

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM011415\
 Data File : BM000193.D
 Acq On : 14 Jan 2015 18:28
 Operator : TP/IZ
 Sample : G1063-06DL 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 ClientSampleId :
 C0AN4DL

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % 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:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-BM122914.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.840	21	26	44	rVB	2699858	3721347	3.06%	2.402%
2	2.987	46	51	59	rVV	281021	427911	0.35%	0.276%
3	3.063	59	64	74	rVB	554924	698880	0.58%	0.451%
4	3.299	97	104	141	rVB2	62067715	121417088	100.00%	78.381%
5	4.151	244	249	257	rBV3	93677	149381	0.12%	0.096%
6	4.575	314	321	335	rBV	10868603	15368609	12.66%	9.921%
7	5.287	436	442	453	rVB	433926	640070	0.53%	0.413%
8	7.163	756	761	768	rVB	83132	129832	0.11%	0.084%
9	7.357	788	794	802	rVB2	92684	151474	0.12%	0.098%
10	7.828	868	874	882	rVB	519571	840283	0.69%	0.542%
11	8.986	1065	1071	1078	rBV	103690	172489	0.14%	0.111%
12	10.239	1276	1284	1290	rBV	109472	182772	0.15%	0.118%
13	10.622	1342	1349	1356	rVB	661923	1105381	0.91%	0.714%
14	13.868	1896	1901	1906	rBV	245629	332977	0.27%	0.215%
15	14.157	1944	1950	1957	rBV	241918	352610	0.29%	0.228%
16	14.462	1994	2002	2009	rBV	909703	1363702	1.12%	0.880%
17	15.457	2165	2171	2182	rVB	346525	504296	0.42%	0.326%
18	17.203	2462	2468	2479	rBV2	1116287	1691743	1.39%	1.092%
19	17.303	2479	2485	2495	rVB2	392189	563514	0.46%	0.364%
20	19.597	2869	2875	2884	rBV	482116	646504	0.53%	0.417%
21	21.391	3174	3180	3185	rBV	1630719	2020744	1.66%	1.304%
22	23.532	3538	3544	3553	rVB2	296096	588572	0.48%	0.380%
23	23.680	3562	3569	3578	rVB	937650	1835583	1.51%	1.185%

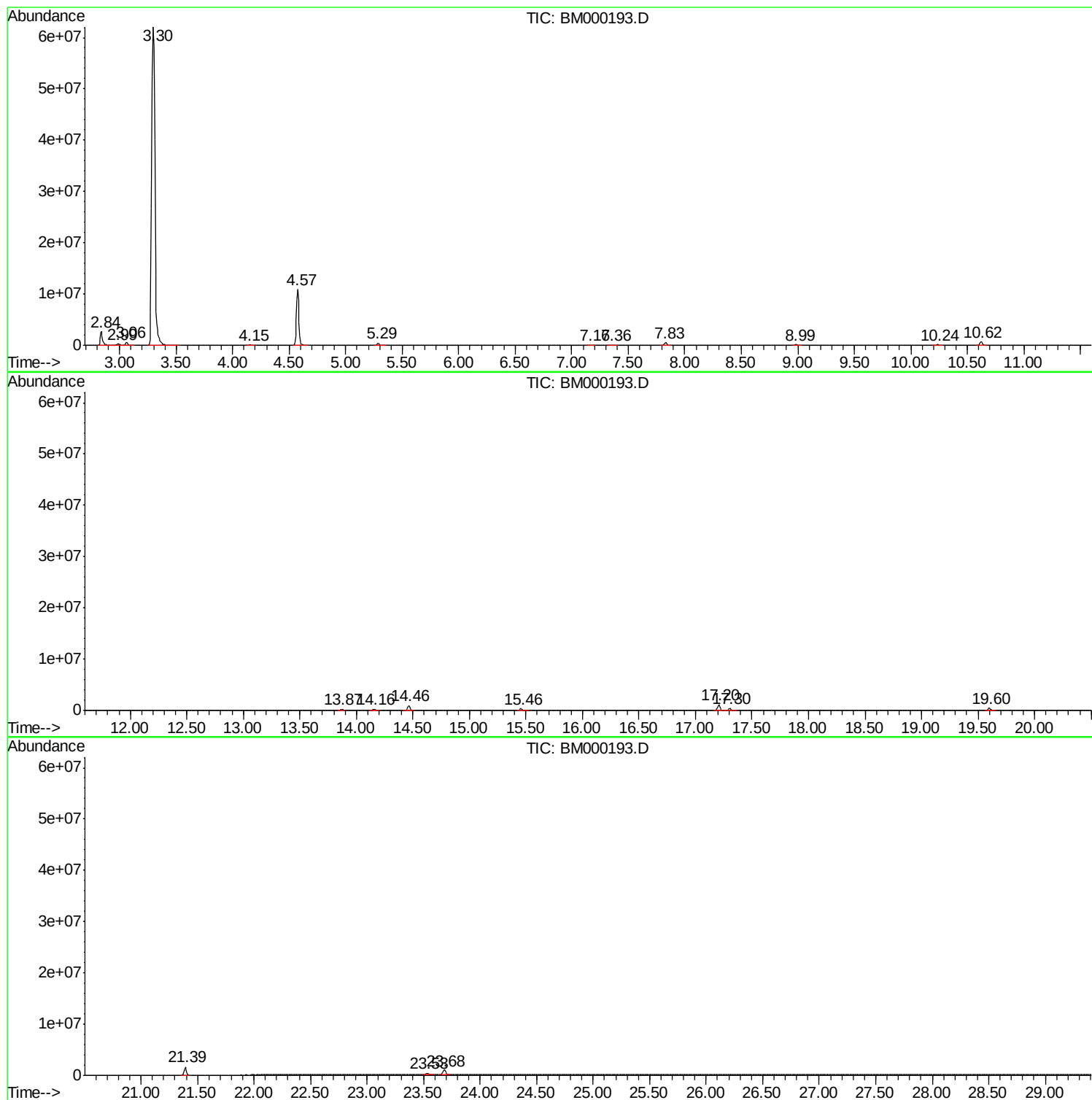
Sum of corrected areas: 154905762

