

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD062519\
 Data File : VD062874.D
 Acq On : 25 Jun 2019 12:58
 Operator : VA/AP
 Sample : K3164-09
 Misc : 4.51g/5ml/MSVOA D/SOIL
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 BAT-B08

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_D\METHOD\82D061919S.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	1.473	3	5	23	rVB	379610	1094457	37.00%	4.901%
2	2.349	89	94	97	rBV2	64385	139496	4.72%	0.625%
3	2.625	119	122	125	rBV2	20447	37324	1.26%	0.167%
4	3.225	179	183	190	rVB	57068	149609	5.06%	0.670%
5	3.471	202	208	217	rVB3	11497	43969	1.49%	0.197%
6	3.865	238	248	259	rBV	450386	1515137	51.23%	6.785%
7	4.682	324	331	337	rBV3	9049	31590	1.07%	0.141%
8	6.169	475	482	490	rBV2	68677	226691	7.66%	1.015%
9	6.927	546	559	566	rBV2	523328	1516461	51.27%	6.791%
10	7.055	566	572	587	rVB	712497	2101383	71.05%	9.410%
11	7.458	607	613	624	rBV	397005	1048374	35.45%	4.694%
12	8.226	684	691	706	rBV	993060	2495045	84.36%	11.172%
13	8.541	718	723	732	rVB2	39043	92581	3.13%	0.415%
14	10.126	880	884	890	rVB2	17577	49382	1.67%	0.221%
15	10.412	907	913	921	rBV	1208553	2957732	100.00%	13.244%
16	10.520	921	924	928	rVB	25266	51347	1.74%	0.230%
17	10.648	932	937	942	rBV4	18130	64079	2.17%	0.287%
18	10.795	948	952	959	rVB4	23325	58709	1.98%	0.263%
19	10.933	959	966	973	rBV3	53569	134525	4.55%	0.602%
20	11.110	979	984	989	rVB3	29223	66486	2.25%	0.298%
21	11.238	989	997	1002	rBV	27453	86644	2.93%	0.388%
22	11.760	1046	1050	1054	rVB2	56073	119292	4.03%	0.534%
23	11.839	1054	1058	1062	rBV	26819	56784	1.92%	0.254%
24	12.371	1108	1112	1119	rBV	1683163	2692895	91.05%	12.058%
25	12.469	1119	1122	1124	rVV2	17614	30995	1.05%	0.139%
26	12.528	1124	1128	1133	rVB	83849	137072	4.63%	0.614%
27	12.666	1138	1142	1148	rBV	320595	549567	18.58%	2.461%
28	13.050	1178	1181	1185	rVB	159410	225594	7.63%	1.010%
29	13.562	1229	1233	1240	rBV	1444856	1962397	66.35%	8.787%
30	13.719	1245	1249	1253	rVB	71609	115520	3.91%	0.517%
31	14.517	1327	1330	1339	rVB	1761591	2397264	81.05%	10.735%
32	15.117	1388	1391	1397	rVB	62015	83690	2.83%	0.375%

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Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D061919S.M
Title : SW846 8260

