

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV052820\
 Data File : VV016297.D
 Acq On : 29 May 2020 00:18
 Operator : SY/MD
 Sample : L2766-07DL 25X
 Misc : 25.0mL/MSVOA V/WATER
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 BE4Z0DL

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\SOMVTR052620WMA.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	1.315	62	70	85	rVB	36984	39770	9.58%	0.987%
2	1.441	91	109	110	rBV7	4008	8695	2.09%	0.216%
3	1.582	146	153	155	rBV	33255	35960	8.66%	0.892%
4	1.968	265	273	283	rBV	14093	19564	4.71%	0.485%
5	2.126	311	322	334	rBV	57465	84790	20.43%	2.104%
6	3.212	650	660	676	rVB2	5291	11676	2.81%	0.290%
7	3.318	676	693	713	rBV2	28297	66947	16.13%	1.661%
8	3.785	827	838	848	rBV7	2904	6124	1.48%	0.152%
9	3.942	871	887	931	rBV4	56257	194908	46.95%	4.837%
10	4.380	1006	1023	1044	rBV2	48184	120643	29.06%	2.994%
11	5.074	1223	1239	1248	rBV2	113043	271648	65.44%	6.741%
12	5.125	1249	1255	1276	rVB	89261	181900	43.82%	4.514%
13	5.643	1404	1416	1445	rBV	109696	216922	52.26%	5.383%
14	6.096	1545	1557	1570	rBV	65431	131644	31.71%	3.267%
15	7.010	1830	1841	1858	rBV	28404	50917	12.27%	1.264%
16	7.338	1932	1943	1955	rBV	131324	223763	53.90%	5.553%
17	7.412	1955	1966	1985	rVB	240848	415116	100.00%	10.301%
18	7.646	2030	2039	2054	rBV2	17050	29648	7.14%	0.736%
19	8.000	2141	2149	2156	rVB3	3522	4972	1.20%	0.123%
20	8.116	2174	2185	2224	rBV	146811	276683	66.65%	6.866%
21	8.874	2409	2421	2426	rBV	178199	294939	71.05%	7.319%
22	9.035	2455	2471	2482	rBV	49010	77429	18.65%	1.921%
23	9.083	2482	2486	2497	rVB3	4998	7558	1.82%	0.188%
24	9.157	2497	2509	2528	rBV	231092	365234	87.98%	9.064%
25	9.566	2623	2636	2652	rBV	45843	74431	17.93%	1.847%
26	9.955	2744	2757	2772	rBV2	17062	26999	6.50%	0.670%
27	10.238	2834	2845	2855	rBV2	46941	71922	17.33%	1.785%
28	10.292	2856	2862	2872	rVB5	5697	7840	1.89%	0.195%
29	10.376	2879	2888	2903	rBV2	16447	26562	6.40%	0.659%
30	10.563	2936	2946	2957	rVB4	3662	5215	1.26%	0.129%
31	10.759	2998	3007	3018	rBV2	14238	21047	5.07%	0.522%
32	10.939	3053	3063	3076	rBV	18466	28082	6.76%	0.697%
33	11.270	3156	3166	3185	rBV	175691	283308	68.25%	7.030%
34	11.363	3185	3195	3208	rVB2	10007	19685	4.74%	0.488%

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Title : TRACE VOA SOM01.0

35	11.553	3245	3254	3272	rVB3	16884	38434	9.26%	0.954%
36	11.646	3272	3283	3308	rVB	139452	230675	55.57%	5.724%
37	11.935	3362	3373	3378	rBV4	3456	4790	1.15%	0.119%
38	12.038	3394	3405	3414	rBV	14735	21330	5.14%	0.529%
39	12.437	3517	3529	3537	rBV	8940	12643	3.05%	0.314%
40	12.492	3537	3546	3556	rVB3	6516	9811	2.36%	0.243%
41	12.900	3668	3673	3682	rVB5	3718	5254	1.27%	0.130%
42	13.533	3860	3870	3880	rBV2	2631	4237	1.02%	0.105%

