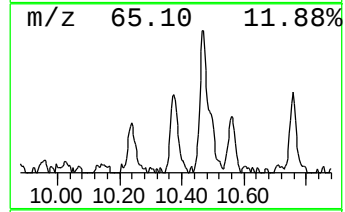
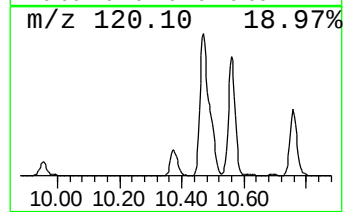
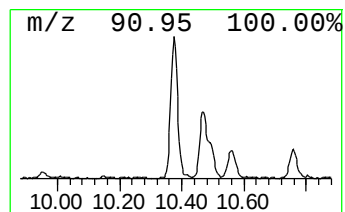
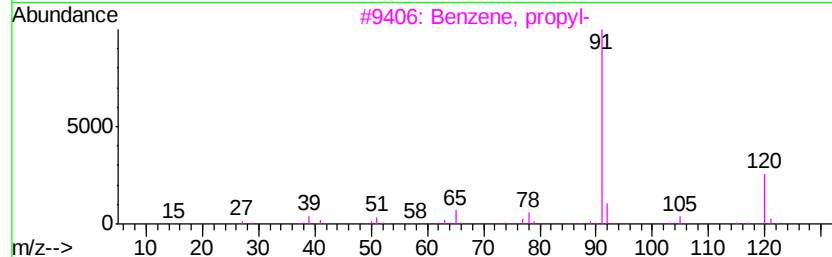
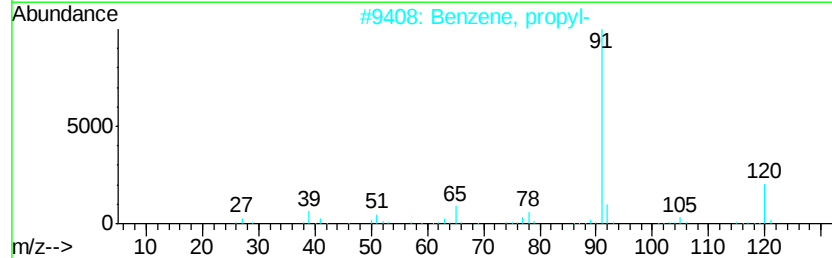
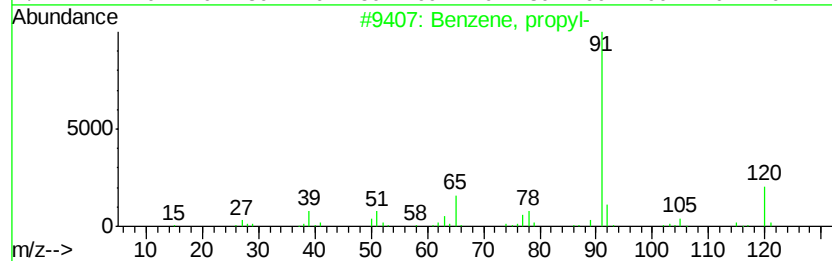
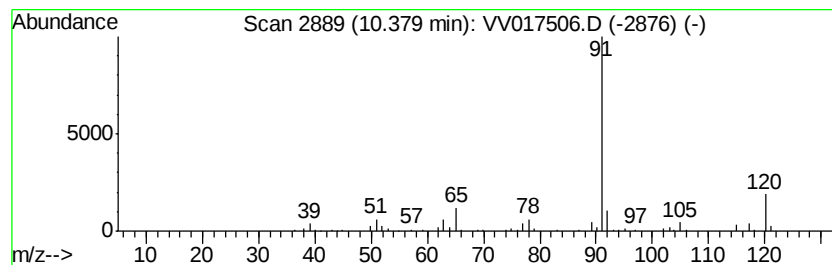
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Benzene, propyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.38	0.70 ug/L	66637	1,4-Dichlorobenzene-d4	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, propyl-	120	C9H12	000103-65-1	94
2		Benzene, propyl-	120	C9H12	000103-65-1	87
3		Benzene, propyl-	120	C9H12	000103-65-1	72
4		2,4,6-Cycloheptatrien-1-one, 4-m...	120	C8H8O	1000327-34-9	53
5		Benzeneacetaldehyde	120	C8H8O	000122-78-1	46



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV071520\
Data File : VV017506.D
Acq On : 15 Jul 2020 12:55
Operator : SY/MD
Sample : L3281-20
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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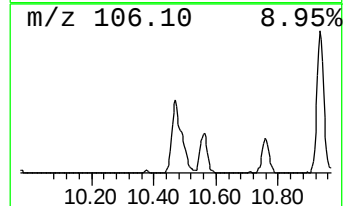
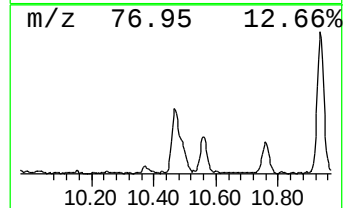
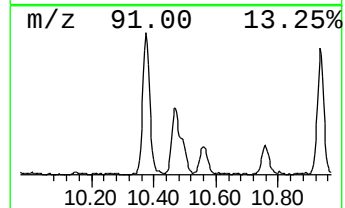
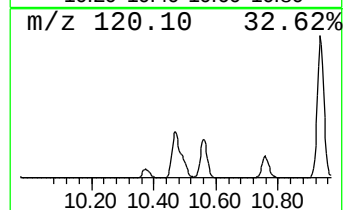
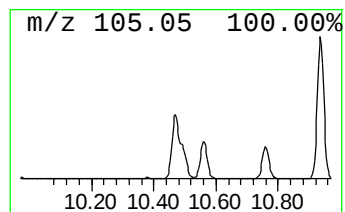
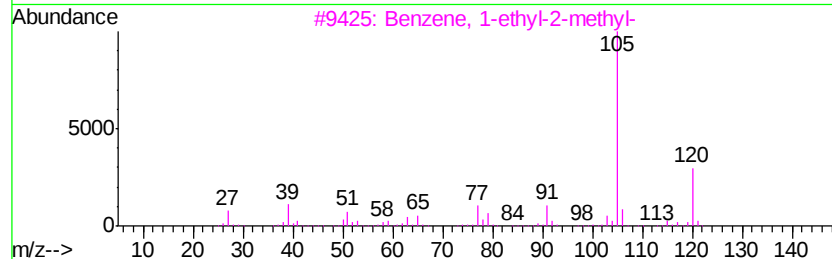
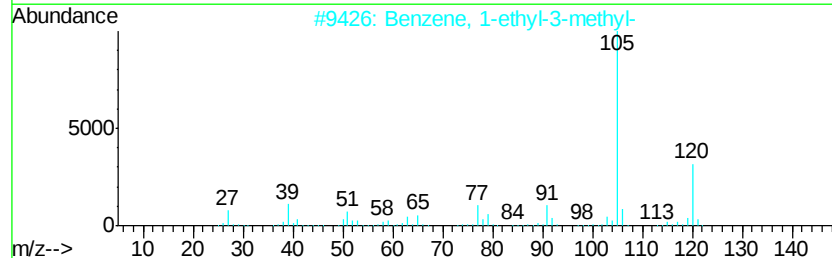
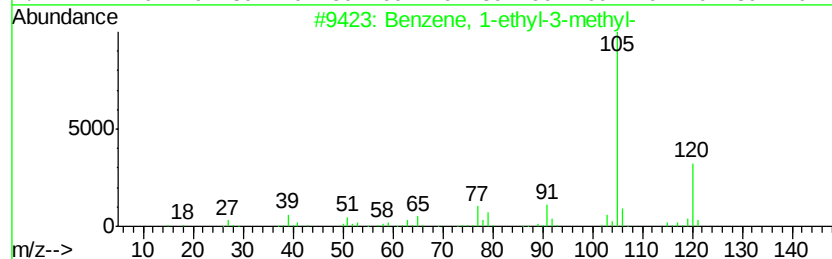
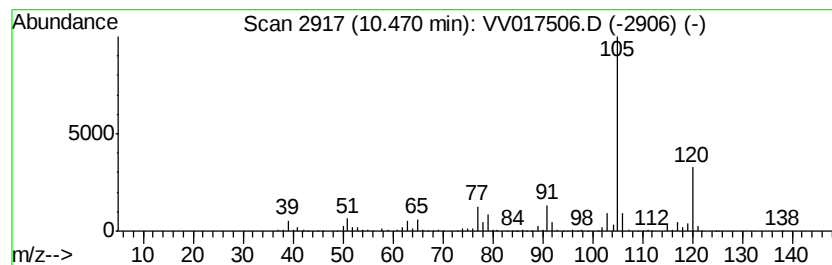
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Benzene, 1-ethyl-3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.47	3.61 ug/L	345784	1,4-Dichlorobenzene-d4	11.27

