

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091219\
Data File : VN057941.D
Acq On : 11 Sep 2019 22:01
Operator : JC/SP
Sample : K4835-03
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PZ-3

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_N\METHODS\82N091019W.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	2.294	152	162	181	rVB	42852	68911	2.19%	0.591%
2	2.426	193	203	215	rBV2	21161	37446	1.19%	0.321%
3	3.802	616	631	654	rBV2	73630	193042	6.14%	1.656%
4	4.046	690	707	735	rBV2	42643	129175	4.11%	1.108%
5	4.767	906	931	968	rBV3	288558	1025146	32.62%	8.795%
6	5.021	982	1010	1043	rBV	854698	3142502	100.00%	26.961%
7	7.577	1787	1805	1817	rBV	283995	705477	22.45%	6.053%
8	7.651	1817	1828	1853	rVB	356376	853406	27.16%	7.322%
9	8.014	1924	1941	1965	rBV	370830	859270	27.34%	7.372%
10	8.577	2100	2116	2145	rVB	441109	927976	29.53%	7.962%
11	11.352	2967	2979	2986	rBV	82437	141056	4.49%	1.210%
12	11.406	2986	2996	3018	rVB	637151	1110497	35.34%	9.528%
13	12.406	3294	3307	3327	rBV	631408	1065997	33.92%	9.146%
14	13.348	3587	3600	3618	rBV	834956	1395748	44.42%	11.975%

