

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU110518\
Data File : VU027967.D
Acq On : 05 Nov 2018 19:01
Operator : MD/SY
Sample : J5811-04
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
DPC-91468-OW-01

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_U\METHOD\82U110218W.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	1.469	83	92	112	rBV	1091900	1397368	2.09%	0.908%
2	2.649	439	459	509	rVB2	6722121	18920661	28.29%	12.288%
3	3.048	544	583	611	rVB	1818641	7497017	11.21%	4.869%
4	4.411	987	1007	1075	rVB2	14846403	41776644	62.47%	27.133%
5	5.482	1307	1340	1369	rVB	4000058	10937757	16.36%	7.104%
6	5.887	1452	1466	1495	rBV	420660	917068	1.37%	0.596%
7	6.254	1565	1580	1613	rBV	523635	1123346	1.68%	0.730%
8	7.498	1941	1967	1982	rBV2	22924800	66874354	100.00%	43.433%
9	7.569	1982	1989	2019	rVB	693742	1388733	2.08%	0.902%
10	9.096	2446	2464	2485	rBV	615191	1149229	1.72%	0.746%
11	10.315	2831	2843	2866	rVB	533724	888189	1.33%	0.577%
12	11.488	3195	3208	3226	rBV	701048	1100534	1.65%	0.715%

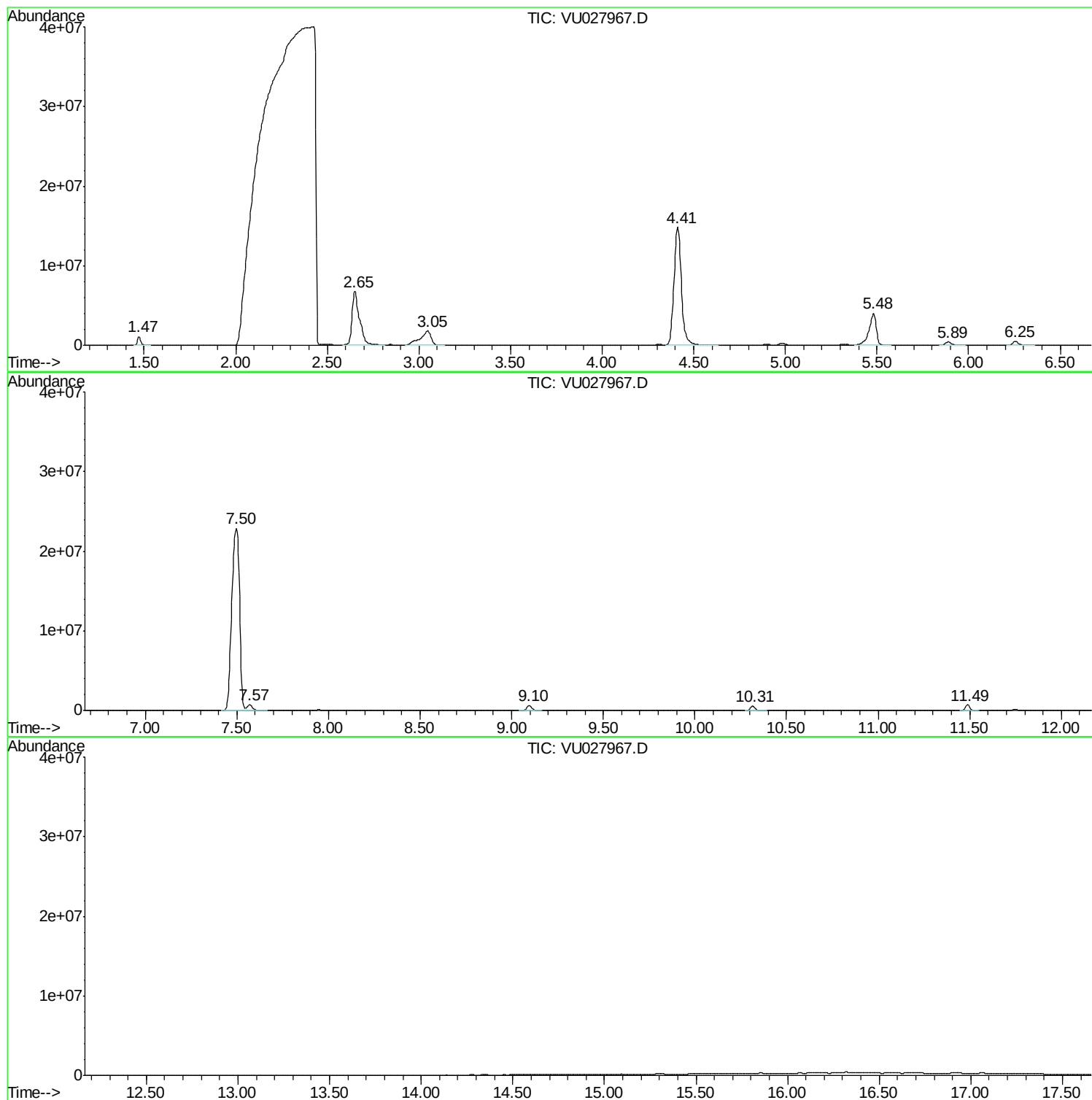
Sum of corrected areas: 153970900

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U110218W.M
Quant Title : SW846 8260

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



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

 Peak Number 1 2-Propanol, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
3.05	34.27 ug/l	7497020	Pentafluorobenzene	4.98	
Hit# of	5	Tentative ID	MW MolForm	CAS#	Qual
1	2-Propanol, 2-methyl-	74	C4H10O	000075-65-0	83
2	Propanoic acid, 2-hydroxy-2-methyl-	104	C4H8O3	000594-61-6	9
3	3-Pentanol	88	C5H12O	000584-02-1	9
4	Guanidine	59	CH5N3	000113-00-8	4
5	Formamide, N-methyl-	59	C2H5NO	000123-39-7	4

 Peak Number 2 1-Propanol, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.48	596.34 ug/l	10937800	1,4-Difluorobenzene	5.89	
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual	
1	1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	90
2	Pentane	72	C5H12	000109-66-0	72
3	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	38
4	2(3H)-Furanone, dihydro-4-methyl-	100	C5H8O2	001679-49-8	38
5	Butane, 2-methyl-	72	C5H12	000078-78-4	36

 Peak Number 3 1-Butanol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.25	61.25 ug/l	1123350	1,4-Difluorobenzene	5.89
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Butanol	74	C4H10O	000071-36-3 91
2	3,4-Dimethyldihydrofuran-2,5-dione	128	C6H8O3	007475-92-5 33
3	Isobutane	58	C4H10	000075-28-5 9
4	Diaziridine,3,3-dimethyl-	72	C3H8N2	004901-76-2 9
5	Formic acid, butyl ester	102	C5H10O2	000592-84-7 9

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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Propanol, 2-met...	3.05	34.3	ug/l	7497020	1	4.98	10937800	50.0
1-Propanol, 2-met...	5.48	596.3	ug/l	10937800	2	5.89	917068	50.0
1-Butanol	6.25	61.3	ug/l	1123350	2	5.89	917068	50.0