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-3-methyl-	120	C9H12	000620-14-4	94
3			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV071520\
Data File : VV017506.D
Acq On : 15 Jul 2020 12:55
Operator : SY/MD
Sample : L3281-20
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAXY0

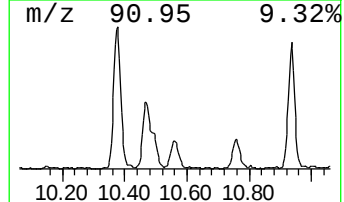
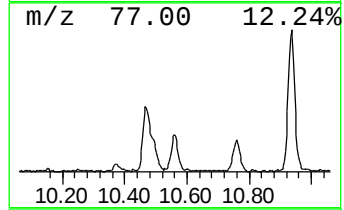
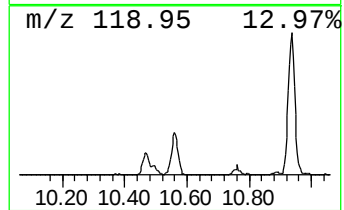
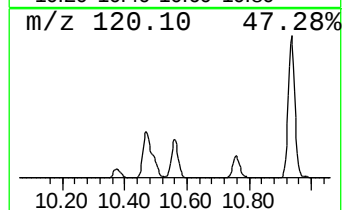
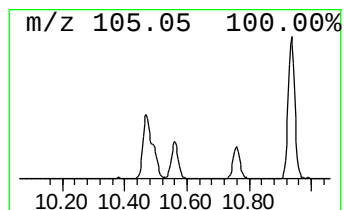
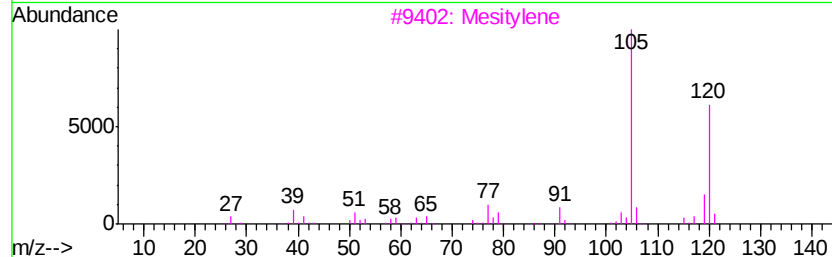
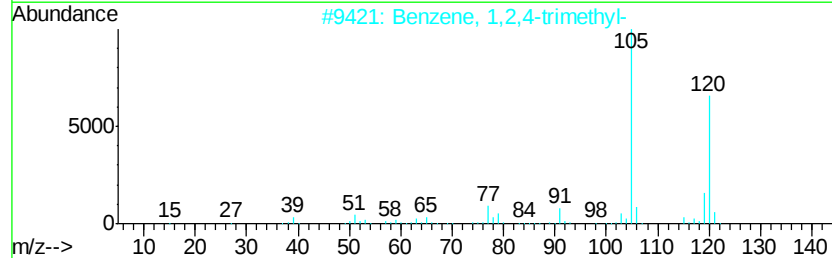
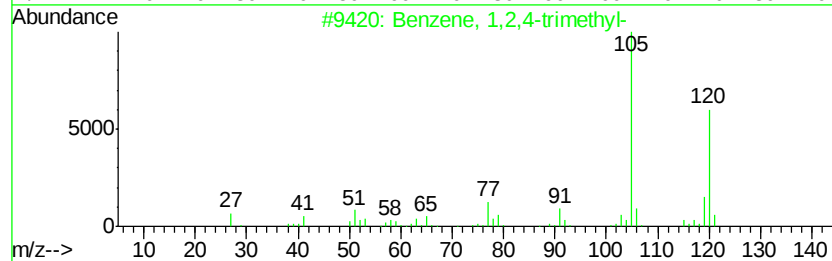
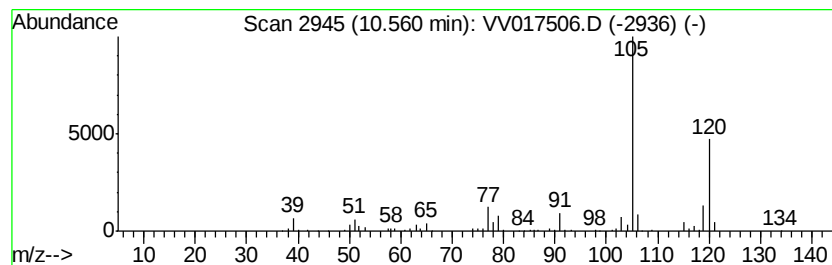
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Benzene, 1,2,4-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.56	1.67 ug/L	160360	1,4-Dichlorobenzene-d4	11.27

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
2			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
3			Mesitylene	120	C9H12	000108-67-8	95
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
5			Mesitylene	120	C9H12	000108-67-8	93



Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV071520\
Data File : VV017506.D
Acq On : 15 Jul 2020 12:55
Operator : SY/MD
Sample : L3281-20
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAXY0

