

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_Y\DATA\VY110619\
 Data File : VY000564.D
 Acq On : 06 Nov 2019 19:58
 Operator : SY/MD
 Sample : K5723-05
 Misc : 7.49G/5ML/MSVOA_Y/SOIL
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 8-B

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_Y\METHODS\82Y103019S.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	1.865	17	21	28	rVB3	12969	19929	1.56%	0.288%
2	3.536	289	295	305	rVB4	5410	14072	1.10%	0.203%
3	4.718	478	489	503	rBV	82341	243525	19.10%	3.518%
4	5.755	645	659	675	rVB2	36235	110529	8.67%	1.597%
5	7.730	973	983	989	rBV	174442	415791	32.60%	6.006%
6	7.803	989	995	1008	rVB	290269	652564	51.17%	9.427%
7	8.151	1041	1052	1063	rBV	182660	398779	31.27%	5.761%
8	8.699	1133	1142	1153	rBV	477624	930101	72.93%	13.436%
9	10.181	1377	1385	1393	rBV	738704	1275266	100.00%	18.422%
10	10.583	1445	1451	1461	rBV3	7492	13241	1.04%	0.191%
11	11.492	1592	1600	1608	rBV	655688	1082401	84.88%	15.636%
12	12.479	1755	1762	1776	rVB	505602	802689	62.94%	11.595%
13	13.424	1910	1917	1924	rVV	608735	932304	73.11%	13.468%
14	13.961	2001	2005	2012	rVB2	10094	15817	1.24%	0.228%
15	16.149	2359	2364	2369	rBV7	7918	15504	1.22%	0.224%

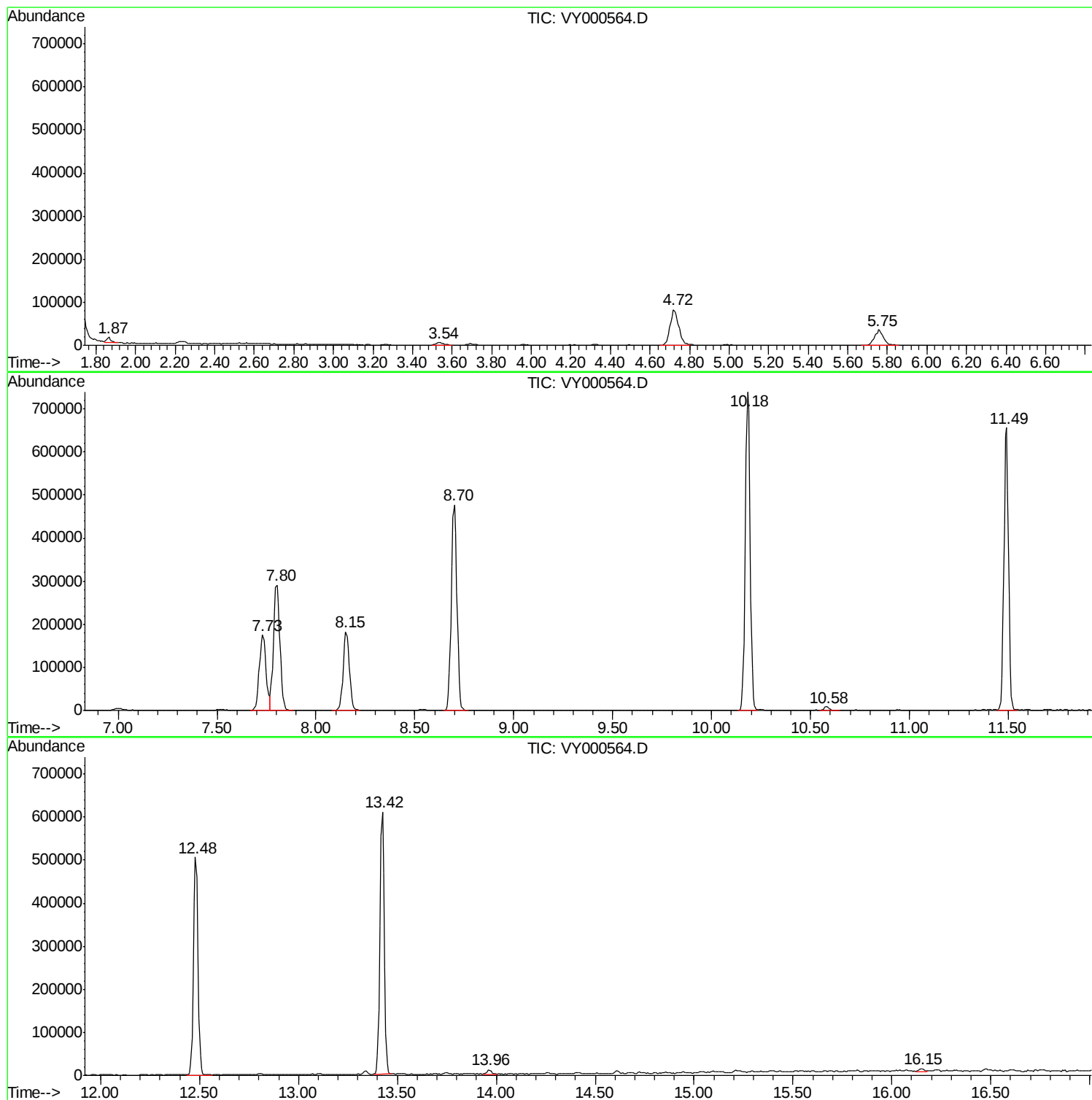
Sum of corrected areas: 6922512

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Quant Method : Z:\V0ASRV\HPCHEM1\MSV0A_Y\METHODS\82Y103019S.M
Quant Title : SW846 8260

