

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD080819\
 Data File : VD063449.D
 Acq On : 8 Aug 2019 20:08
 Operator : JC/SY
 Sample : K4223-18
 Misc : 6.38g/5ml/MSVOA D/SOIL
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 SP-9

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\82D072919S.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	1.483	3	6	25	rVB	2077821	5884350	100.00%	22.284%
2	3.255	177	186	197	rBV2	140447	532302	9.05%	2.016%
3	6.080	466	473	483	rBV2	22269	65338	1.11%	0.247%
4	6.927	552	559	566	rBV	536839	1540907	26.19%	5.835%
5	7.045	566	571	585	rVB	943066	2810664	47.77%	10.644%
6	7.458	607	613	625	rBV	367475	1053710	17.91%	3.990%
7	8.226	684	691	706	rBV	1286401	3186558	54.15%	12.067%
8	10.412	907	913	920	rBV	1103681	2536651	43.11%	9.606%
9	10.510	920	923	932	rVB2	23119	59696	1.01%	0.226%
10	12.371	1108	1112	1121	rBV	1582856	2628832	44.67%	9.955%
11	13.227	1193	1199	1202	rBV3	37364	74218	1.26%	0.281%
12	13.316	1202	1208	1209	rBV2	34011	77594	1.32%	0.294%
13	13.483	1220	1225	1229	rBV3	99787	172933	2.94%	0.655%
14	13.562	1229	1233	1237	rBV	1087757	1515935	25.76%	5.741%
15	13.719	1246	1249	1253	rVB2	76616	115734	1.97%	0.438%
16	13.808	1255	1258	1264	rBV5	26318	97440	1.66%	0.369%
17	13.926	1264	1270	1273	rVV6	68632	206981	3.52%	0.784%
18	13.975	1273	1275	1277	rVB2	47999	64777	1.10%	0.245%
19	14.212	1293	1299	1301	rVV2	78157	153906	2.62%	0.583%
20	14.261	1301	1304	1309	rVB6	39077	79658	1.35%	0.302%
21	14.379	1309	1316	1319	rBV5	43848	132877	2.26%	0.503%
22	14.438	1319	1322	1324	rVV2	106620	168072	2.86%	0.636%
23	14.517	1324	1330	1334	rVV	1254440	2005217	34.08%	7.594%
24	14.783	1355	1357	1358	rBV	44010	62351	1.06%	0.236%
25	14.812	1358	1360	1366	rVV6	88368	195256	3.32%	0.739%
26	14.891	1366	1368	1370	rVV2	37650	73311	1.25%	0.278%
27	14.950	1372	1374	1377	rVV2	51815	63605	1.08%	0.241%
28	15.019	1377	1381	1383	rVV4	37657	82629	1.40%	0.313%
29	15.058	1383	1385	1387	rVB2	43539	63724	1.08%	0.241%
30	15.147	1390	1394	1398	rVB6	43855	94668	1.61%	0.359%
31	15.226	1398	1402	1405	rVB6	38905	76856	1.31%	0.291%
32	15.285	1405	1408	1410	rBV4	83908	136301	2.32%	0.516%
33	15.324	1410	1412	1416	rVB3	30849	62835	1.07%	0.238%
34	15.432	1420	1423	1426	rBV2	87653	119394	2.03%	0.452%

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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 >
Peak separation: 5

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Title : SW846 8260

35	15.678	1444	1448	1450	rBV4	31179	66741	1.13%	0.253%
36	15.806	1458	1461	1468	rVB7	34547	68097	1.16%	0.258%
37	16.230	1500	1504	1507	rVV6	39153	76472	1.30%	0.290%

