

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX060120\
 Data File : VX016517.D
 Acq On : 01 Jun 2020 10:27
 Operator : JC/SP
 Sample : VX0601WBL01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 VX0601WBL01

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_X\METHOD\82X052720W.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	5.464	706	718	730	rBV	174935	503230	29.32%	5.181%
2	5.623	733	744	761	rVB	338968	947271	55.19%	9.753%
3	6.031	799	811	826	rBV	191986	518661	30.22%	5.340%
4	6.824	931	941	952	rBV	500324	1202883	70.08%	12.385%
5	8.696	1240	1248	1255	rBV	1116480	1716331	100.00%	17.672%
6	10.098	1467	1478	1496	rBV	1080231	1535080	89.44%	15.806%
7	11.122	1640	1646	1655	rBV	933764	1211204	70.57%	12.471%
8	12.061	1795	1800	1807	rBV	1172125	1440540	83.93%	14.832%
9	12.951	1903	1946	1947	rBV2	110833	636958	37.11%	6.558%

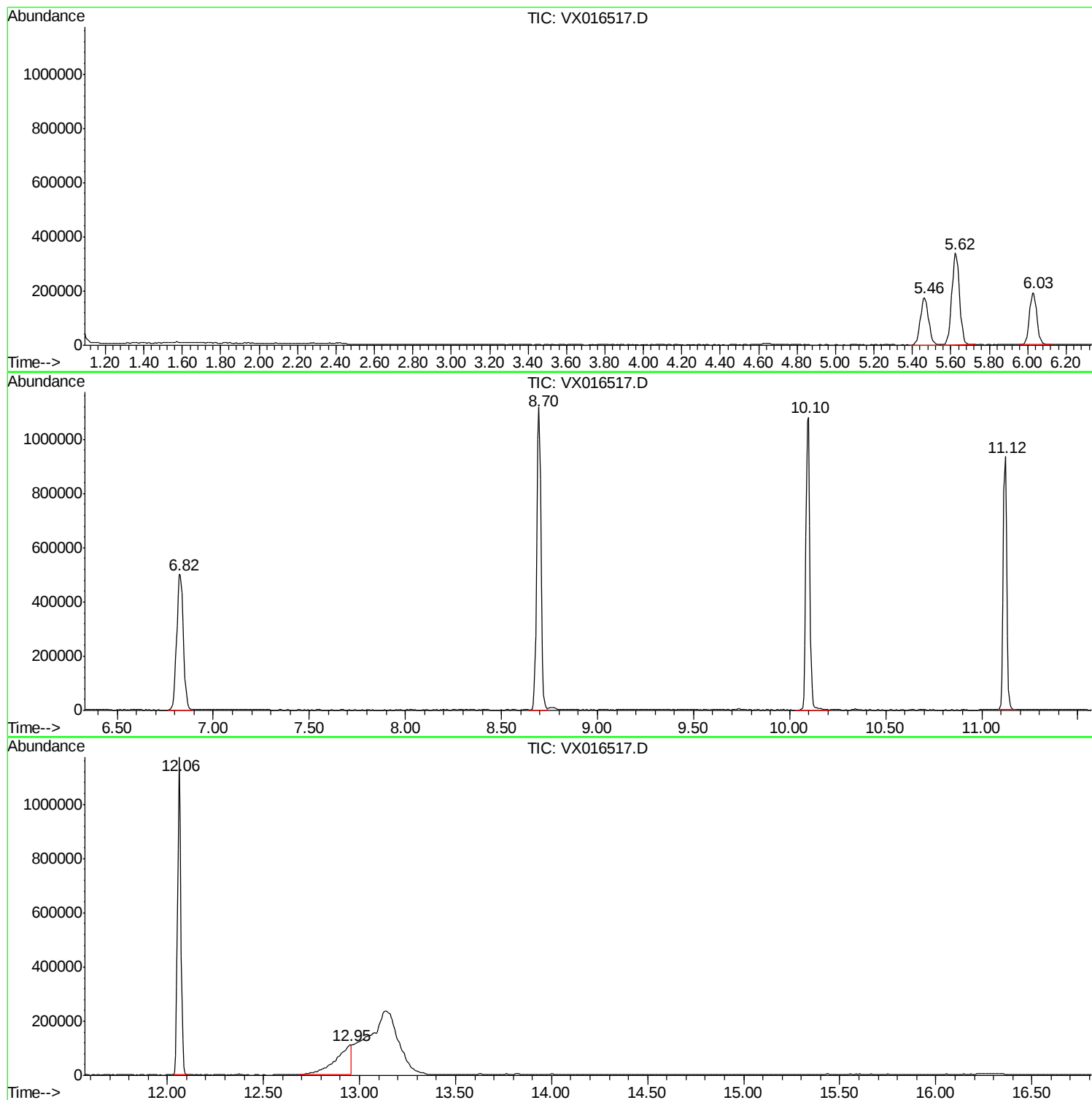
Sum of corrected areas: 9712158

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X052720W.M
Quant Title : SW846 8260