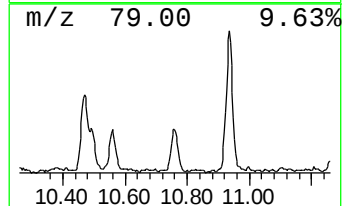
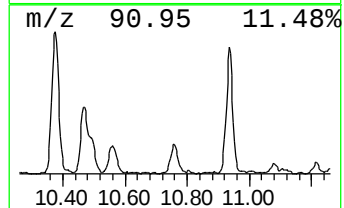
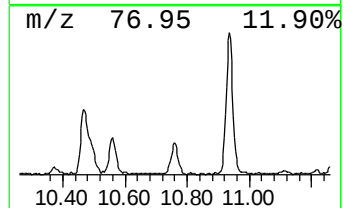
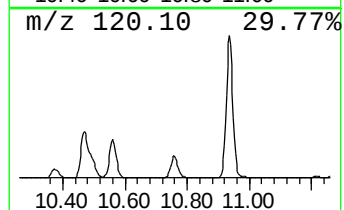
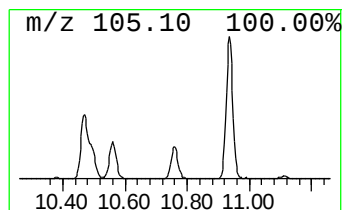
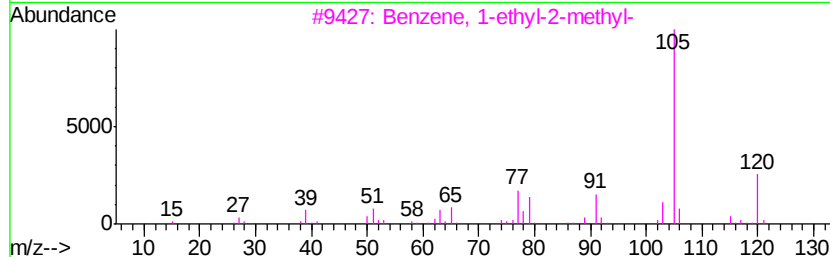
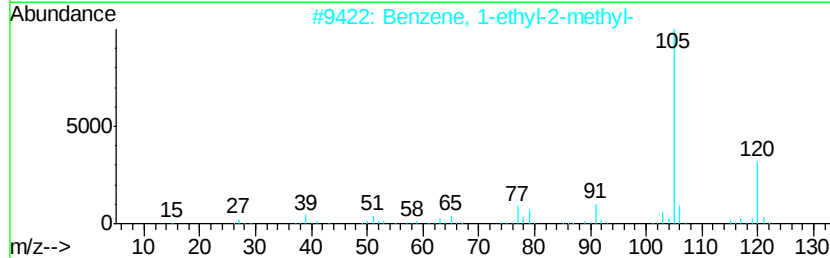
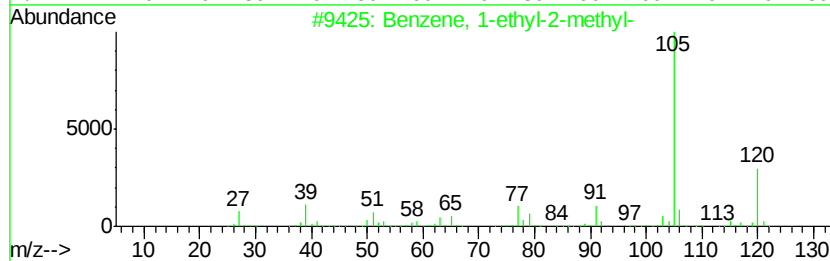
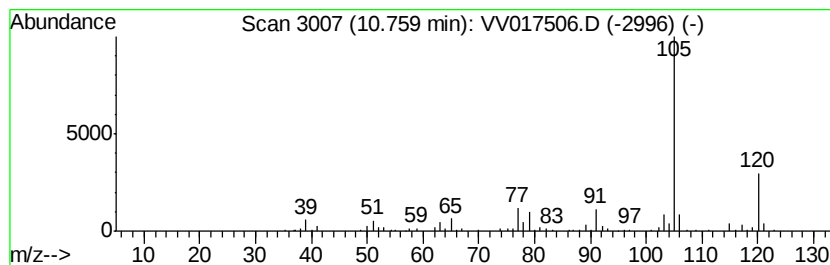
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Benzene, 1-ethyl-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.76	1.40 ug/L	133953	1,4-Dichlorobenzene-d4	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
3		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
5		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV071520\
Data File : VV017506.D
Acq On : 15 Jul 2020 12:55
Operator : SY/MD
Sample : L3281-20
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAXY0

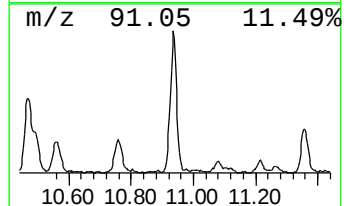
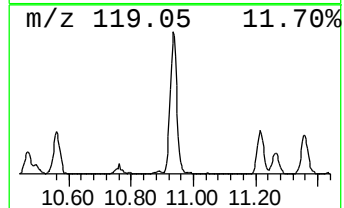
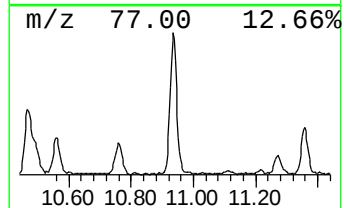
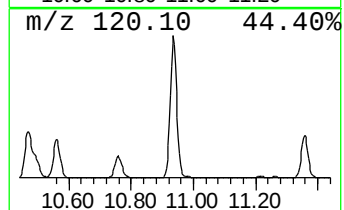
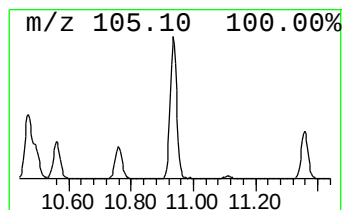
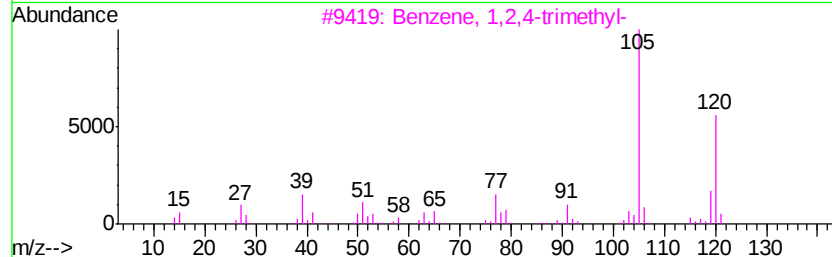
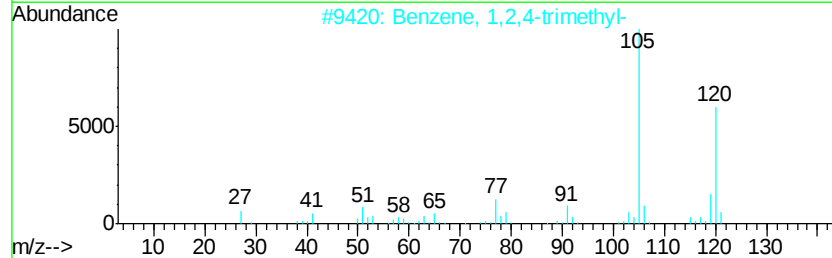
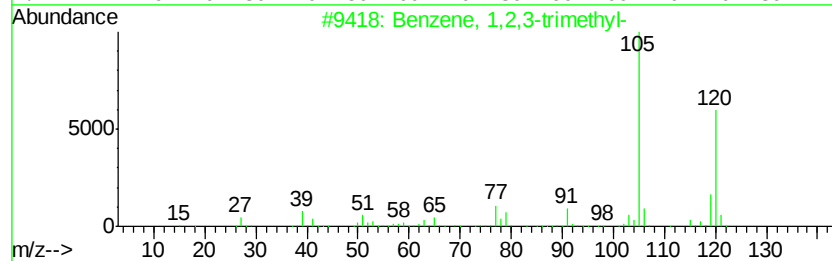
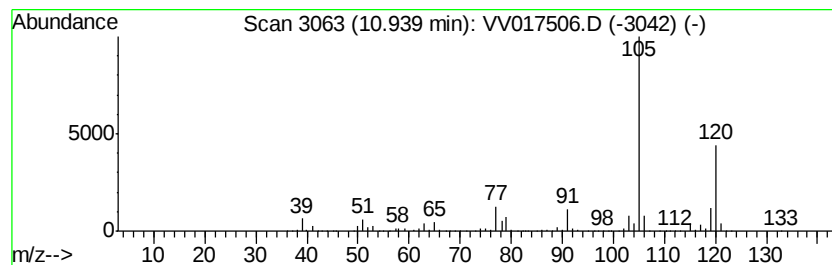
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Benzene, 1,2,3-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.94	5.85 ug/L	560425	1,4-Dichlorobenzene-d4	11.27

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV071520\
Data File : VV017506.D
Acq On : 15 Jul 2020 12:55
Operator : SY/MD
Sample : L3281-20
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAXY0

