

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN081720\
 Data File : VN062990.D
 Acq On : 17 Aug 2020 23:57
 Operator : JC/MD
 Sample : L3696-06
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 W-1

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

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_N\METHODS\82N081720W.M

Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.429	197	204	215	rVB7	11882	20059	2.11%	0.322%
2	7.548	1778	1796	1808	rBV	126174	317898	33.37%	5.098%
3	7.625	1808	1820	1848	rVB	210876	501489	52.64%	8.042%
4	7.989	1914	1933	1953	rBV	126113	293669	30.82%	4.709%
5	8.551	2091	2108	2133	rBV	315880	664336	69.73%	10.654%
6	10.059	2563	2577	2603	rBV	522010	952700	100.00%	15.278%
7	10.828	2806	2816	2832	rBV5	7148	15174	1.59%	0.243%
8	11.381	2972	2988	3011	rBV	501716	872458	91.58%	13.991%
9	12.226	3242	3251	3262	rBV2	9432	13960	1.47%	0.224%
10	12.378	3284	3298	3318	rBV	387943	655204	68.77%	10.507%
11	12.567	3350	3357	3366	rVB2	12654	18128	1.90%	0.291%
12	13.014	3484	3496	3508	rBV	24077	38883	4.08%	0.624%
13	13.146	3525	3537	3545	rBV2	10042	17410	1.83%	0.279%
14	13.320	3579	3591	3612	rVB	447551	798808	83.85%	12.810%
15	13.487	3635	3643	3646	rBV4	7690	10036	1.05%	0.161%
16	13.519	3646	3653	3661	rVV2	19196	31957	3.35%	0.512%
17	13.570	3661	3669	3682	rVV5	7333	17550	1.84%	0.281%
18	13.644	3682	3692	3702	rVB6	7672	12986	1.36%	0.208%
19	13.776	3723	3733	3738	rBV	9339	14228	1.49%	0.228%
20	13.805	3739	3742	3751	rVB2	8387	10605	1.11%	0.170%
21	13.866	3751	3761	3771	rVB2	21264	32110	3.37%	0.515%
22	13.972	3783	3794	3807	rVB4	23787	42927	4.51%	0.688%
23	14.107	3826	3836	3845	rBV4	9338	16800	1.76%	0.269%
24	14.185	3853	3860	3866	rBV2	10423	14764	1.55%	0.237%
25	14.223	3866	3872	3879	rVB3	16190	23003	2.41%	0.369%
26	14.451	3934	3943	3951	rBV2	23838	39050	4.10%	0.626%
27	14.500	3952	3958	3966	rVB2	13985	19494	2.05%	0.313%
28	14.567	3966	3979	3990	rBV4	61691	128082	13.44%	2.054%
29	14.625	3990	3997	4007	rVB5	11165	19014	2.00%	0.305%
30	14.718	4017	4026	4032	rVB6	6888	10447	1.10%	0.168%
31	14.824	4040	4059	4065	rBV6	28064	65039	6.83%	1.043%
32	14.931	4083	4092	4104	rVB4	25537	50895	5.34%	0.816%
33	15.162	4155	4164	4172	rBV2	20269	33953	3.56%	0.544%
34	15.217	4172	4181	4191	rBV7	13466	35947	3.77%	0.576%

LSC Area Percent Report

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Integration Parameters: RTEINT.P

Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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Title : SW846 8260

35	15.387	4222	4234	4248	rBV3	19680	37866	3.97%	0.607%
36	15.545	4264	4283	4286	rBV7	17514	38609	4.05%	0.619%
37	15.580	4287	4294	4311	rVV3	40178	75037	7.88%	1.203%
38	15.670	4314	4322	4329	rVV8	8505	13551	1.42%	0.217%
39	15.738	4332	4343	4357	rVV6	12815	30382	3.19%	0.487%
40	15.876	4371	4386	4398	rBV3	27350	58873	6.18%	0.944%
41	16.066	4435	4445	4454	rBV4	19722	38230	4.01%	0.613%
42	16.127	4454	4464	4476	rBV	27030	49820	5.23%	0.799%
43	16.316	4508	4523	4536	rBV3	25923	66398	6.97%	1.065%
44	16.458	4557	4567	4580	rVB8	7793	17925	1.88%	0.287%

