

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD051519\
 Data File : VD062400.D
 Acq On : 15 May 2019 20:49
 Operator : FY/SY
 Sample : K2865-02
 Misc : 3.02g/5mL/MSVOA D/SOIL
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 TP-21-S2

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 >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D050719S.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	1.841	39	42	45	rVB2	7491	13589	1.61%	0.303%
2	2.087	64	67	68	rBV2	6101	9867	1.17%	0.220%
3	2.363	90	95	98	rBV2	5358	16231	1.92%	0.362%
4	2.874	141	147	149	rBV5	3790	12474	1.47%	0.278%
5	2.904	149	150	158	rVB2	5715	10764	1.27%	0.240%
6	3.219	178	182	191	rVB	21714	53412	6.32%	1.190%
7	3.829	238	244	245	rBV	14416	40377	4.77%	0.900%
8	3.868	245	248	254	rVB	39588	98303	11.62%	2.190%
9	6.161	475	481	487	rBV	9546	31063	3.67%	0.692%
10	6.929	552	559	565	rBV	251119	681971	80.63%	15.196%
11	7.047	565	571	584	rVB	280227	782524	92.52%	17.436%
12	7.460	607	613	628	rBV	149113	423522	50.07%	9.437%
13	8.218	684	690	701	rBV	330913	845785	100.00%	18.846%
14	10.412	908	913	919	rBV	310765	657090	77.69%	14.641%
15	10.511	919	923	931	rVB	55433	122675	14.50%	2.733%
16	11.229	992	996	1002	rBV	10725	31164	3.68%	0.694%
17	12.371	1108	1112	1120	rVB	269322	418335	49.46%	9.321%
18	13.561	1229	1233	1237	rBV	80193	110330	13.04%	2.458%
19	13.719	1245	1249	1252	rVB	15733	26099	3.09%	0.582%
20	14.516	1327	1330	1333	rBV	57291	71482	8.45%	1.593%
21	15.116	1388	1391	1398	rVB	14664	22029	2.60%	0.491%
22	16.297	1508	1511	1515	rVB	6558	8796	1.04%	0.196%