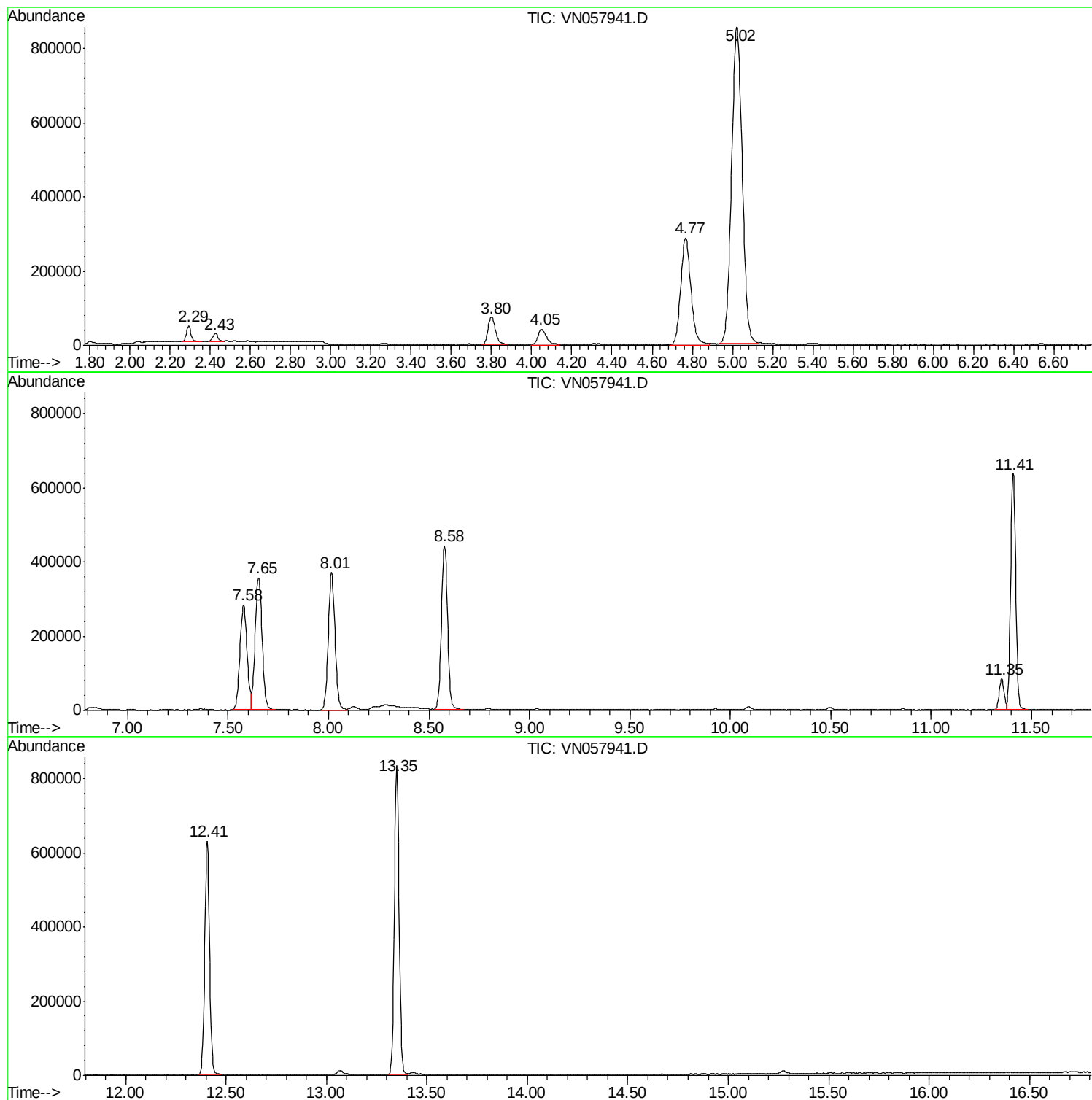
Sum of corrected areas: 11655649

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Quant Title : SW846 8260

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



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Instrument :
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ClientSampled :
PZ-3

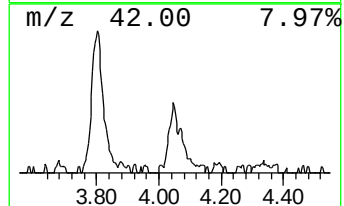
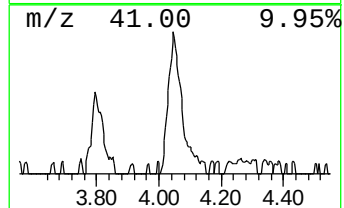