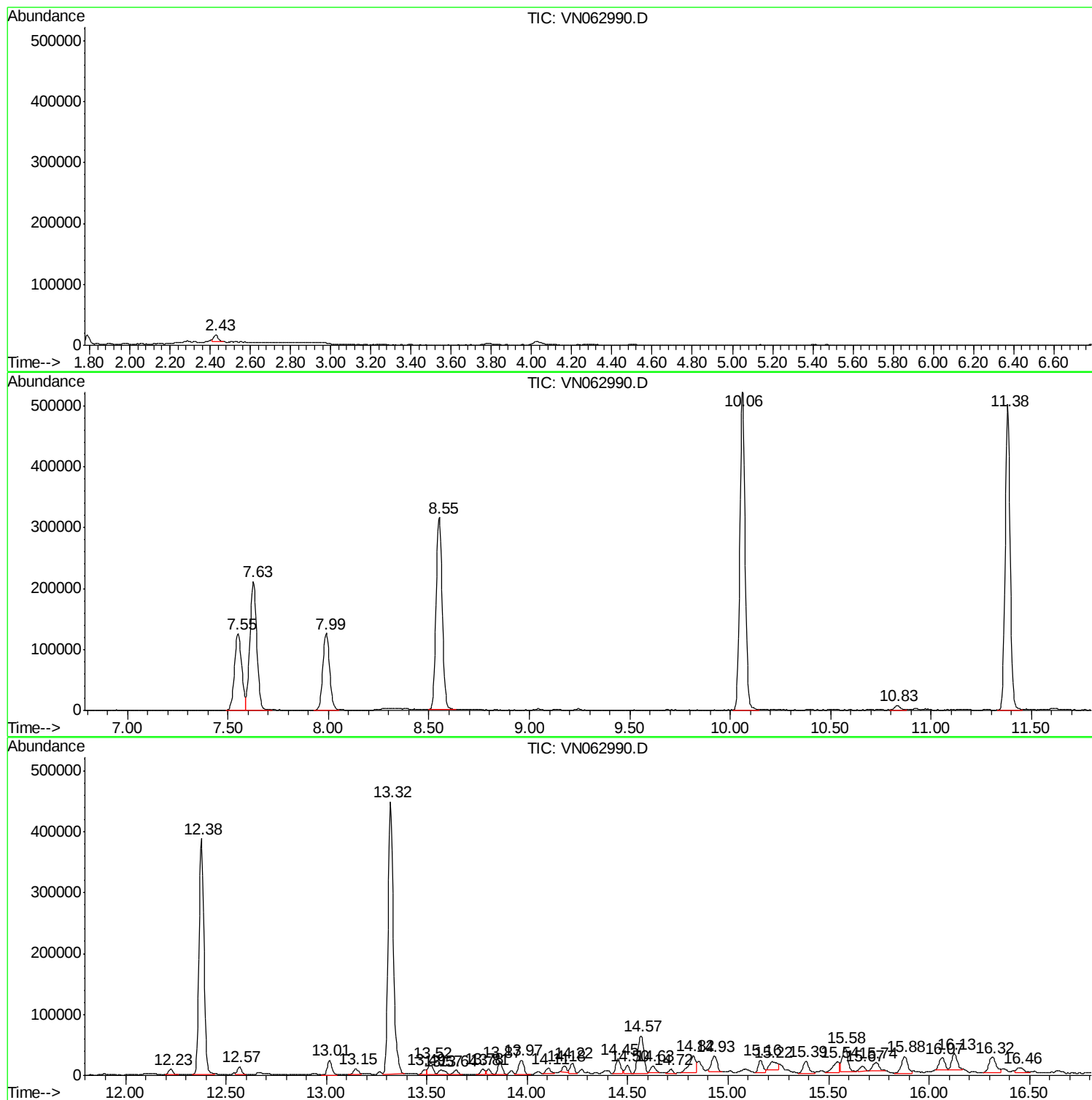
Sum of corrected areas: 6235754

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N081720W.M
Quant Title : SW846 8260

