

Data Path : Z:\HPCHEM1\BNA M\DATA\BM011615\
 Data File : BM000234.D
 Acq On : 16 Jan 2015 15:37
 Operator : TP/IZ
 Sample : G1063-09RXDL 2X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 ClientSampleId :
 C0AN8RXDL

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	20	26	45	rVB	4581446	6544534	18.07%	7.305%
2	2.987	46	51	59	rVV	432296	677840	1.87%	0.757%
3	3.058	59	63	74	rVB	1376561	1686840	4.66%	1.883%
4	3.293	97	103	111	rBV	18363703	25041253	69.15%	27.951%
5	4.146	244	248	258	rVB4	24536	41846	0.12%	0.047%
6	4.575	314	321	340	rVB	25778892	36213663	100.00%	40.422%
7	4.957	378	386	392	rBV	44377	70510	0.19%	0.079%
8	5.281	436	441	448	rBV	65854	98704	0.27%	0.110%
9	6.987	725	731	738	rBV	69463	109960	0.30%	0.123%
10	7.157	754	760	769	rBV	212025	323059	0.89%	0.361%
11	7.357	788	794	804	rVB	240413	378262	1.04%	0.422%
12	7.828	868	874	882	rVB	510057	805579	2.22%	0.899%
13	8.522	986	992	999	rBV2	170962	275904	0.76%	0.308%
14	8.981	1063	1070	1078	rBV	271400	446260	1.23%	0.498%
15	9.704	1186	1193	1199	rBV	205326	342144	0.94%	0.382%
16	10.234	1276	1283	1293	rBV	311386	525805	1.45%	0.587%
17	10.616	1341	1348	1356	rBV	613043	1057908	2.92%	1.181%
18	13.869	1895	1901	1906	rBV	619407	869251	2.40%	0.970%
19	14.157	1943	1950	1957	rBV	648591	958814	2.65%	1.070%
20	14.463	1995	2002	2014	rBV	939251	1354857	3.74%	1.512%
21	14.651	2029	2034	2039	rBV5	26486	42298	0.12%	0.047%
22	15.451	2164	2170	2181	rVB	869540	1251634	3.46%	1.397%
23	15.569	2184	2190	2196	rBV	180474	267213	0.74%	0.298%
24	17.204	2462	2468	2476	rBV	1155141	1659759	4.58%	1.853%
25	17.304	2479	2485	2493	rBV	943114	1363117	3.76%	1.522%
26	18.092	2615	2619	2623	rBV3	41942	56445	0.16%	0.063%
27	19.598	2869	2875	2881	rBV	1156701	1547310	4.27%	1.727%
28	21.098	3126	3130	3134	rBV4	40702	61836	0.17%	0.069%
29	21.392	3175	3180	3186	rBV	1542323	1939509	5.36%	2.165%
30	22.586	3379	3383	3390	rVB	135418	195575	0.54%	0.218%
31	23.539	3538	3545	3552	rBV2	827953	1527119	4.22%	1.705%
32	23.686	3563	3570	3581	rVB2	942150	1853901	5.12%	2.069%

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Instrument :
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Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-BM122914.M
Title : SVOA CALIBRATION