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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-BM122914.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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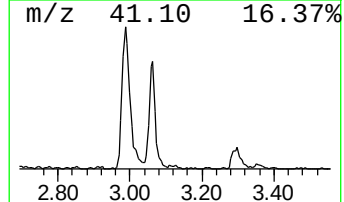
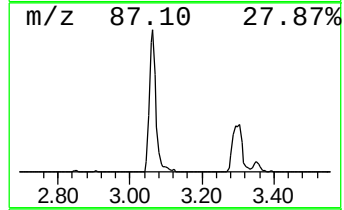
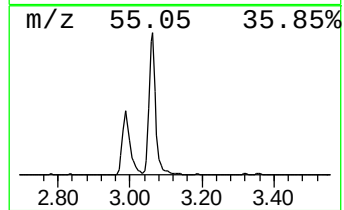
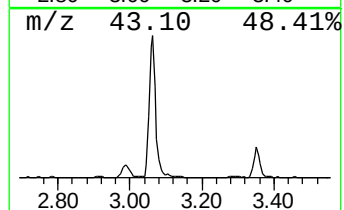
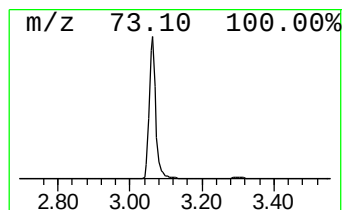
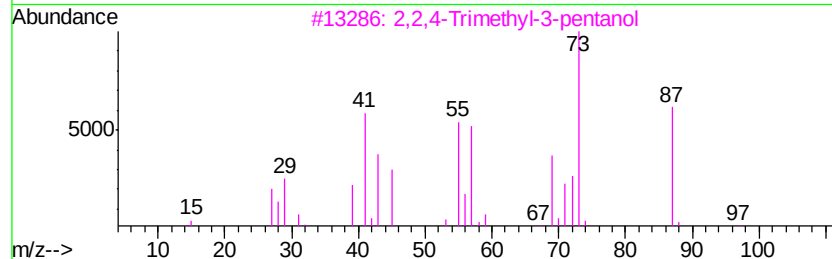
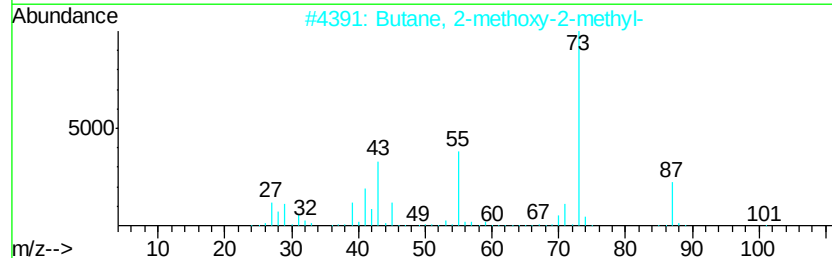
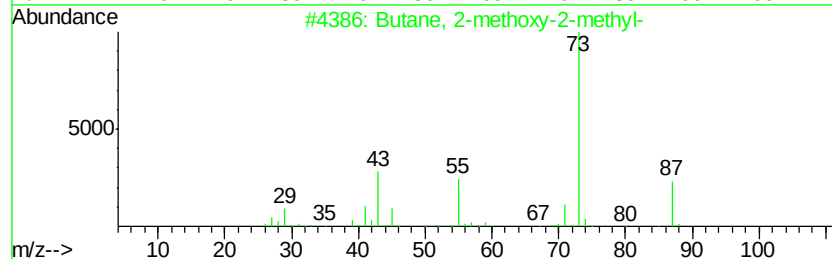
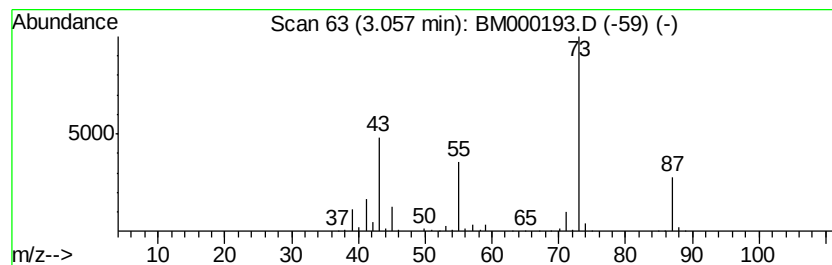
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.06	16.63 ng/ul	698880	1,4-Dichlorobenzene-d4	7.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	74
3		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4		Silane, tetramethyl-	88	C4H12Si	000075-76-3	23
5		2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	17



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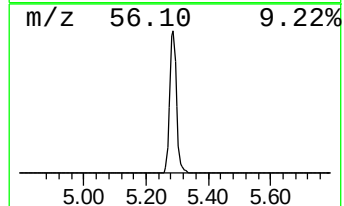
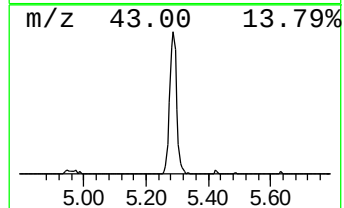