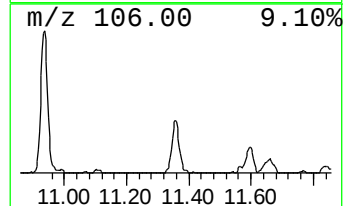
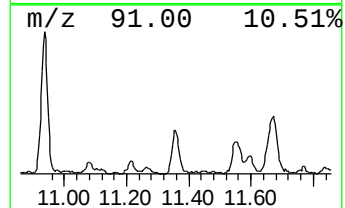
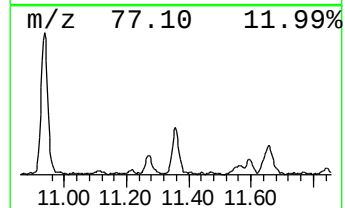
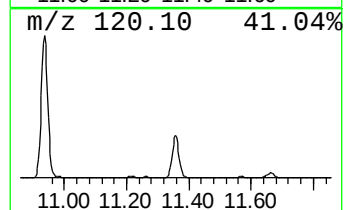
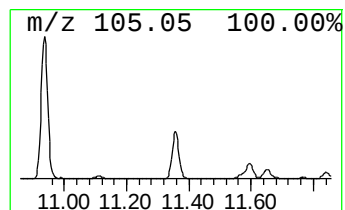
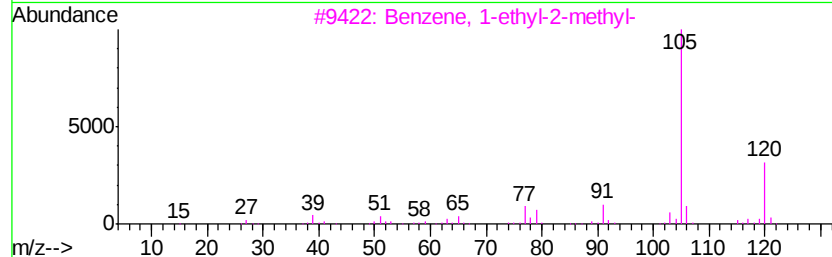
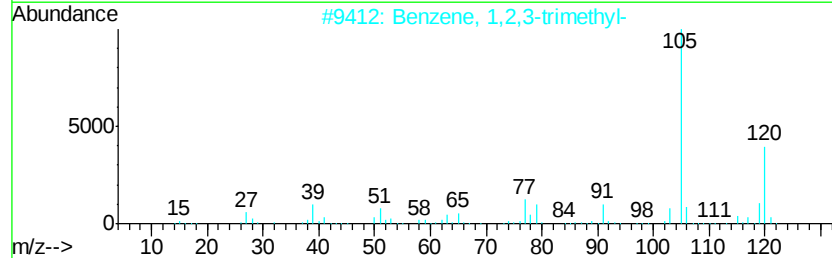
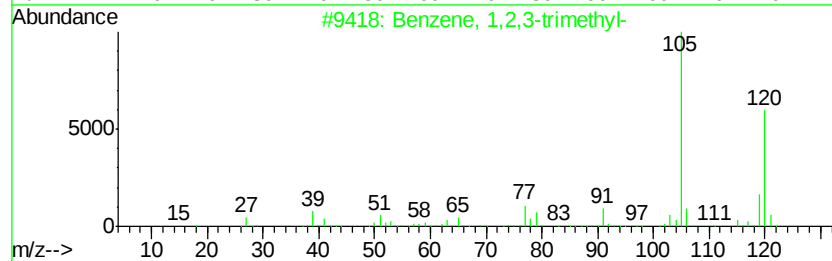
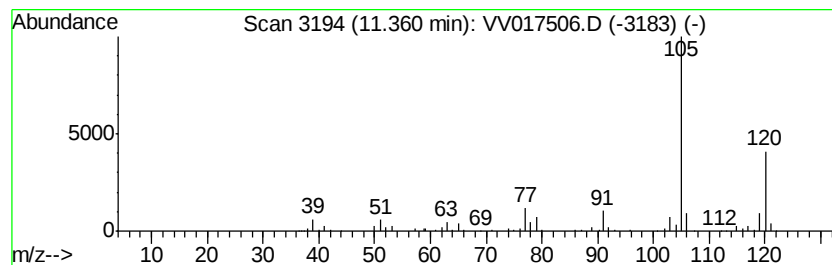
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Mesitylene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.36	1.89 ug/L	181043	1,4-Dichlorobenzene-d4	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
4		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
5		Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV071520\
Data File : VV017506.D
Acq On : 15 Jul 2020 12:55
Operator : SY/MD
Sample : L3281-20
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAXY0

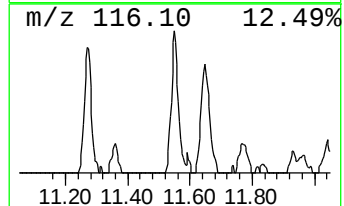
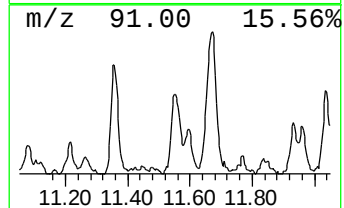
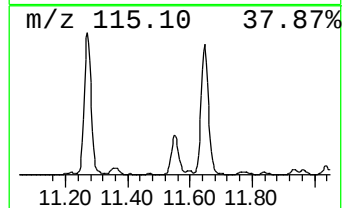
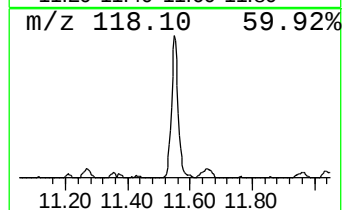
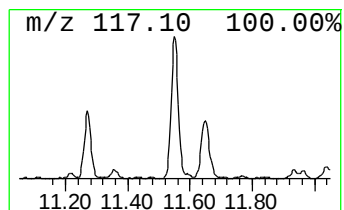
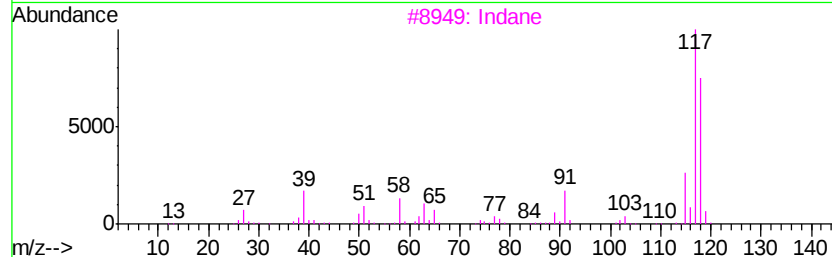
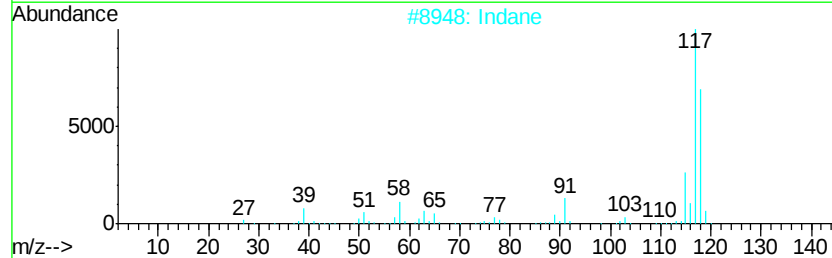
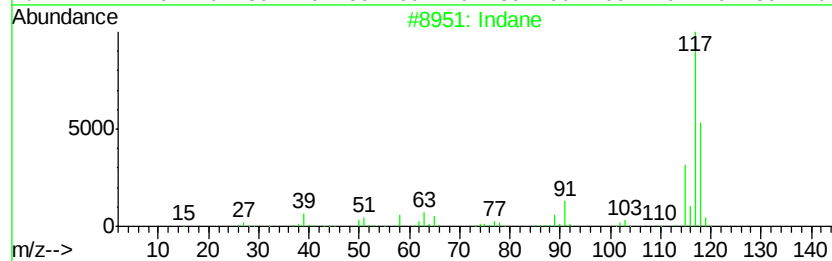
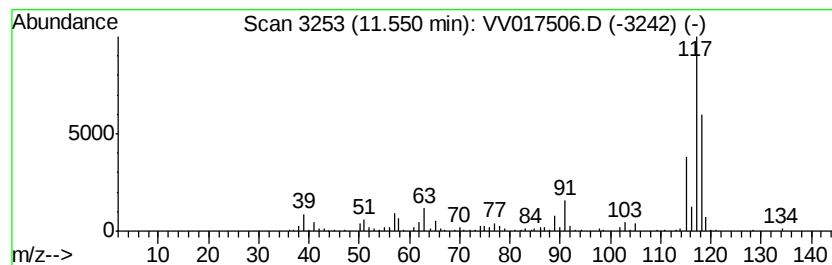
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 Indane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.55	1.57 ug/L	150569	1,4-Dichlorobenzene-d4	11.27

