

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD070924\
Data File : VD079367.D
Acq On : 09 Jul 2024 17:54
Operator : RP/MD
Sample : P3170-03
Misc : 5.05G/5.0ml/MSVOA_D/SOIL/A
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
SOIL-DRUMS-2

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\82D070824S.M

Title : SW846 8260

Signal : TIC: VD079367.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.740	18	26	36	rVB	229323	359539	52.67%	32.044%
2	7.799	1049	1056	1063	rBV5	2666	6857	1.00%	0.611%
3	8.222	1122	1128	1138	rVB3	3442	7689	1.13%	0.685%
4	10.263	1469	1475	1482	rVV7	9516	22086	3.24%	1.968%
5	11.581	1695	1699	1703	rBV6	10121	17904	2.62%	1.596%
6	12.204	1803	1805	1807	rBV3	6977	8171	1.20%	0.728%
7	13.616	2035	2045	2057	rVB5	169515	682677	100.00%	60.843%
8	13.763	2068	2070	2071	rBV	21653	17101	2.50%	1.524%

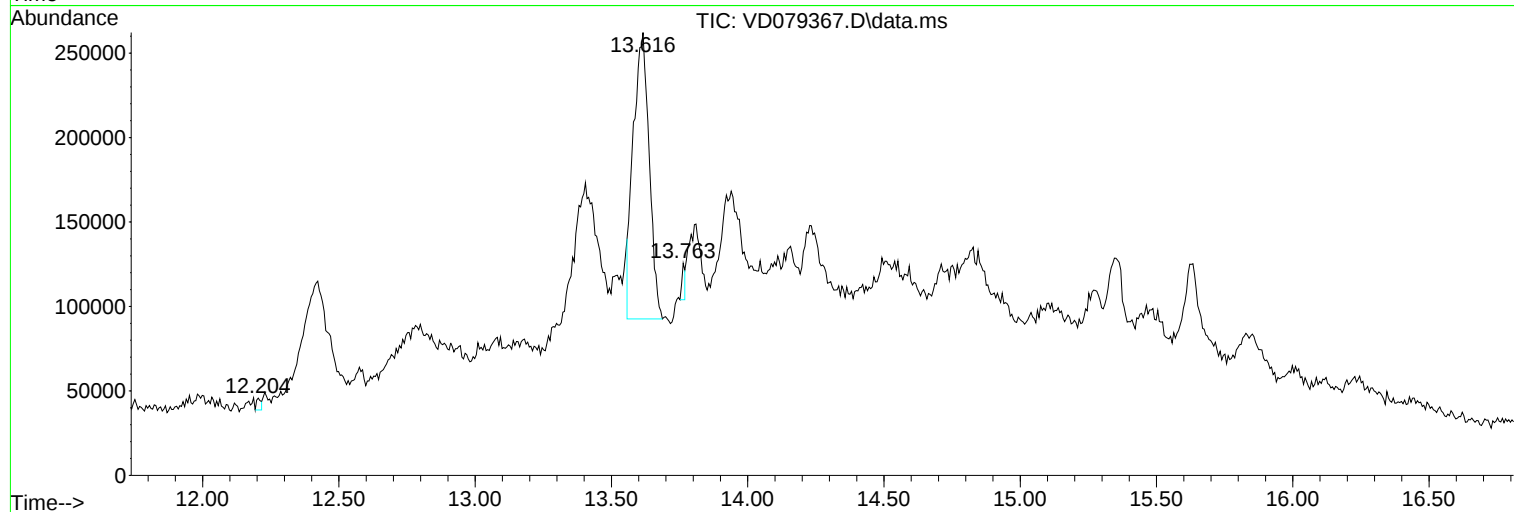
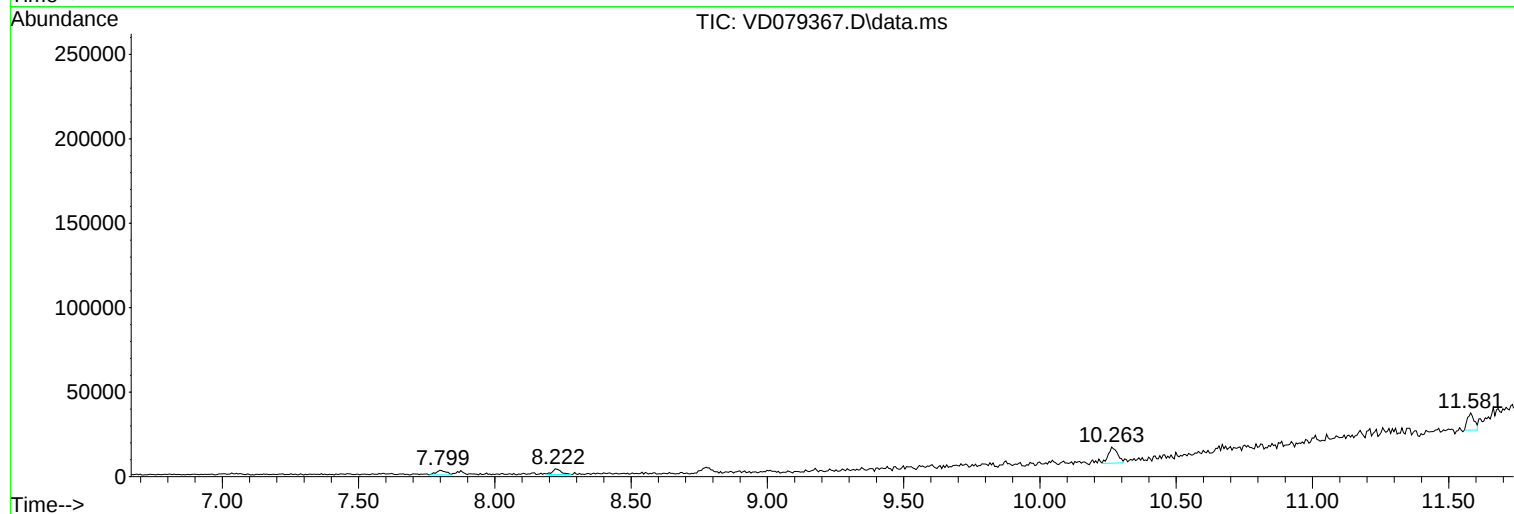
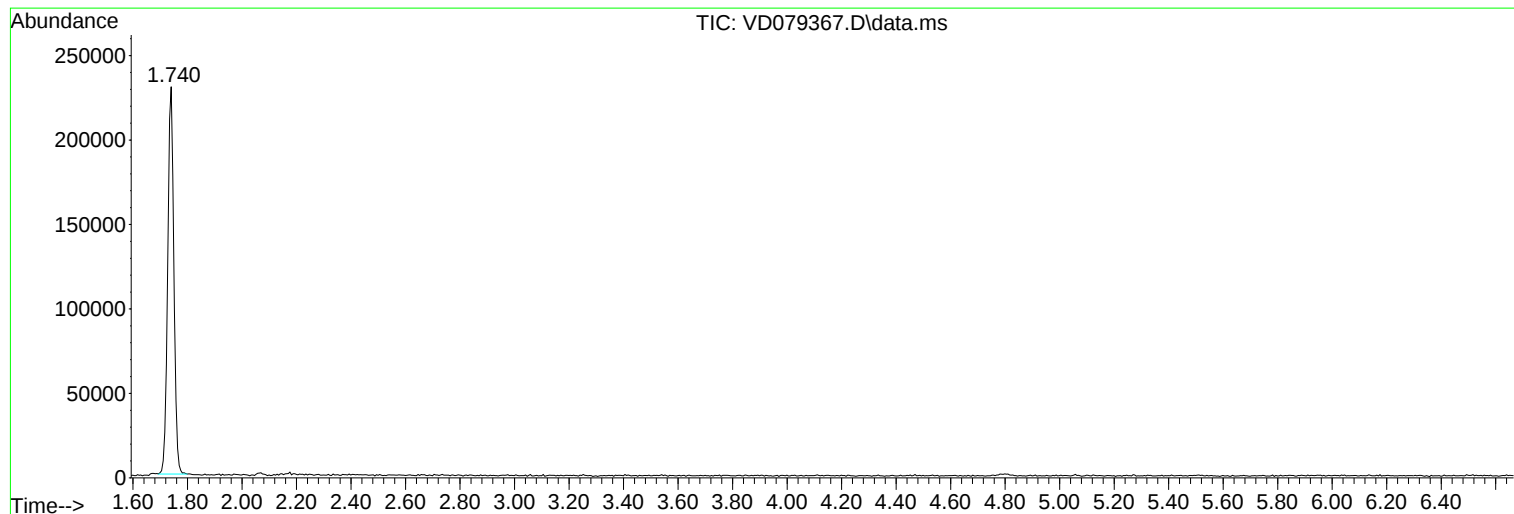
Sum of corrected areas: 1122024