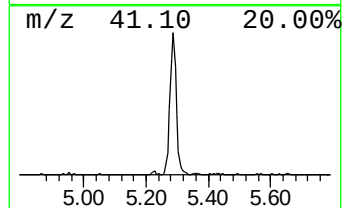
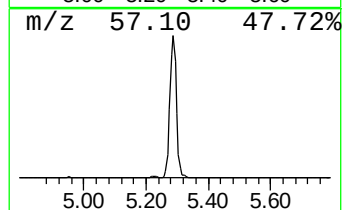
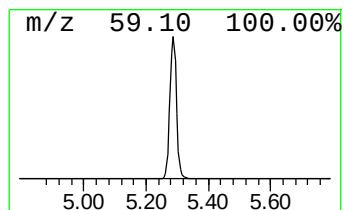
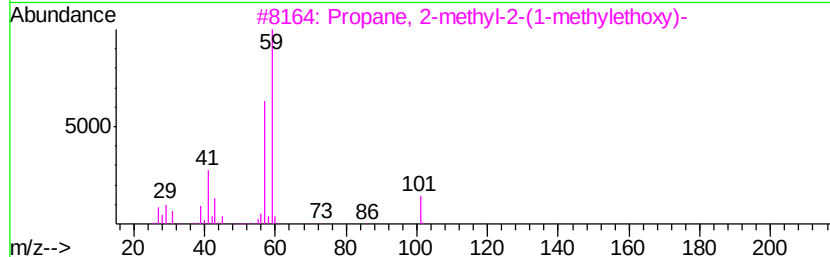
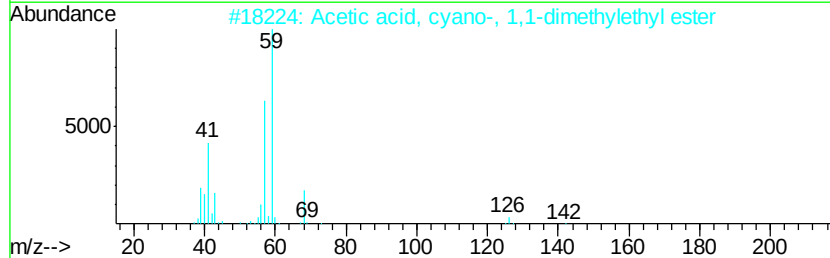
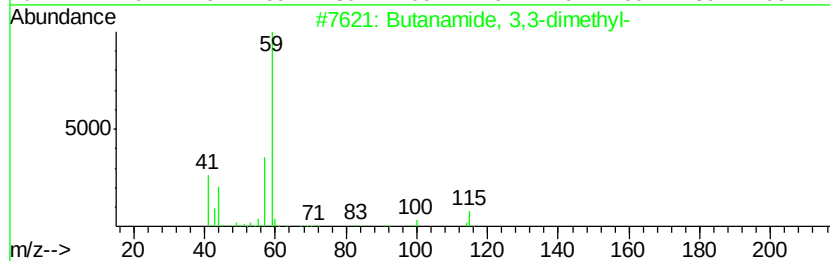
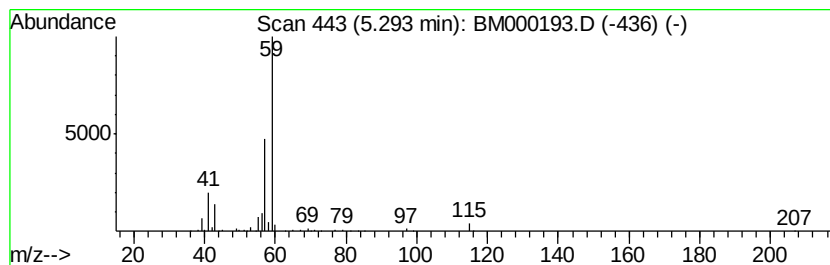
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Butanamide, 3,3-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.29	15.23 ng/ul	640070	1,4-Dichlorobenzene-d4	7.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butanamide, 3,3-dimethyl-	115	C6H13NO	000926-04-5	64
2		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	56
3		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	39
4		2-Propanol, 2-methyl-	74	C4H10O	000075-65-0	9
5		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	9



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	3.06	16.6	ng/ul	698880	1	7.83	840283	20.0
Butanamide, 3,3-d...	5.29	15.2	ng/ul	640070	1	7.83	840283	20.0