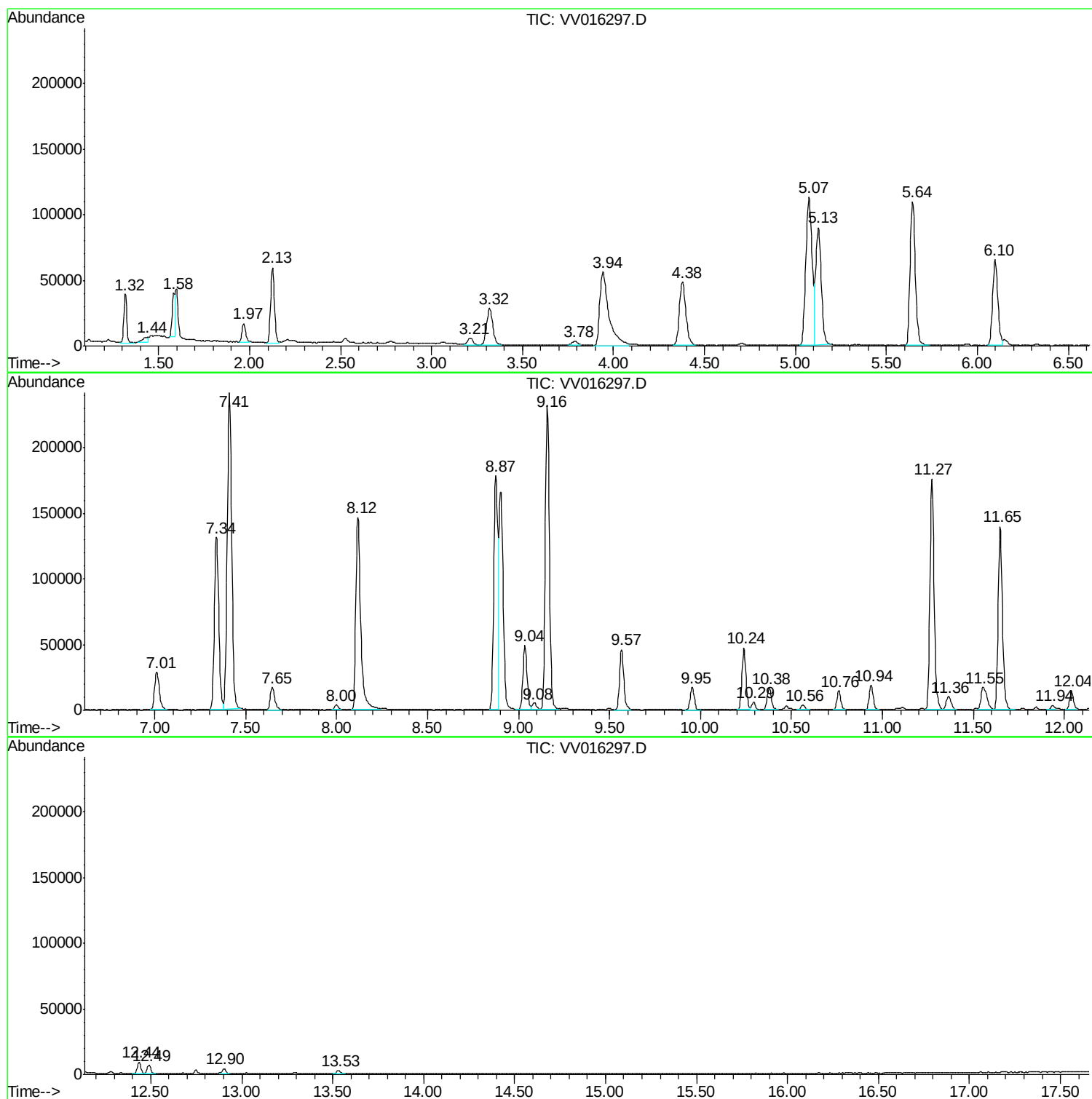
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR052620WMA.M
Quant Title : TRACE VOA SOM01.0