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



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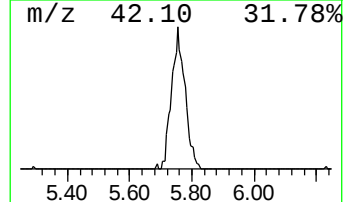
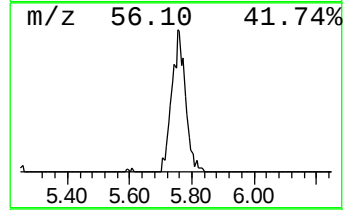
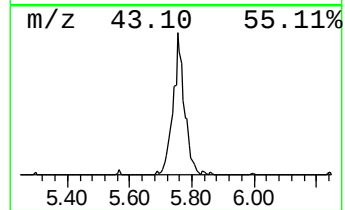
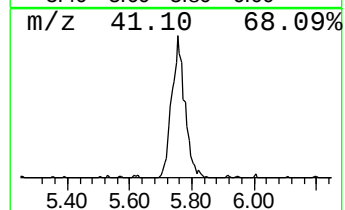
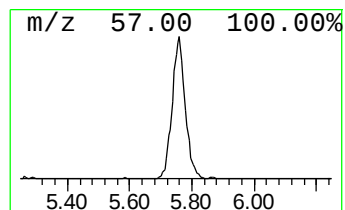
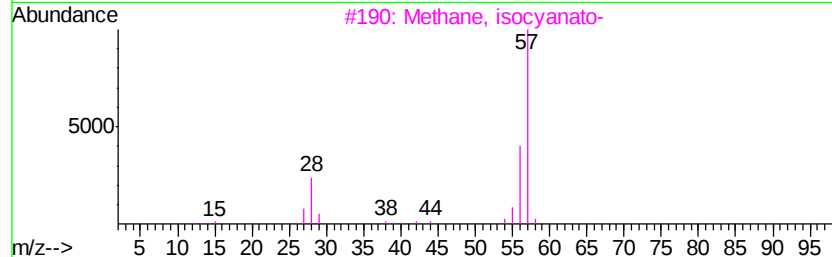
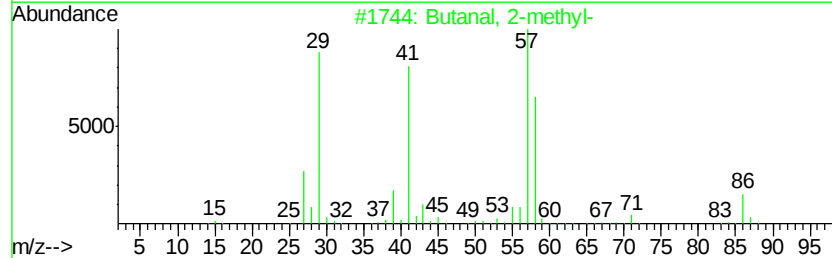
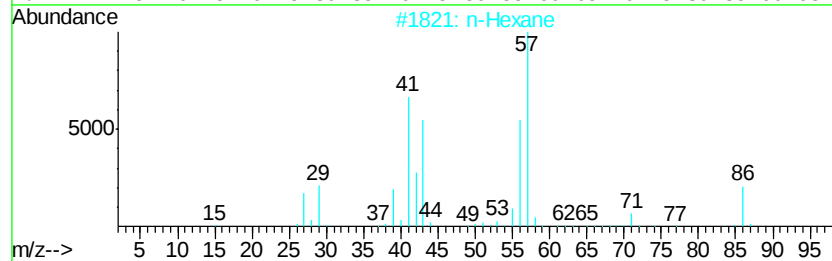
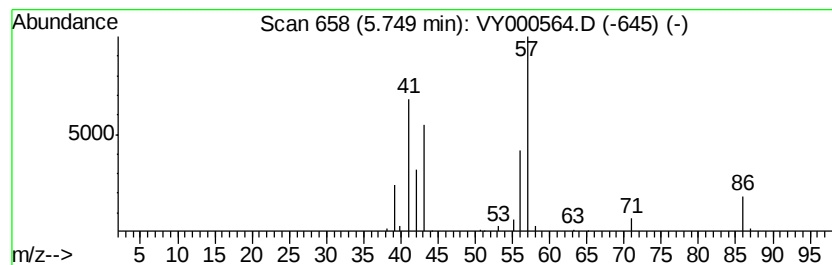
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Peak Number 1 n-Hexane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.75	8.47 ug/l	110529	Pentafluorobenzene	7.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexane	86	C6H14	000110-54-3	87
2		Butanal, 2-methyl-	86	C5H10O	000096-17-3	38
3		Methane, isocyanato-	57	C2H3NO	000624-83-9	35
4		Propanal, 2,2-dimethyl-	86	C5H10O	000630-19-3	32
5		1-Pentene, 4-methyl-	84	C6H12	000691-37-2	25



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
n-Hexane	5.75	8.5	ug/l	110529	1	7.80	652564	50.0