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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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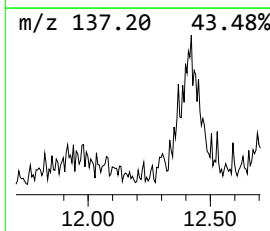
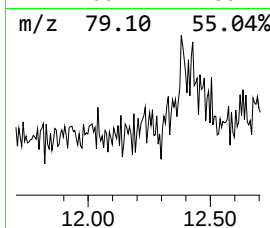
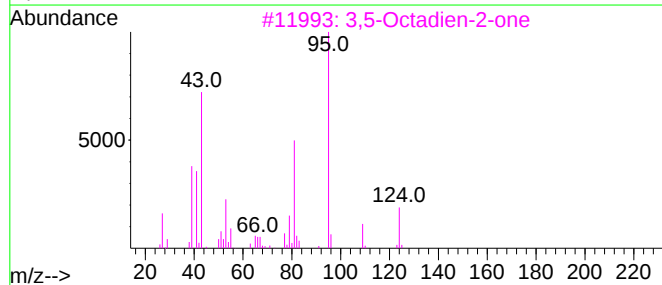
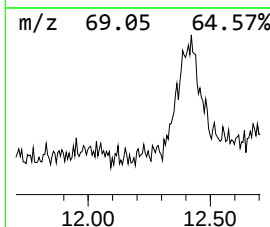
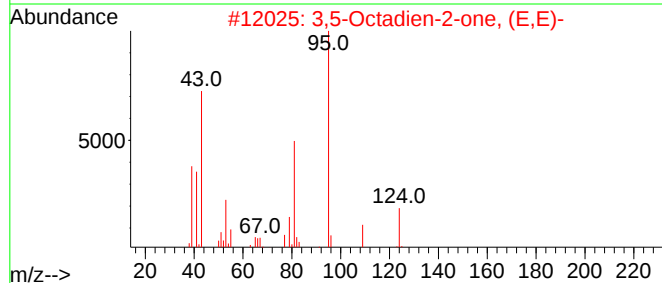
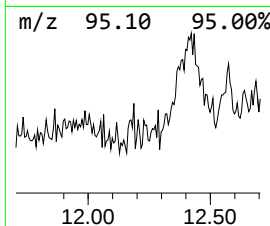
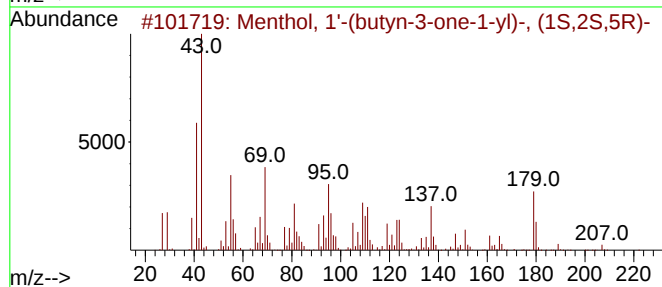
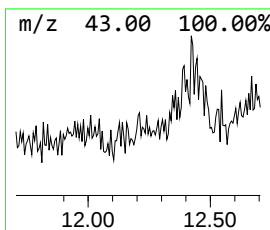
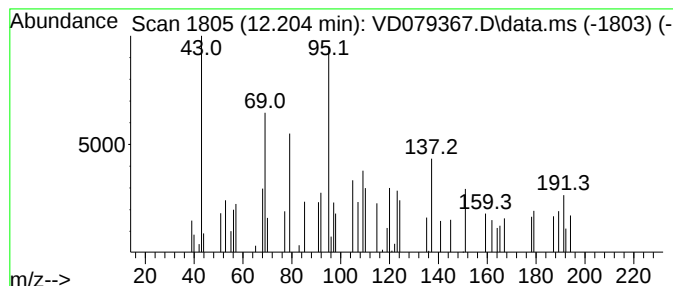
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D070824S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown12.204 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.204	22.82 ug/l	8171	Chlorobenzene-d5	11.581

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Menthol, 1'-(butyn-3-one-1-yl)-,...	222	C14H22O2	1000156-95-3	32
2			3,5-Octadien-2-one, (E,E)-	124	C8H12O	030086-02-3	27
3			3,5-Octadien-2-one	124	C8H12O	038284-27-4	27
4			Pentaleno[1,2-b]oxirene, octahyd...	124	C8H12O	055449-71-3	22
5			1,3-Hexadiene, 2,5-dimethyl-	110	C8H14	029253-64-3	14