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



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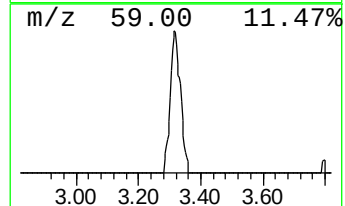
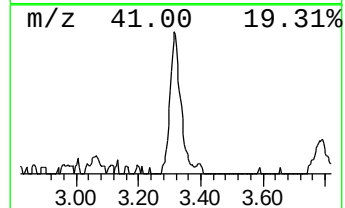
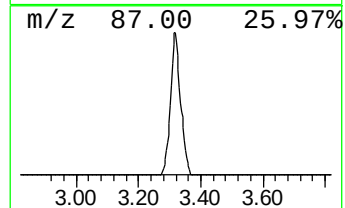
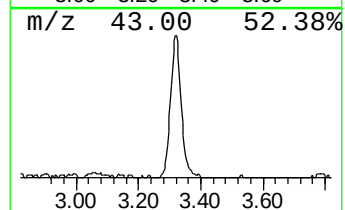
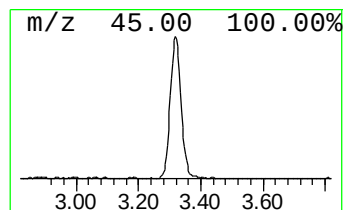
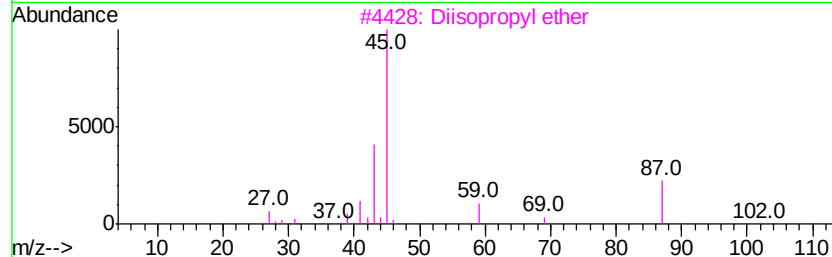
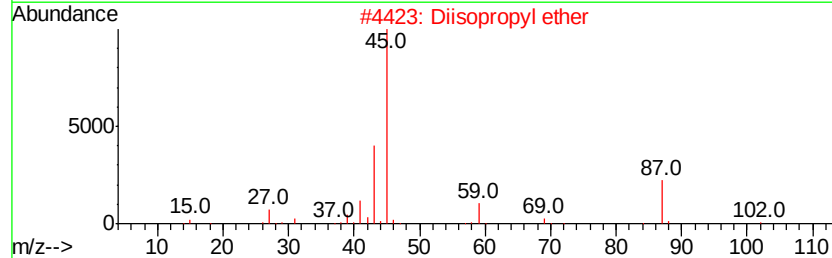
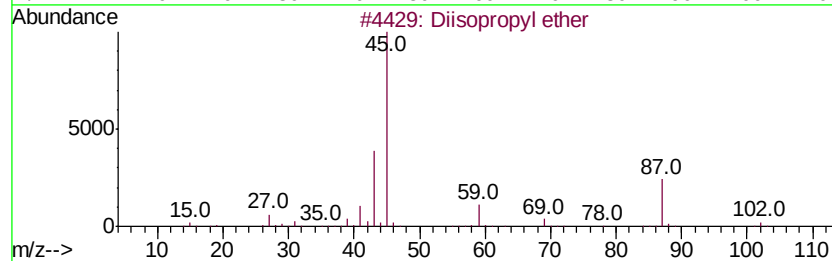
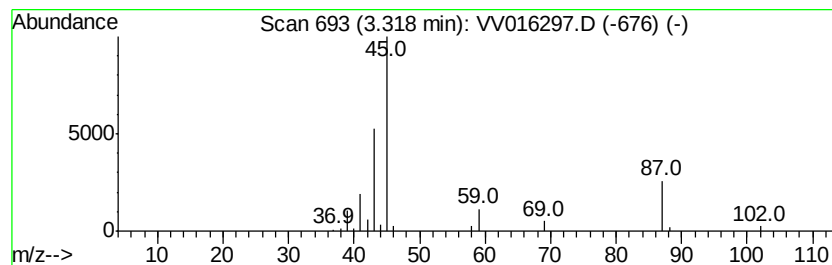
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR052620WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 1 Diisopropyl ether Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.32	1.54 ug/L	66947	1,4-Difluorobenzene	5.64

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Diisopropyl ether	102	C6H14O	000108-20-3	91
2			Diisopropyl ether	102	C6H14O	000108-20-3	83
3			Diisopropyl ether	102	C6H14O	000108-20-3	74
4			Ether, 2-chloro-1-methylethyl is...	136	C6H13ClO	098277-76-0	64
5			Propane, 2-ethoxy-	88	C5H12O	000625-54-7	59