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	94
2	Indane		118	C9H10	000496-11-7	93
3	Indane		118	C9H10	000496-11-7	90
4	Indane		118	C9H10	000496-11-7	87
5	(Z)-1-Phenylpropene		118	C9H10	000766-90-5	70



Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV071520\
Data File : VV017506.D
Acq On : 15 Jul 2020 12:55
Operator : SY/MD
Sample : L3281-20
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAXY0

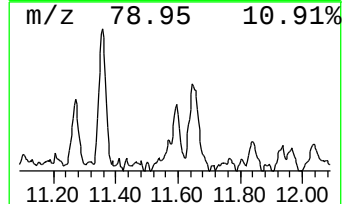
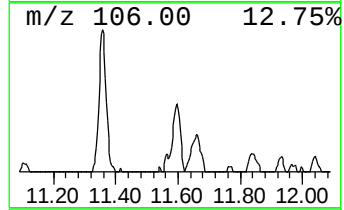
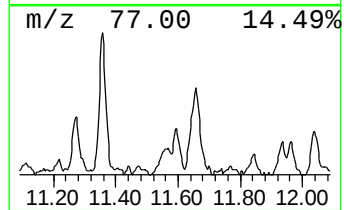
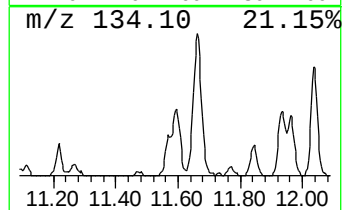
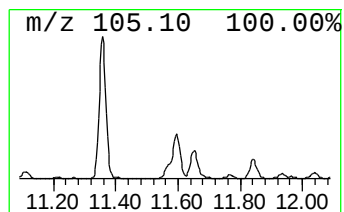
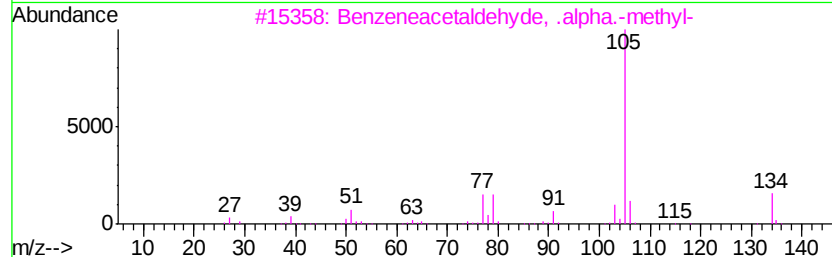
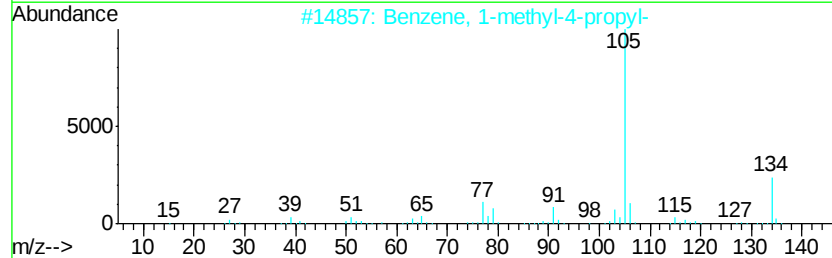
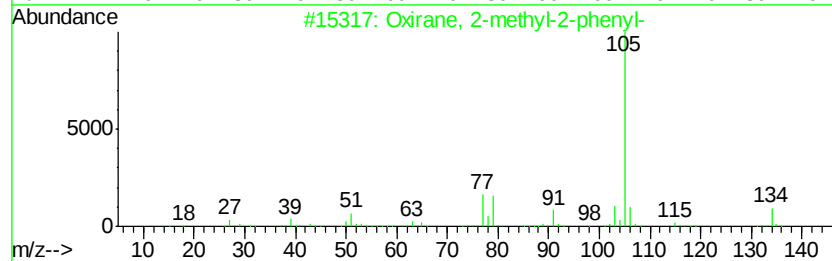
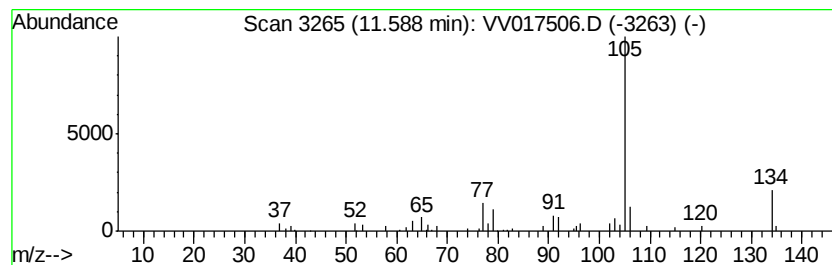
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 Oxirane, 2-methyl-2-phenyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.59	0.59 ug/L	56739	1,4-Dichlorobenzene-d4	11.27

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxirane, 2-methyl-2-phenyl-	134	C9H10O	002085-88-3	72
2			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	68
3			Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	64
4			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	64
5			Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	59



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV071520\
Data File : VV071506.D
Acq On : 15 Jul 2020 12:55
Operator : SY/MD
Sample : L3281-20
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAXY0

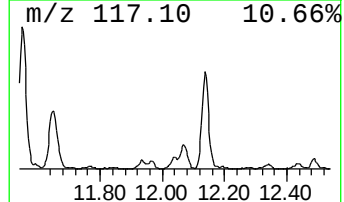
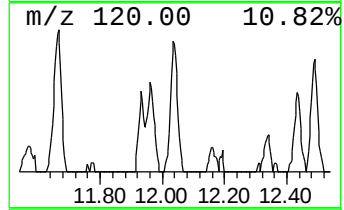
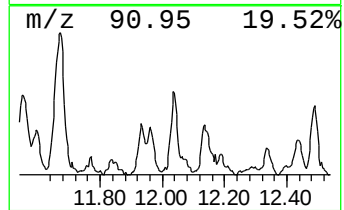
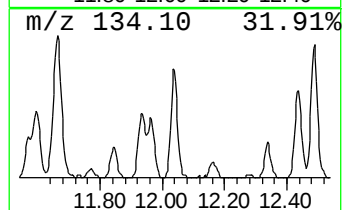
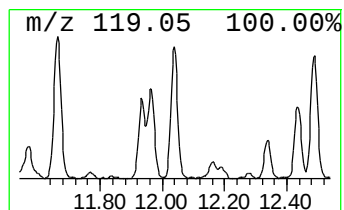
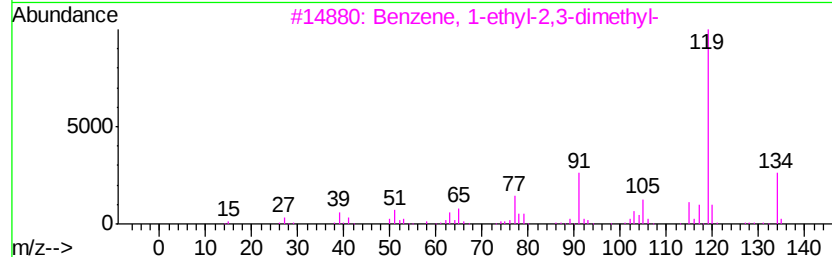
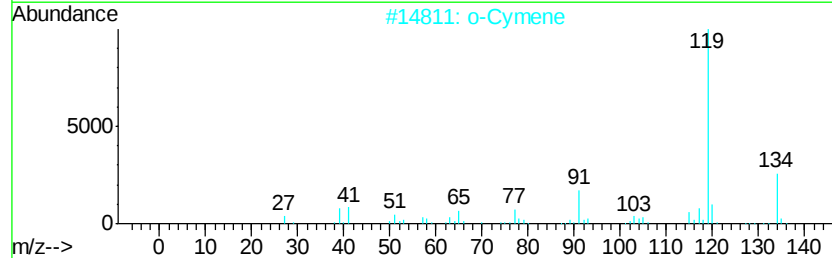
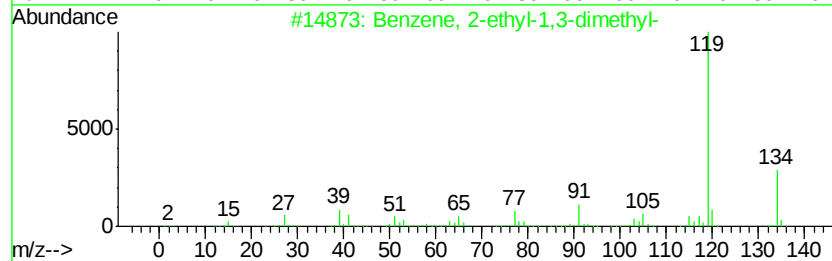
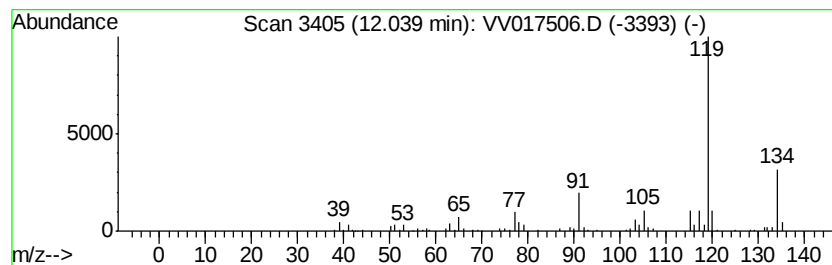
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 18 Benzene, 2-ethyl-1,3-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.04	0.83 ug/L	79535	1,4-Dichlorobenzene-d4	11.27