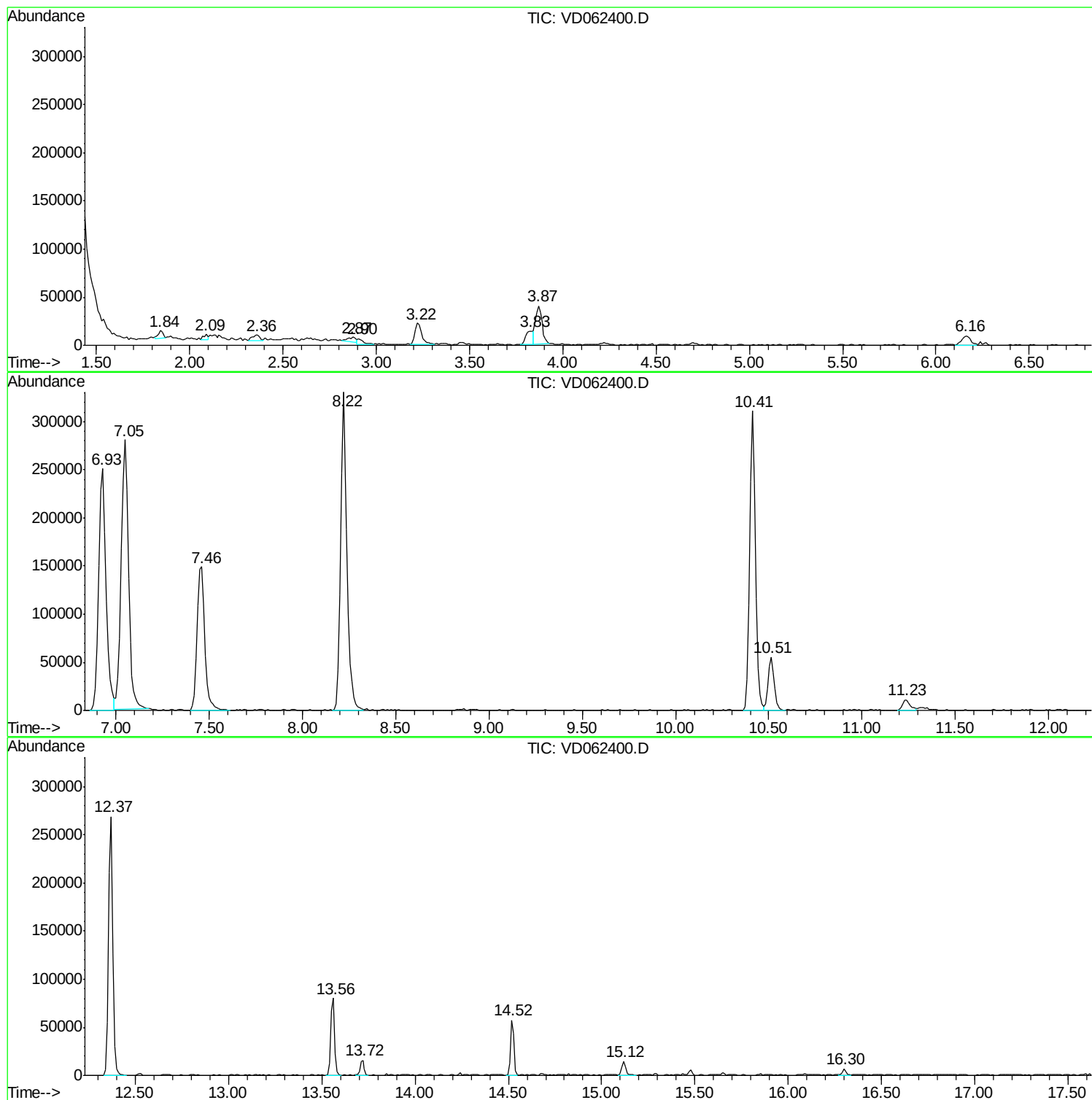
Sum of corrected areas: 4487882

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D050719S.M
Quant Title : SW846 8260