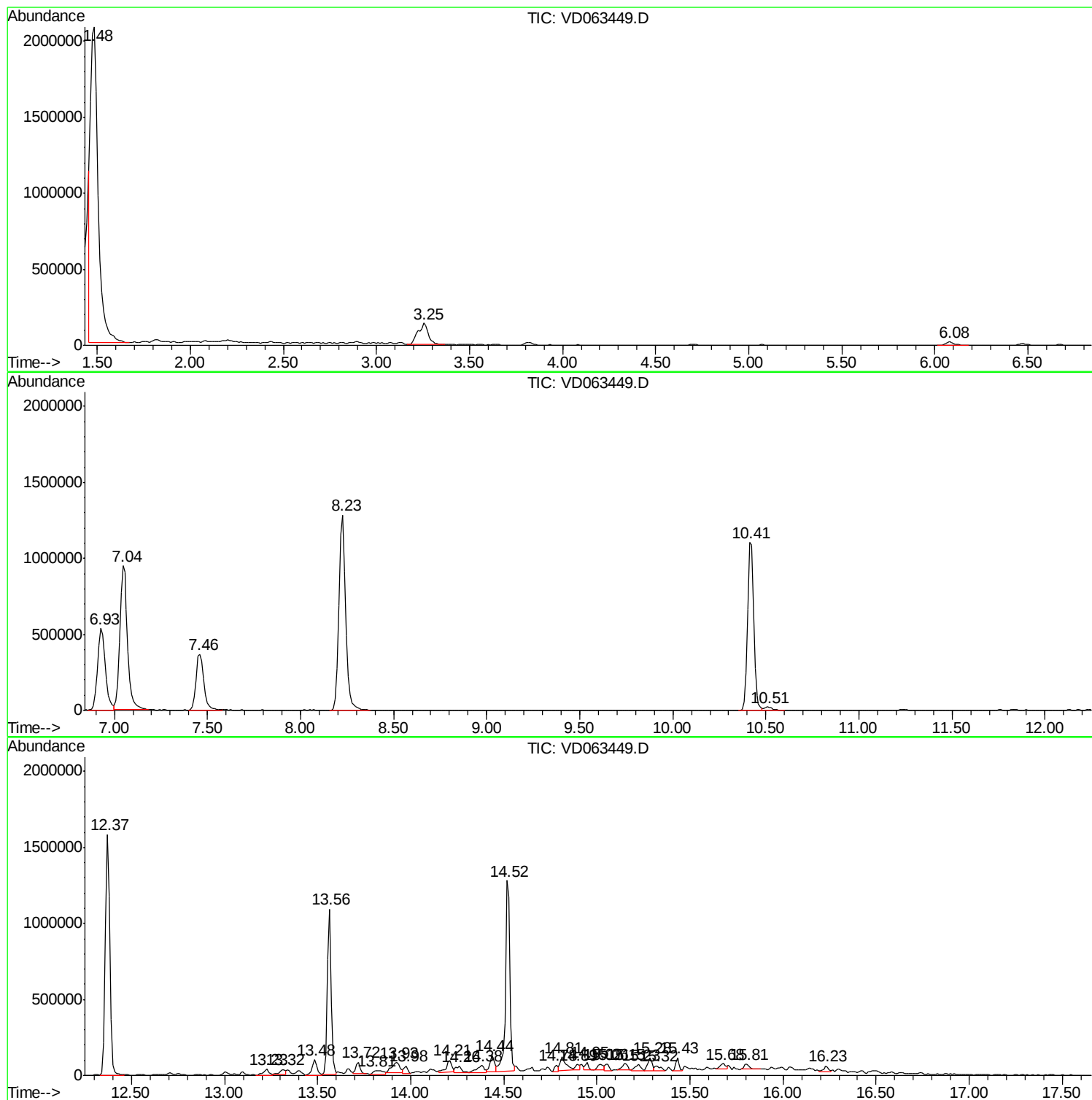
Sum of corrected areas: 26406590

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D072919S.M
Quant Title : SW846 8260