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
2			o-Cymene	134	C10H14	000527-84-4	95
3			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
5			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV071520\
Data File : VV017506.D
Acq On : 15 Jul 2020 12:55
Operator : SY/MD
Sample : L3281-20
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAXY0

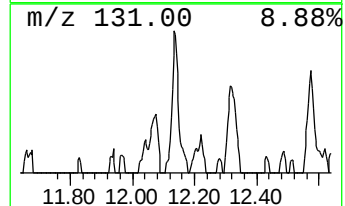
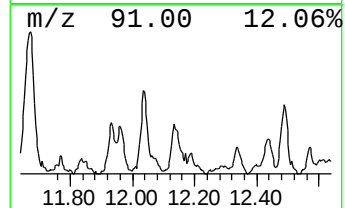
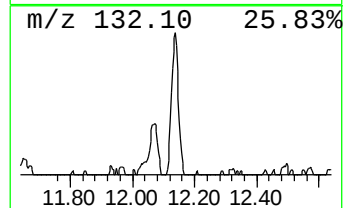
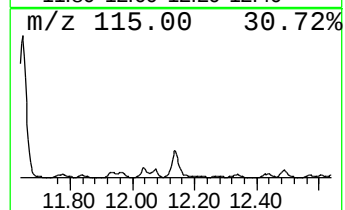
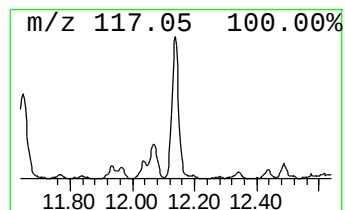
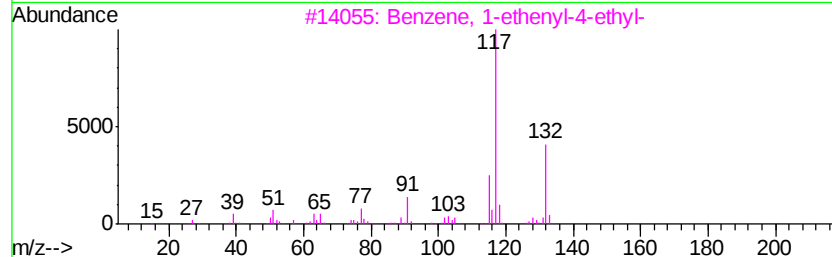
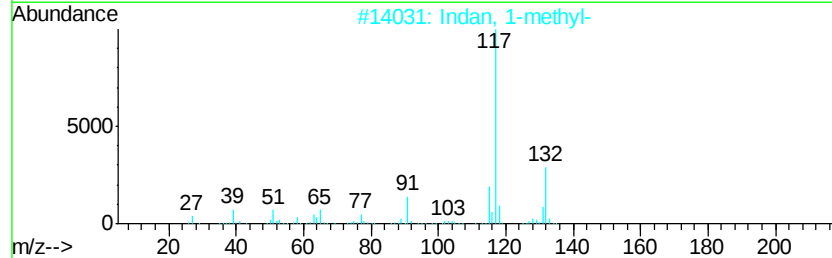
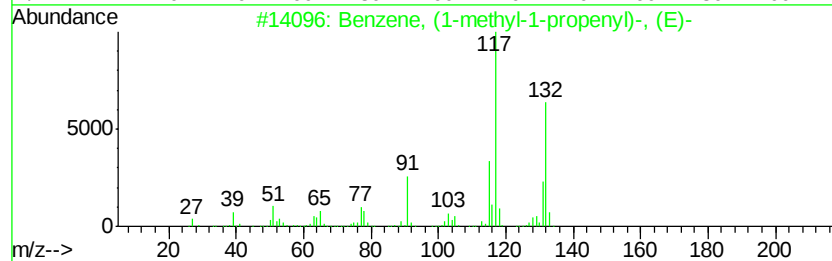
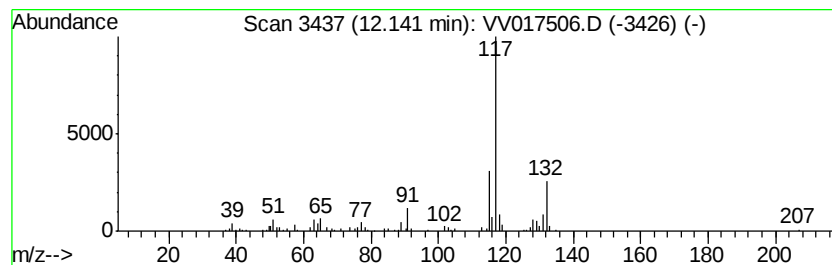
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 19 Benzene, (1-methyl-1-propen... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	0.89 ug/L	84909	1,4-Dichlorobenzene-d4	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1-methyl-1-propenyl)-, ...	132	C10H12	000768-00-3	90
2		Indan, 1-methyl-	132	C10H12	000767-58-8	87
3		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	86
4		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	86
5		Indan, 1-methyl-	132	C10H12	000767-58-8	81



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV071520\
Data File : VV071506.D
Acq On : 15 Jul 2020 12:55
Operator : SY/MD
Sample : L3281-20
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAXY0