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



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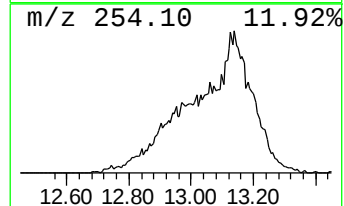
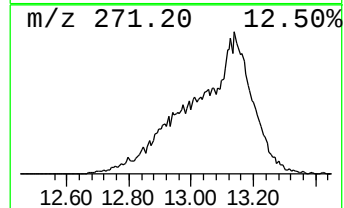
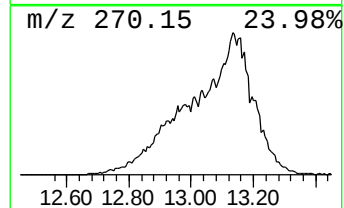
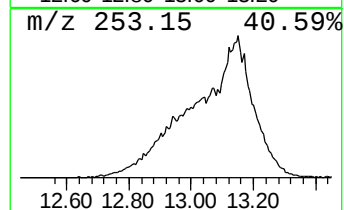
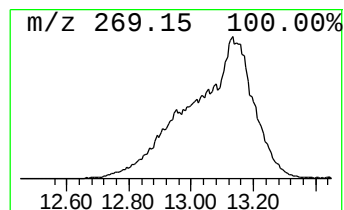
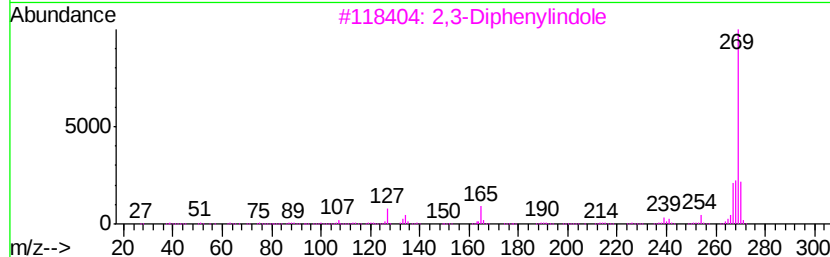
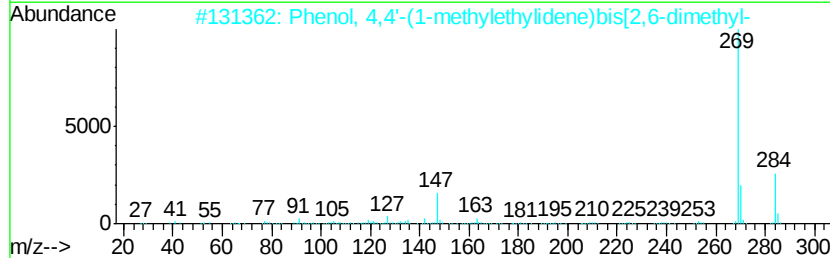
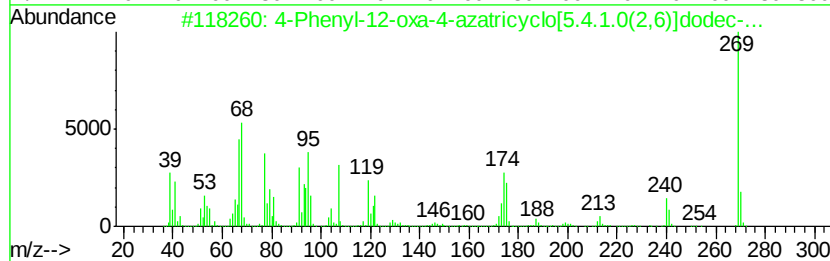
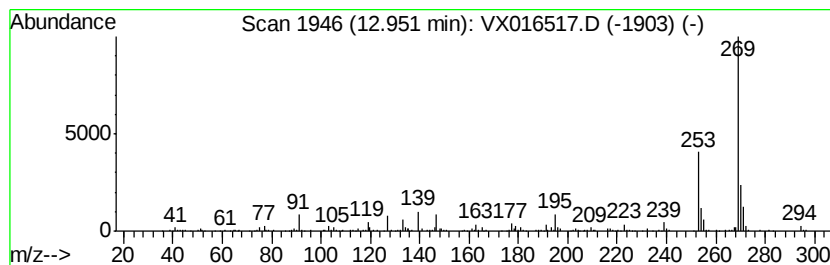
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X052720W.M
Quant Title : SW846 8260

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

Peak Number 1 unknown12.95 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.95	22.11 ug/l	636958	1,4-Dichlorobenzene-d4	12.06	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Phenyl-12-oxa-4-azatricyclo[5.4.1.0(2,6)]dodec-	269	C16H15N03	1000210-56-6	43
2	Phenol, 4,4'-(1-methylethylidene)bis[2,6-dimethyl-	284	C19H24O2	005613-46-7	43
3	2,3-Diphenylindole	269	C20H15N	003469-20-3	37
4	Morpholine, 4,4',4''-borvlydynet...	269	C12H24BN3O3	038761-87-4	35
5	Physcion-10,10'-bianthrone	538	C32H26O8	021871-90-9	32



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown12.95	12.95	22.1	ug/l	636958	4	12.06	1440540	50.0