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



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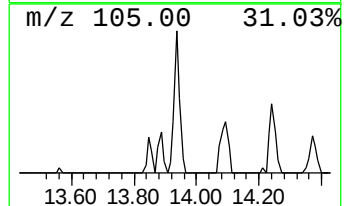
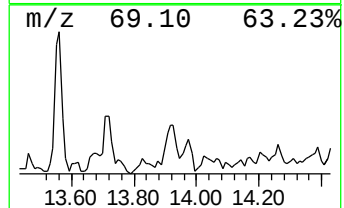
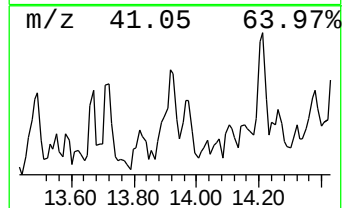
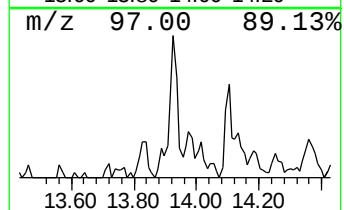
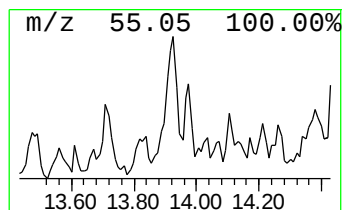
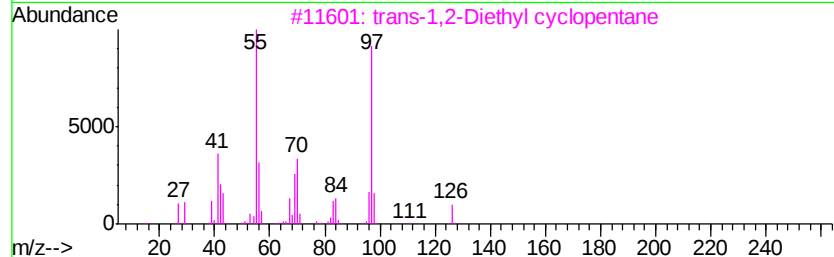
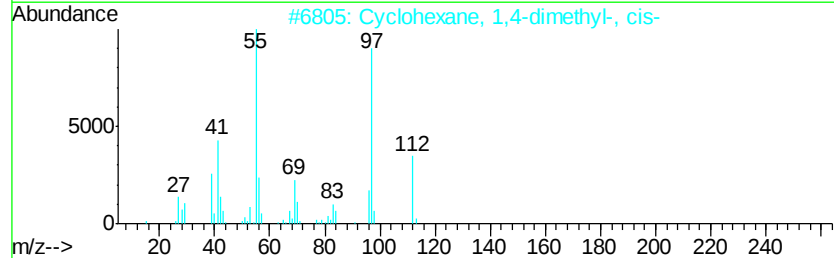
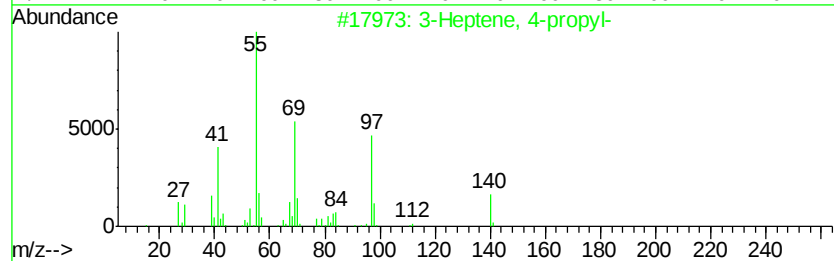
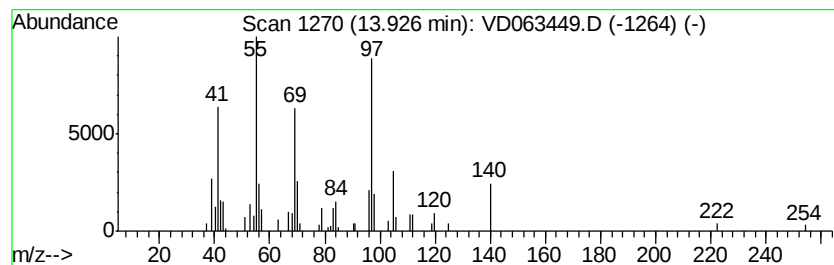
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TIC Integration Parameters: LSCINT.P

Peak Number 2 3-Heptene, 4-propyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.93	5.16 ug/l	206981	1,4-Dichlorobenzene-d4	14.53

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Heptene, 4-propyl-	140	C10H20	004485-13-6	64
2			Cyclohexane, 1,4-dimethyl-, cis-	112	C8H16	000624-29-3	52
3			trans-1,2-Diethyl cyclopentane	126	C9H18	000932-40-1	50
4			3-Hexene, 3-ethyl-2,5-dimethyl-	140	C10H20	062338-08-3	46
5			Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	46



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3-Heptene, 4-propyl-	13.93	5.2	ug/l	206981	4	14.53	2005220	50.0