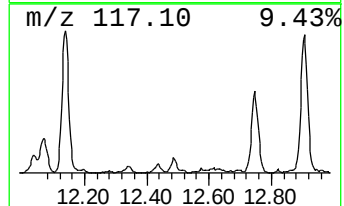
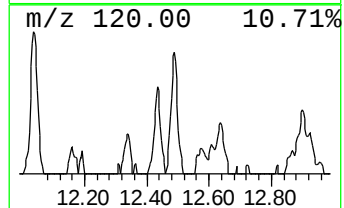
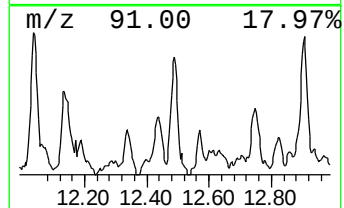
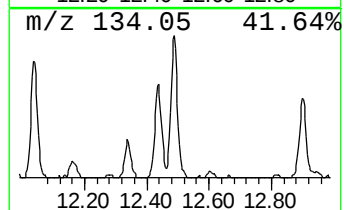
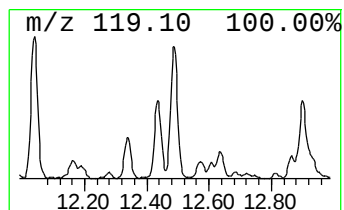
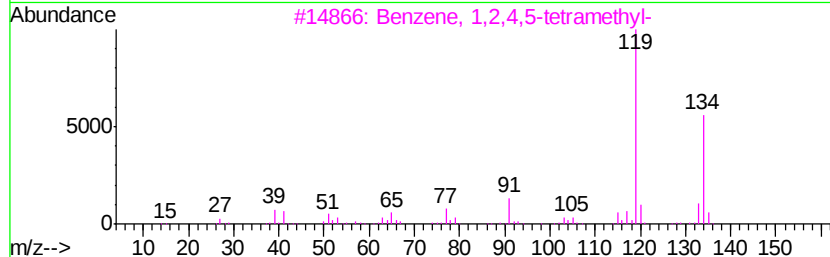
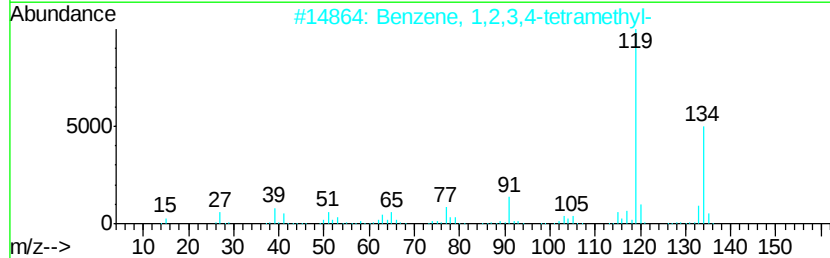
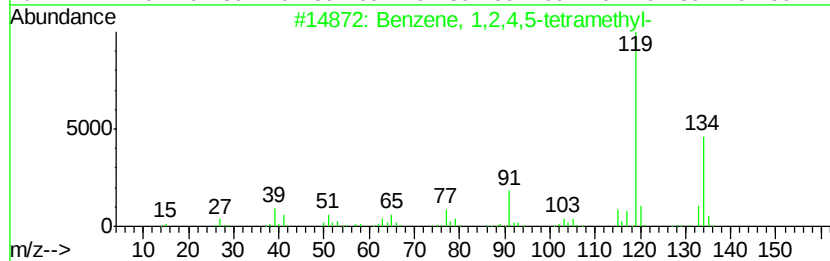
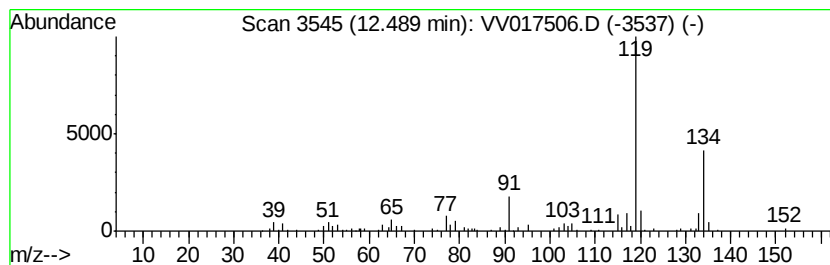
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 20 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.49	0.80 ug/L	76960	1,4-Dichlorobenzene-d4	11.27

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
3			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94
4			o-Cymene	134	C10H14	000527-84-4	94
5			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94



Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV071520\
Data File : VV071506.D
Acq On : 15 Jul 2020 12:55
Operator : SY/MD
Sample : L3281-20
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAXY0

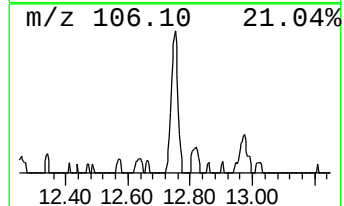
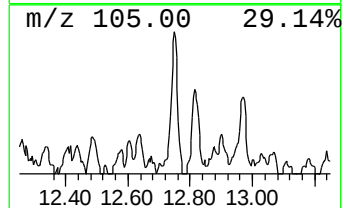
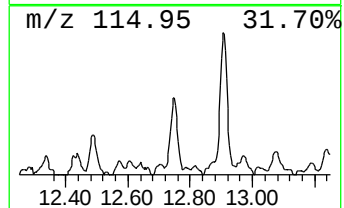
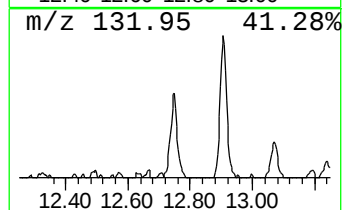
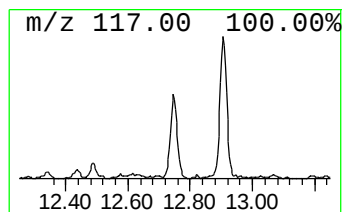
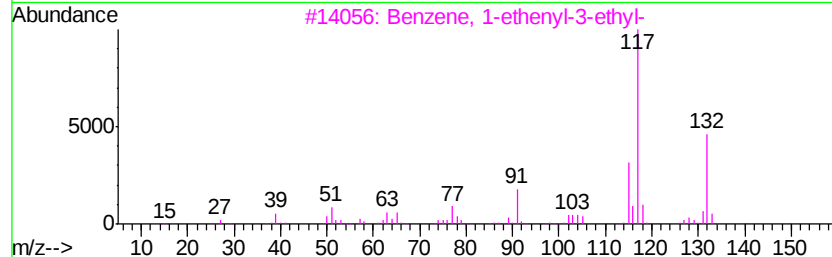
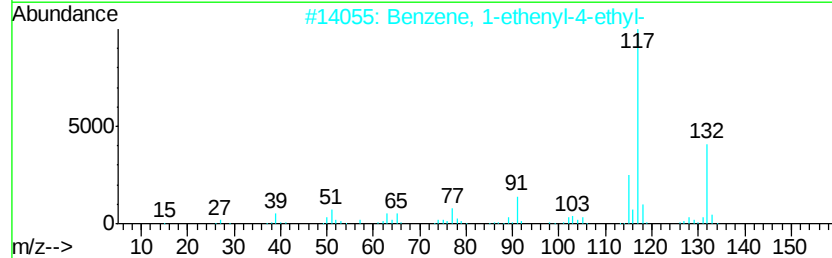
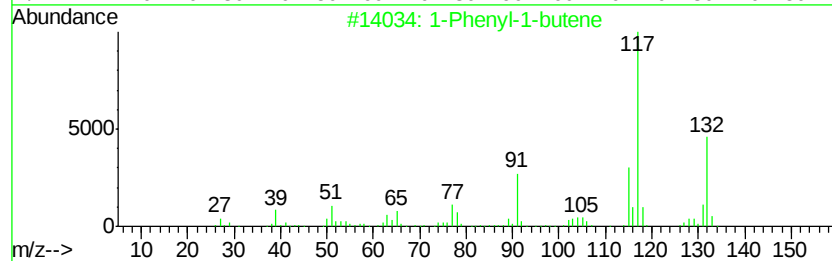
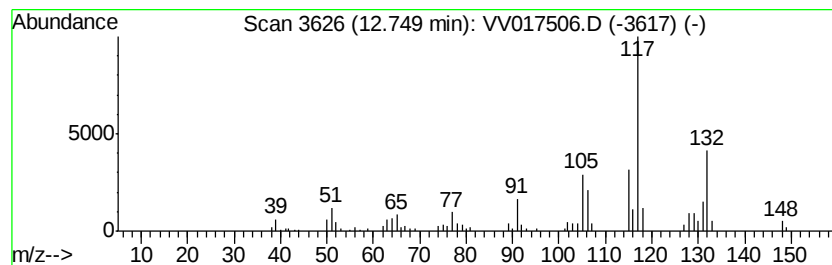
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 21 1-Phenyl-1-butene Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.75	0.64 ug/L	61144	1,4-Dichlorobenzene-d4	11.27