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



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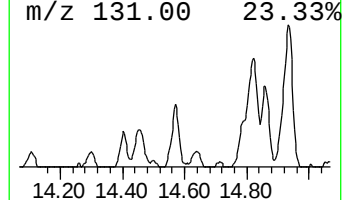
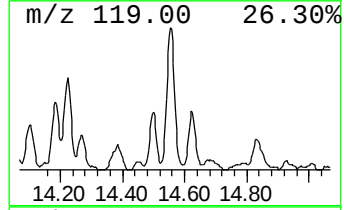
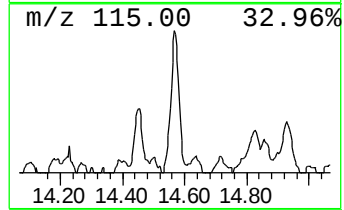
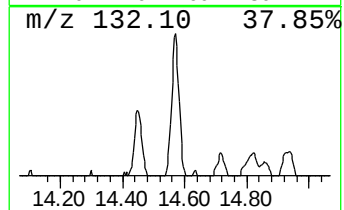
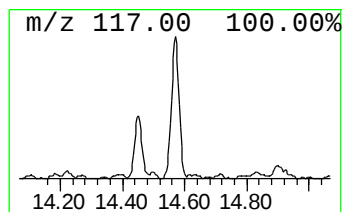
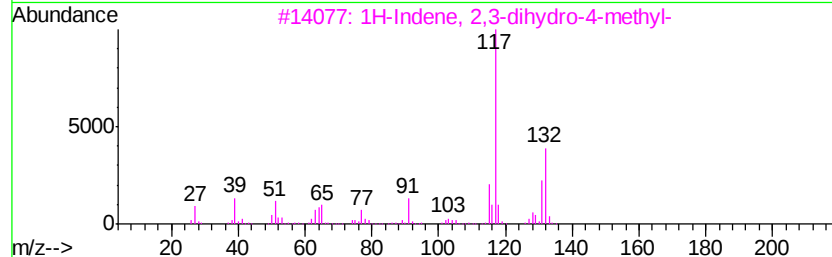
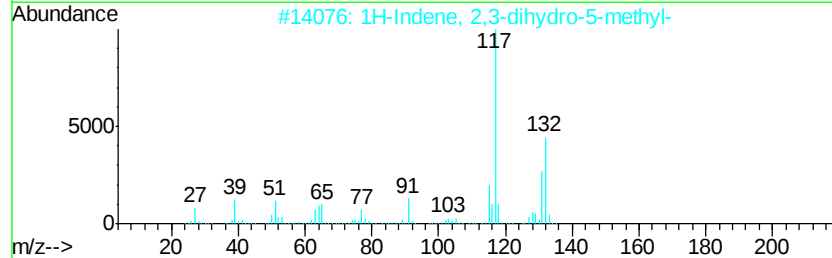
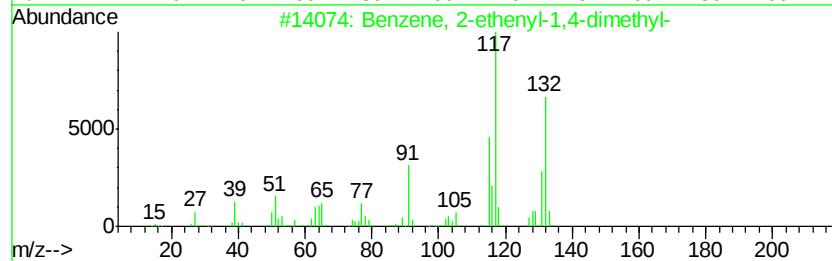
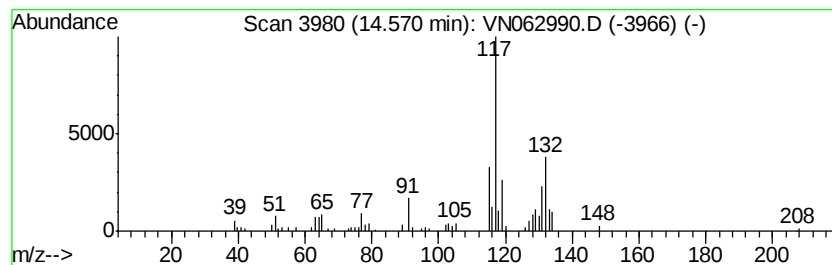
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.		
14.57	8.02 ug/l	128082	1,4-Dichlorobenzene-d4	13.32		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	90
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	70
3		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	70
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
5		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	64



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc

Benzene, 2-etheny...	14.57	8.0	ug/l	128082	4	13.32	798808	50.0