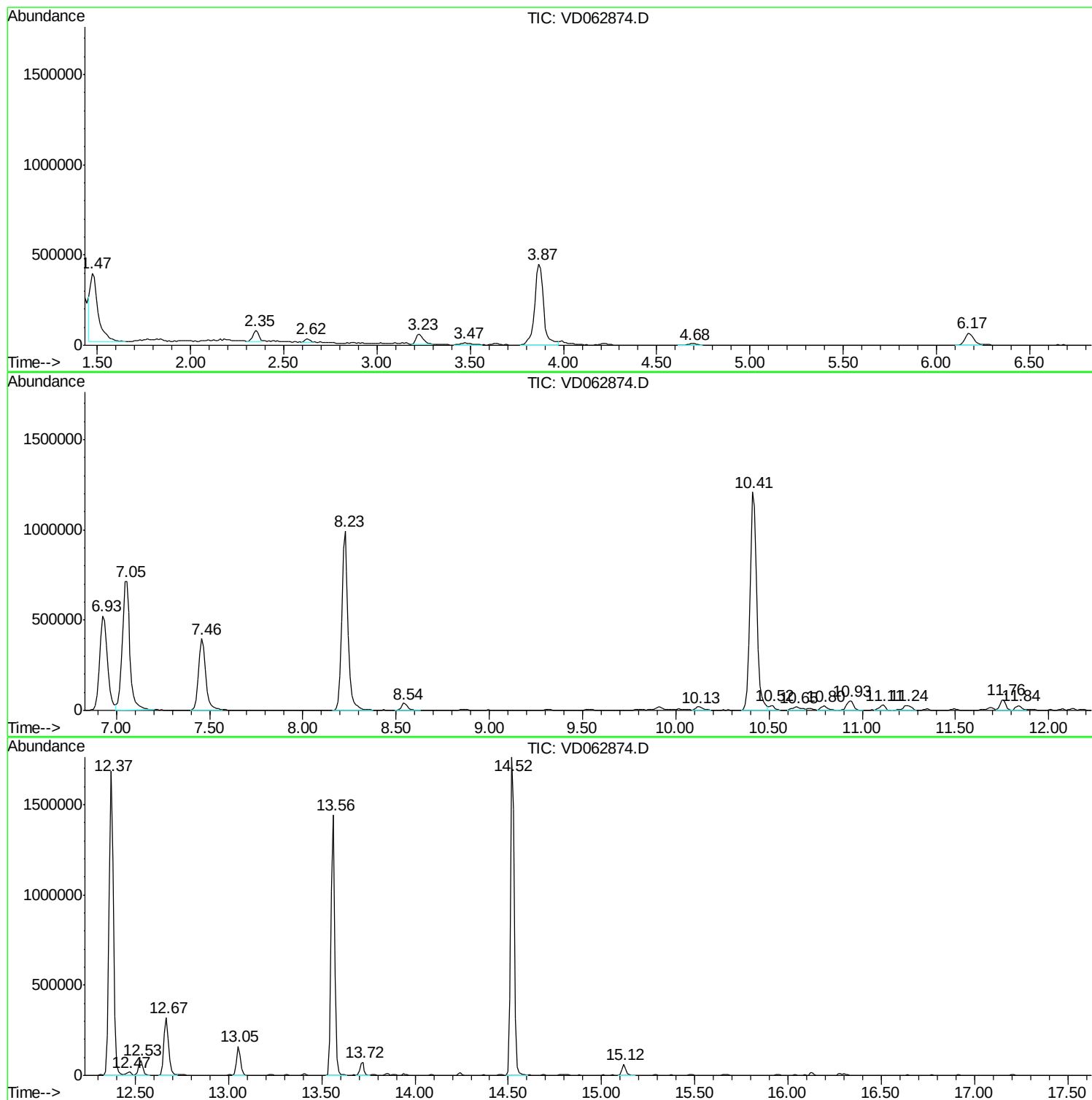
Sum of corrected areas: 22332091

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D061919S.M
Quant Title : SW846 8260