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenyl-1-butene	132	C10H12	000824-90-8	94
2			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	92
3			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	92
4			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
5			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA V\DATA\VV071520\
Data File : VV017506.D
Acq On : 15 Jul 2020 12:55
Operator : SY/MD
Sample : L3281-20
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GAXY0

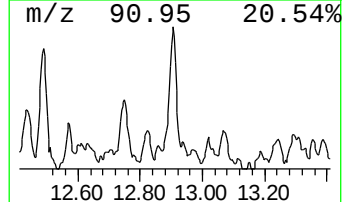
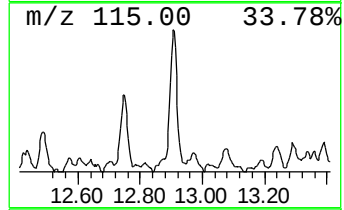
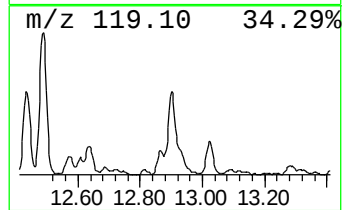
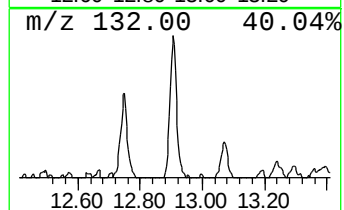
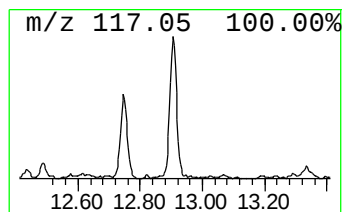
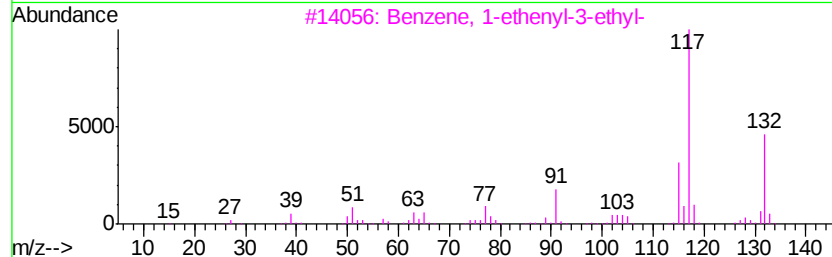
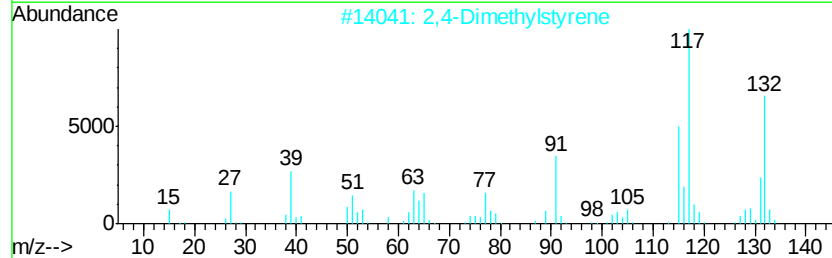
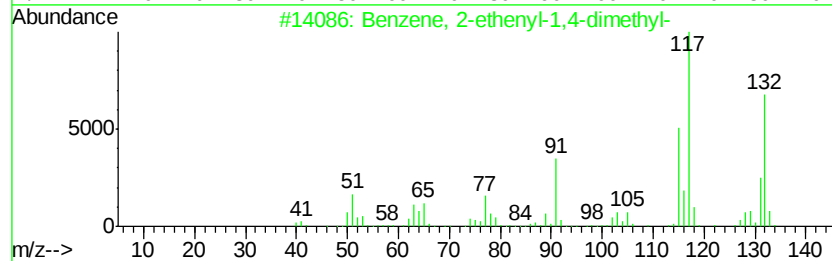
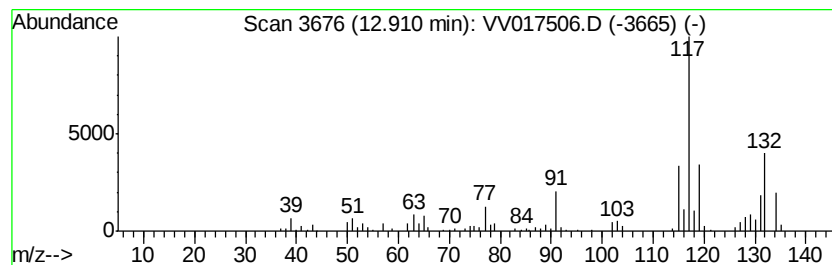
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 22 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.91	1.10 ug/L	105220	1,4-Dichlorobenzene-d4	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	87
2		2,4-Dimethylstyrene	132	C10H12	002234-20-0	70
3		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	64
4		1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	64
5		1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	64



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV071520\
 Data File : VV017506.D
 Acq On : 15 Jul 2020 12:55
 Operator : SY/MD
 Sample : L3281-20
 Misc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 GAXY0

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR071420WMA.M
 Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Butenal, (E)-	1.12	0.9	ug/L	68895	1	5.64	373572	5.0
(DEL) Alkane: Str...	1.65	0.9	ug/L	70780	1	5.64	373572	5.0
(DEL) Alkane: Cyc...	3.79	0.6	ug/L	42015	1	5.64	373572	5.0
(DEL) Alkane: Str...	4.94	0.5	ug/L	38600	1	5.64	373572	5.0
(DEL) Alkane: Str...	5.52	0.7	ug/L	49372	1	5.64	373572	5.0
(DEL) Alkane: Cyc...	7.29	0.5	ug/L	43172	2	8.87	428199	5.0
(DEL) Alkane: Str...	7.59	0.7	ug/L	60199	2	8.87	428199	5.0
(DEL) Alkane: Str...	9.23	1.1	ug/L	91308	2	8.87	428199	5.0
Benzene, propyl-	10.38	0.7	ug/L	66637	3	11.27	479125	5.0
Benzene, 1-ethyl-...	10.47	3.6	ug/L	345784	3	11.27	479125	5.0
Benzene, 1,2,4-tr...	10.56	1.7	ug/L	160360	3	11.27	479125	5.0
Benzene, 1-ethyl-...	10.76	1.4	ug/L	133953	3	11.27	479125	5.0
Benzene, 1,2,3-tr...	10.94	5.8	ug/L	560425	3	11.27	479125	5.0
Mesitylene	11.36	1.9	ug/L	181043	3	11.27	479125	5.0
Indane	11.55	1.6	ug/L	150569	3	11.27	479125	5.0
Oxirane, 2-methyl...	11.59	0.6	ug/L	56739	3	11.27	479125	5.0
Benzene, 2-ethyl-...	12.04	0.8	ug/L	79535	3	11.27	479125	5.0
Benzene, (1-methy...	12.14	0.9	ug/L	84909	3	11.27	479125	5.0
Benzene, 1,2,4,5-...	12.49	0.8	ug/L	76960	3	11.27	479125	5.0
1-Phenyl-1-butene	12.75	0.6	ug/L	61144	3	11.27	479125	5.0
Benzene, 2-etheny...	12.91	1.1	ug/L	105220	3	11.27	479125	5.0