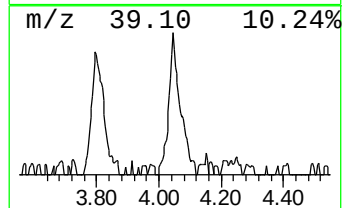
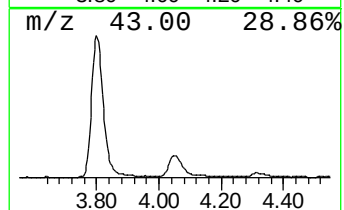
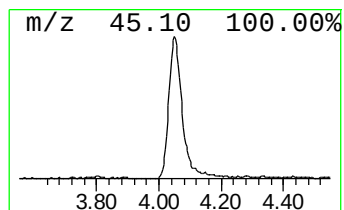
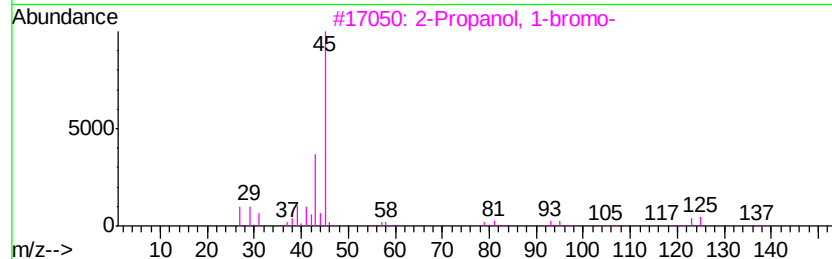
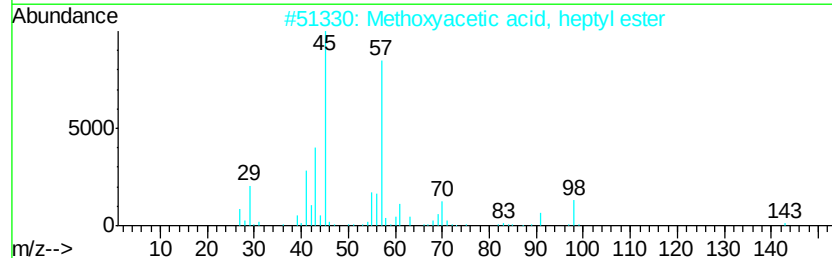
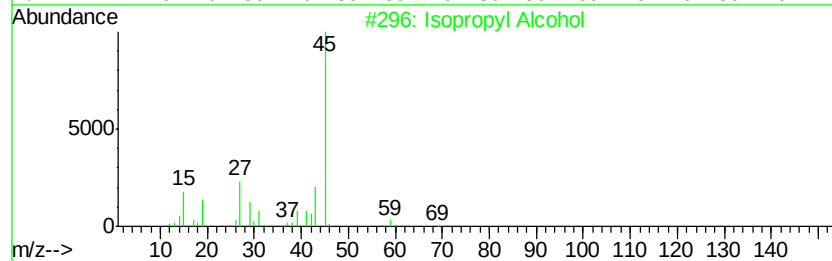
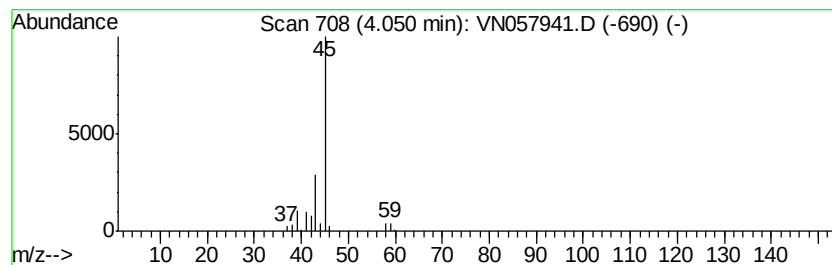
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Isopropyl Alcohol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.05	7.57 ug/l	129175	Pentafluorobenzene	7.65