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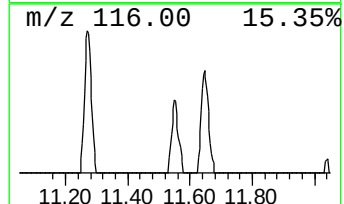
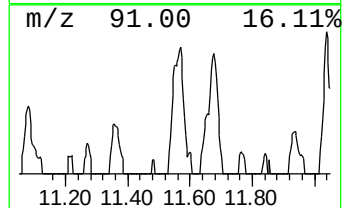
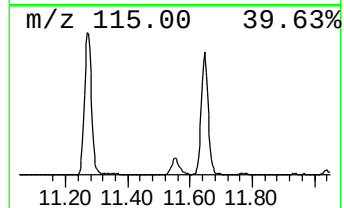
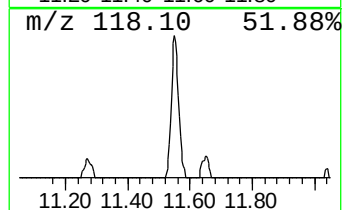
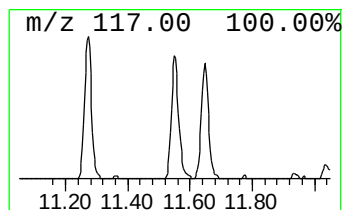
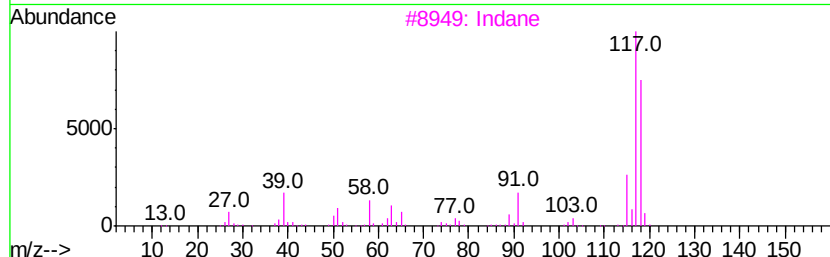
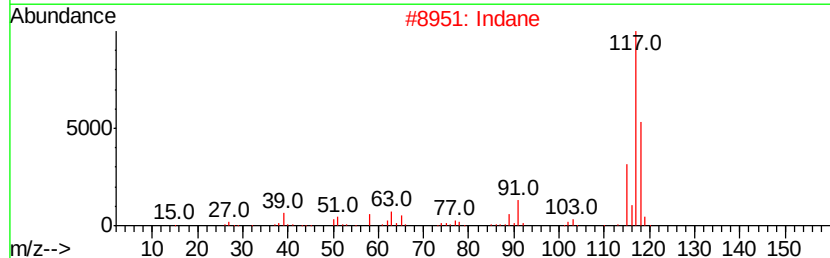
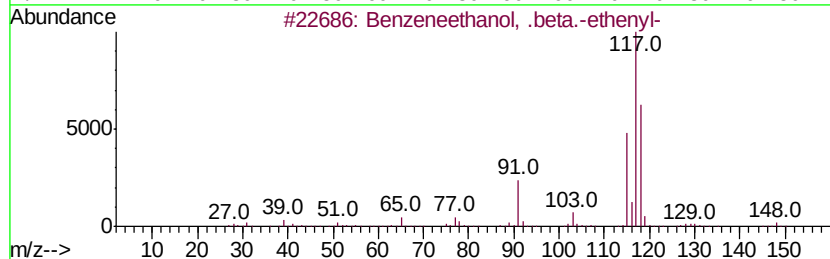
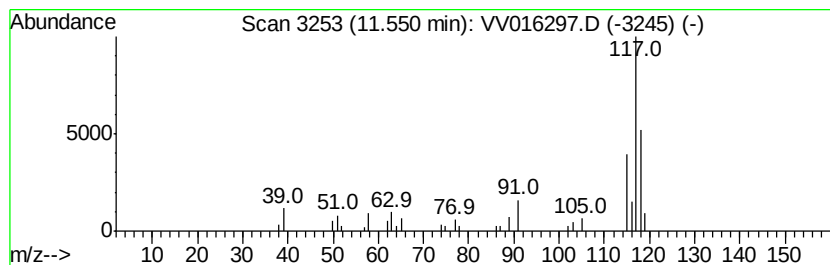
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Peak Number 3 Benzeneethanol, .beta.-ethe... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.55	0.68 ug/L	38434	1,4-Dichlorobenzene-d4	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzeneethanol, .beta.-ethenyl-	148	C10H12O	006052-63-7	78
2		Indane	118	C9H10	000496-11-7	70
3		Indane	118	C9H10	000496-11-7	70
4		Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
5		Indane	118	C9H10	000496-11-7	64



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Diisopropyl ether	3.32	1.5	ug/L	66947	1	5.64	216922	5.0
Benzeneethanol,	11.55	0.7	ug/L	38434	3	11.27	283308	5.0