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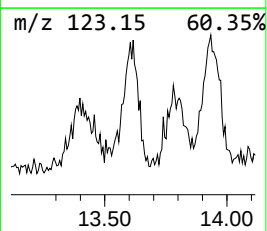
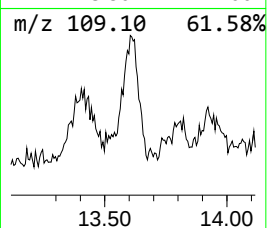
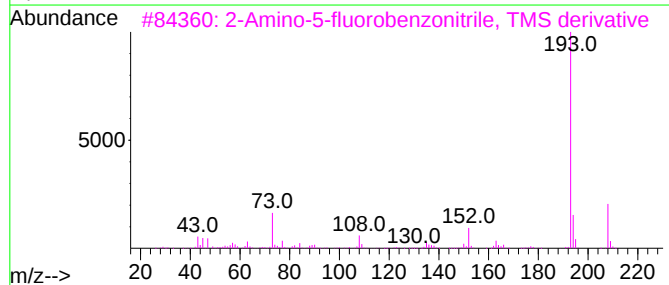
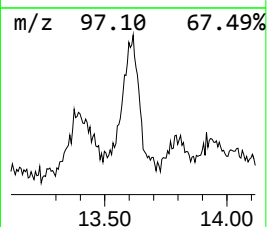
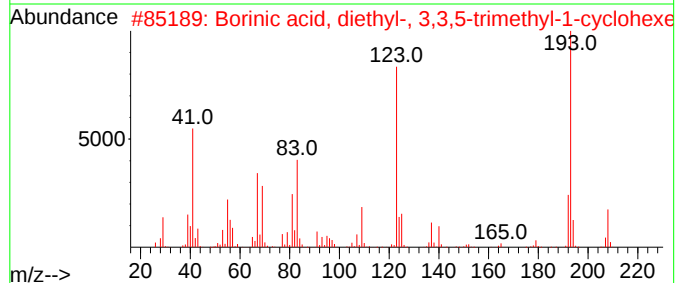
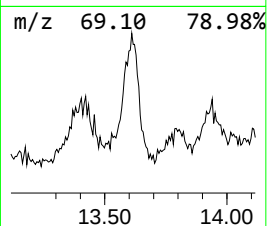
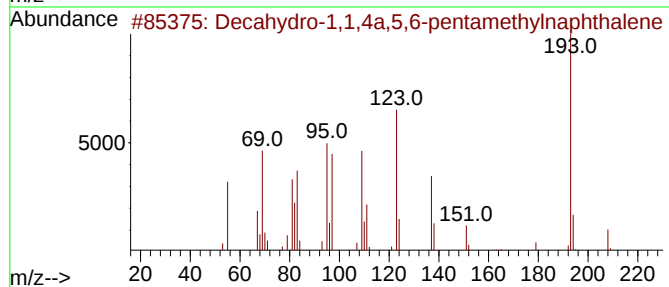
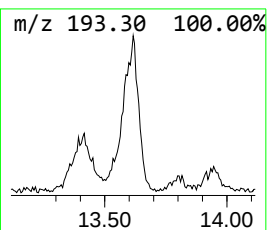
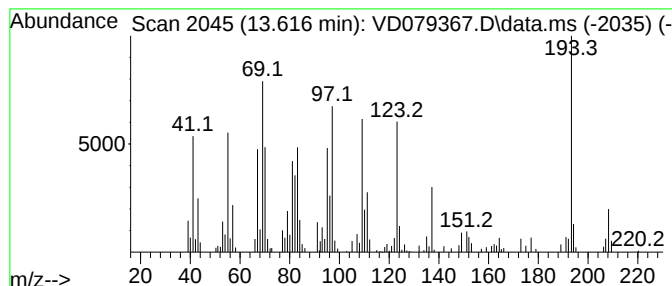
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ClientSampleId :
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TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown13.616 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.616	50.00 ug/l	682677	1,4-Dichlorobenzene-d4	13.251	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	47
2	Borinic acid, diethyl-, 3,3,5-tr...	208	C13H25BO	057387-76-5	38
3	2-Amino-5-fluorobenzonitrile, TM...	208	C10H13FN2Si	1000479-05-6	25
4	1-Cyclopentyl-4-(1-methylethyl)c...	194	C14H26	1000215-44-4	25
5	2(5H)-Furanone, 5,5-dimethyl-	112	C6H8O2	020019-64-1	18



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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown12.204	12.204	22.8	ug/l	8171	3	11.581	17904	50.0
unknown13.616	13.616	50.0	ug/l	682677	4	13.251	682677	50.0