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



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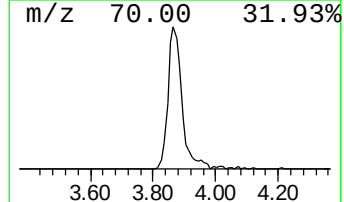
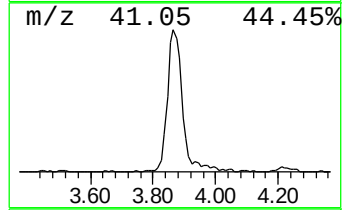
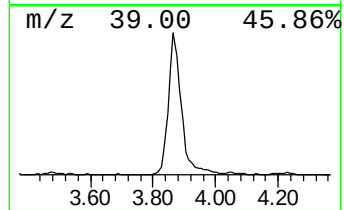
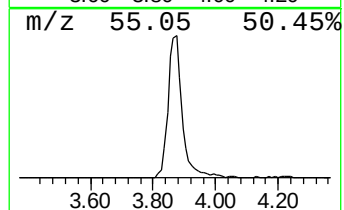
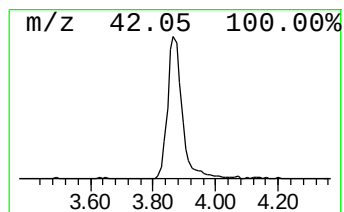
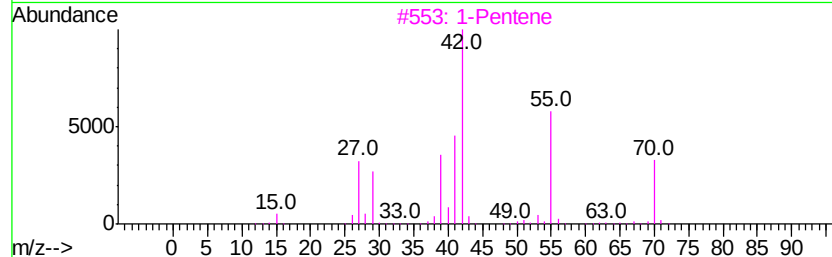
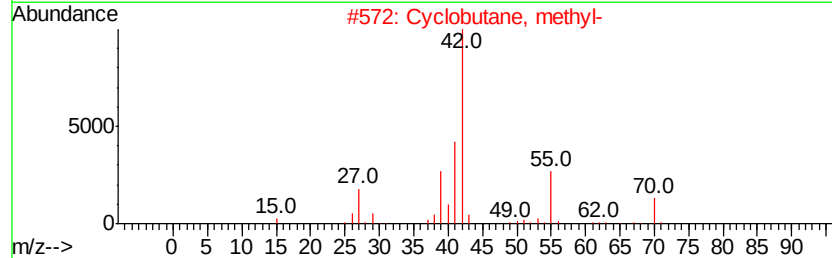
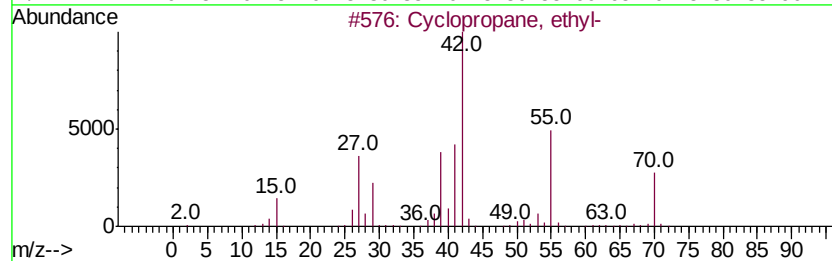
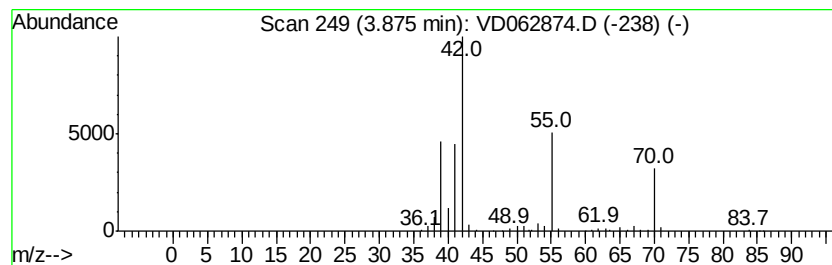
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TIC Integration Parameters: LSCINT.P

Peak Number 2 Cyclopropane, ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.87	36.05 ug/l	1515140	Pentafluorobenzene	7.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopropane, ethyl-	70	C5H10	001191-96-4	78
2		Cyclobutane, methyl-	70	C5H10	000598-61-8	59
3		1-Pentene	70	C5H10	000109-67-1	58
4		Aziridine, 2-ethyl-	71	C4H9N	002549-67-9	33
5		4-Pentenal	84	C5H8O	002100-17-6	12



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Cyclopropane, ethyl-	3.87	36.0	ug/l	1515140	1	7.05	2101380	50.0