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



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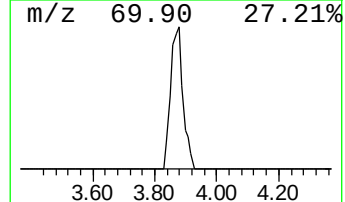
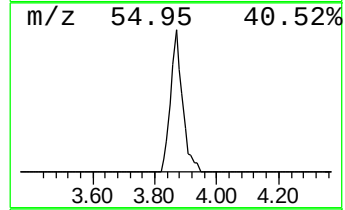
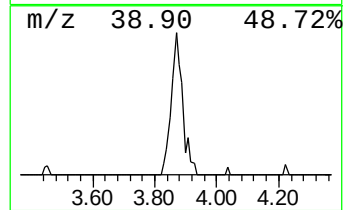
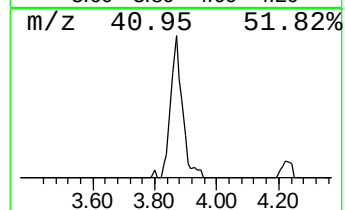
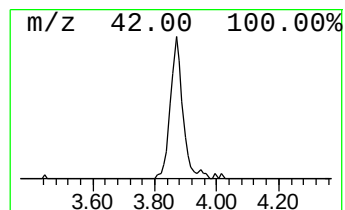
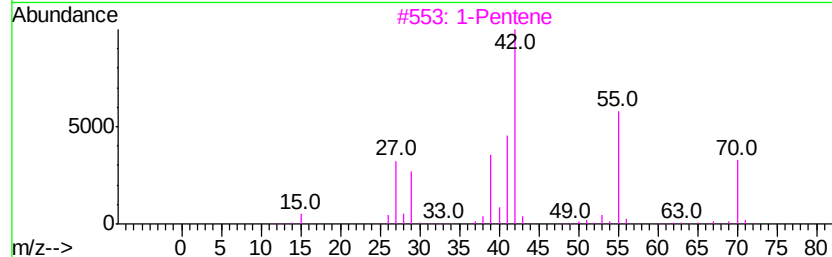
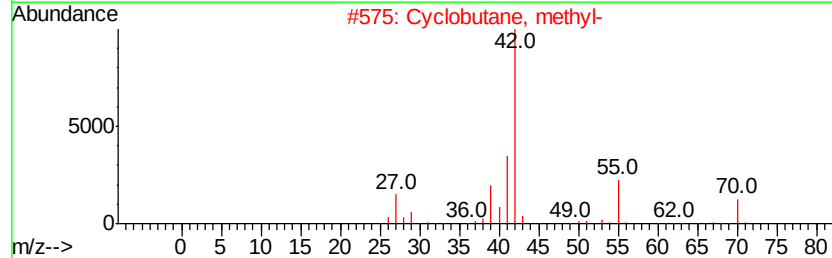
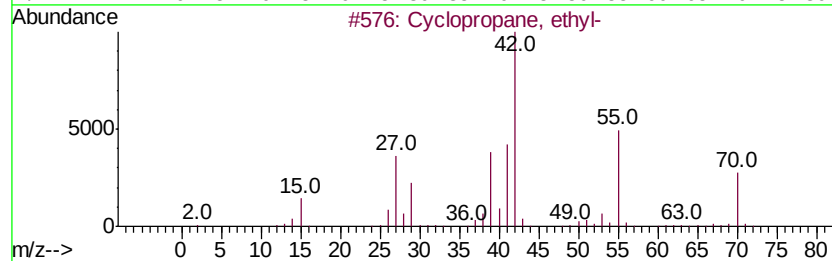
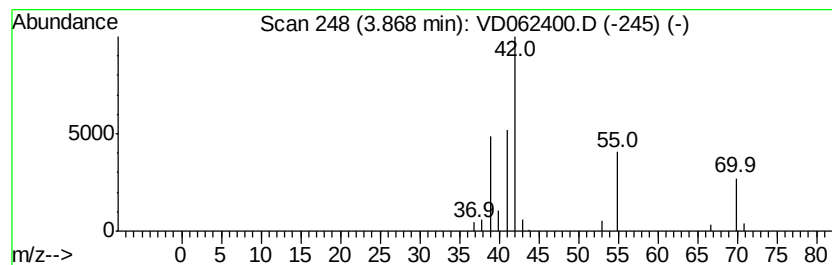
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D050719S.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclopropane, ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.		
3.87	6.28 ug/l	98303	Pentafluorobenzene	7.05		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopropane, ethyl-	70	C5H10	001191-96-4	59
2		Cyclobutane, methyl-	70	C5H10	000598-61-8	58
3		1-Pentene	70	C5H10	000109-67-1	50
4		Methacrolein	70	C4H6O	000078-85-3	43
5		Cyclobutanone	70	C4H6O	001191-95-3	38