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isopropyl Alcohol	60	C3H8O	000067-63-0	72
2			Methoxyacetic acid, heptyl ester	188	C10H20O3	168920-36-3	40
3			2-Propanol, 1-bromo-	138	C3H7BrO	019686-73-8	40
4			Formic acid, 1-methylethyl ester	88	C4H8O2	000625-55-8	9
5			2-Methyl-4-methoxy-1-butanol	118	C6H14O2	071052-94-3	9



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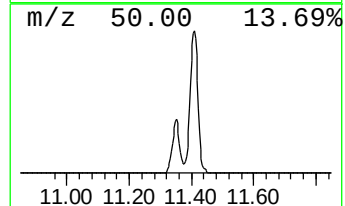
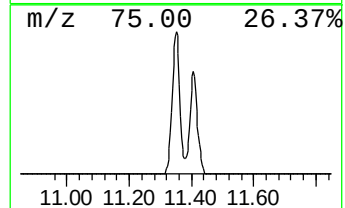
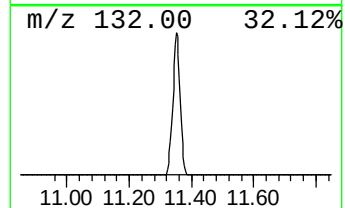
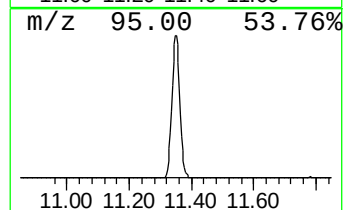
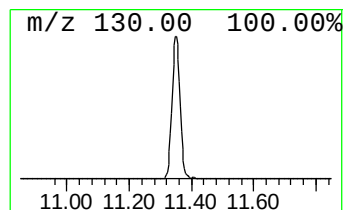
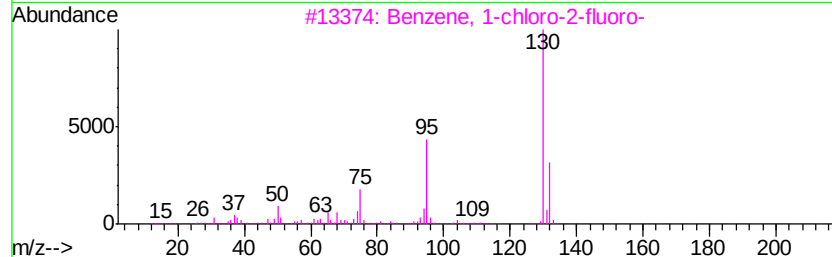
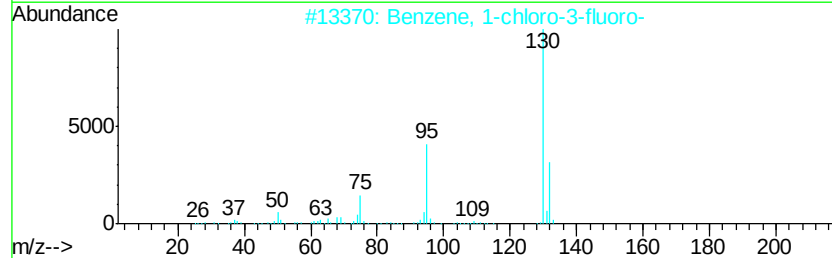
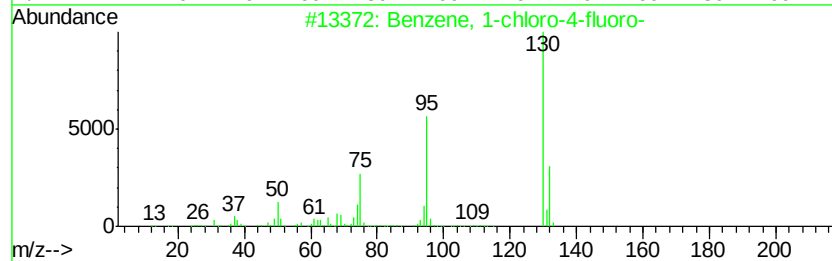
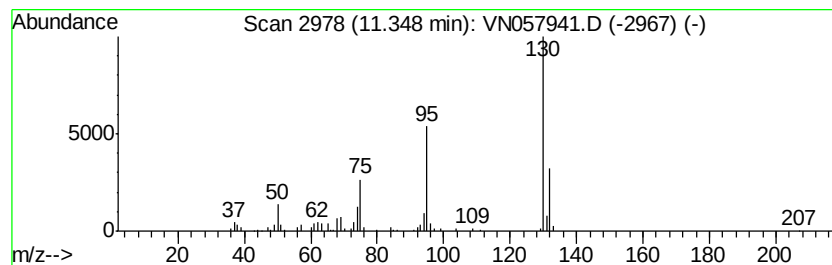
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Peak Number 2 Benzene, 1-chloro-4-fluoro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.35	6.35 ug/l	141056	Chlorobenzene-d5	11.41	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-chloro-4-fluoro-	130	C6H4ClF	000352-33-0	97
2	Benzene, 1-chloro-3-fluoro-	130	C6H4ClF	000625-98-9	96
3	Benzene, 1-chloro-2-fluoro-	130	C6H4ClF	000348-51-6	94
4	3-Chloropyrazine 1-oxide	130	C4H3ClN2O	006863-76-9	40
5	3(2H)-Pyridazinone, 6-chloro-	130	C4H3ClN2O	019064-67-6	38



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Isopropyl Alcohol	4.05	7.6	ug/l	129175	1	7.65	853406	50.0
Benzene, 1-chloro...	11.35	6.3	ug/l	141056	3	11.41	1110500	50.0