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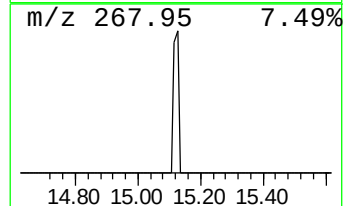
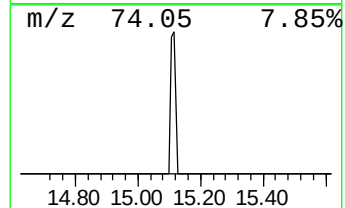
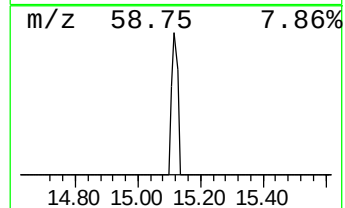
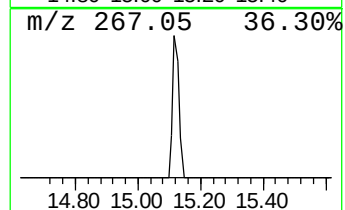
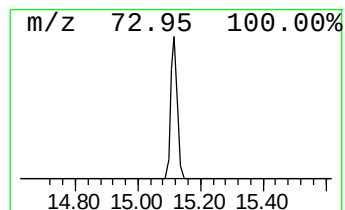
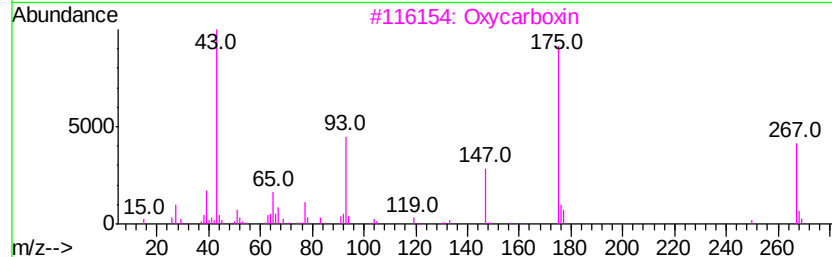
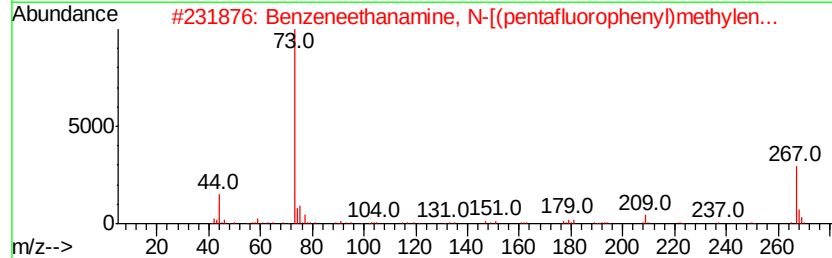
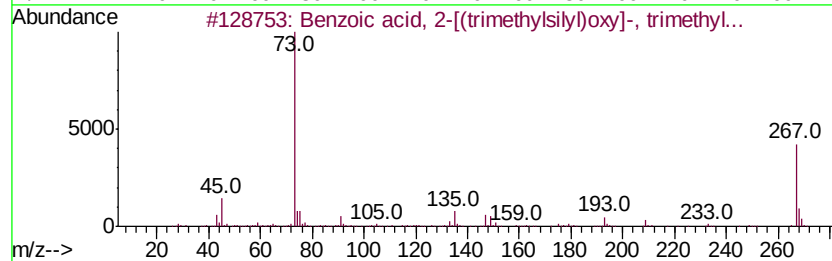
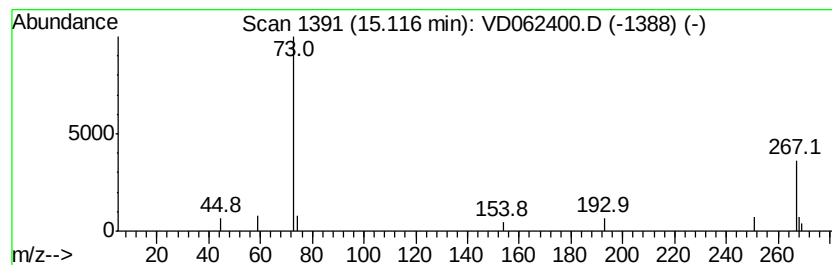
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Peak Number 3 unknown15.12 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.		
15.12	15.41 ug/l	22029	1,4-Dichlorobenzene-d4	14.52		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid, 2-[(trimethylsilyl)oxy]-, trimethyl...	282	C13H22O3Si2	003789-85-3	43	
2	Benzenethanamine, N-[(pentafluorophenyl)methyl]-	475	C21H26F5N02Si2	055429-85-1	9	
3	Oxycarboxin	267	C12H13N04S	005259-88-1	5	
4	Acetamide, N-methyl-	73	C3H7NO	000079-16-3	4	
5	N-Ethylformamide	73	C3H7NO	000627-45-2	4	



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Cyclopropane, ethyl-	3.87	6.3	ug/l	98303	1	7.05	782524	50.0
unknown15.12	15.12	15.4	ug/l	22029	4	14.52	71482	50.0