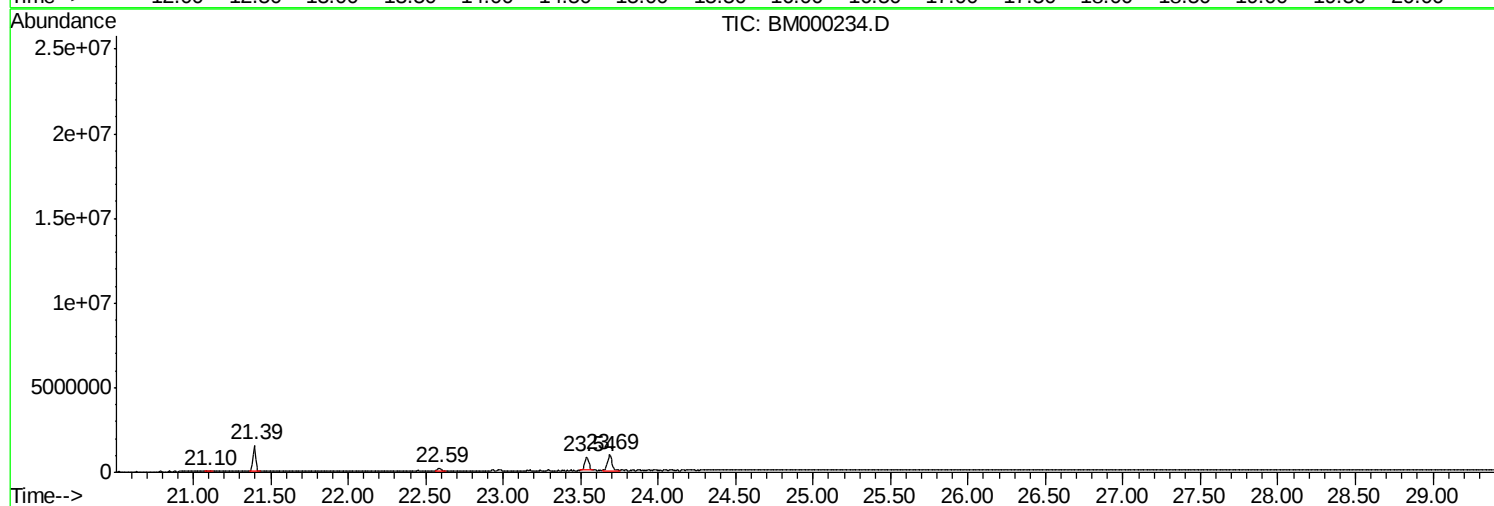
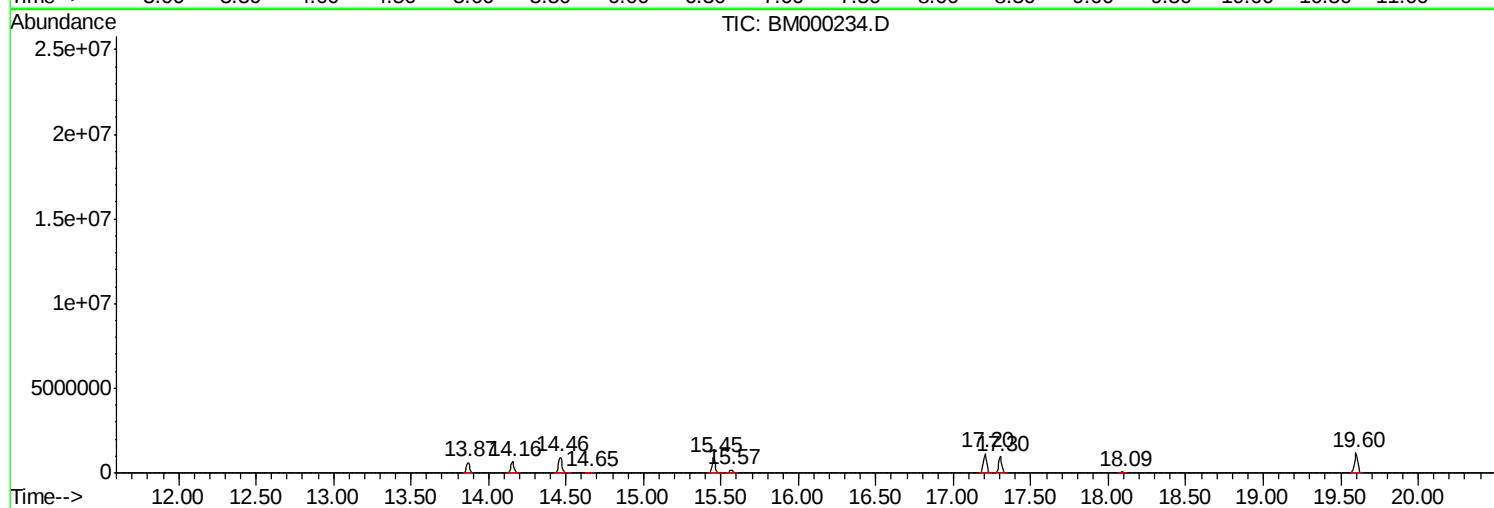
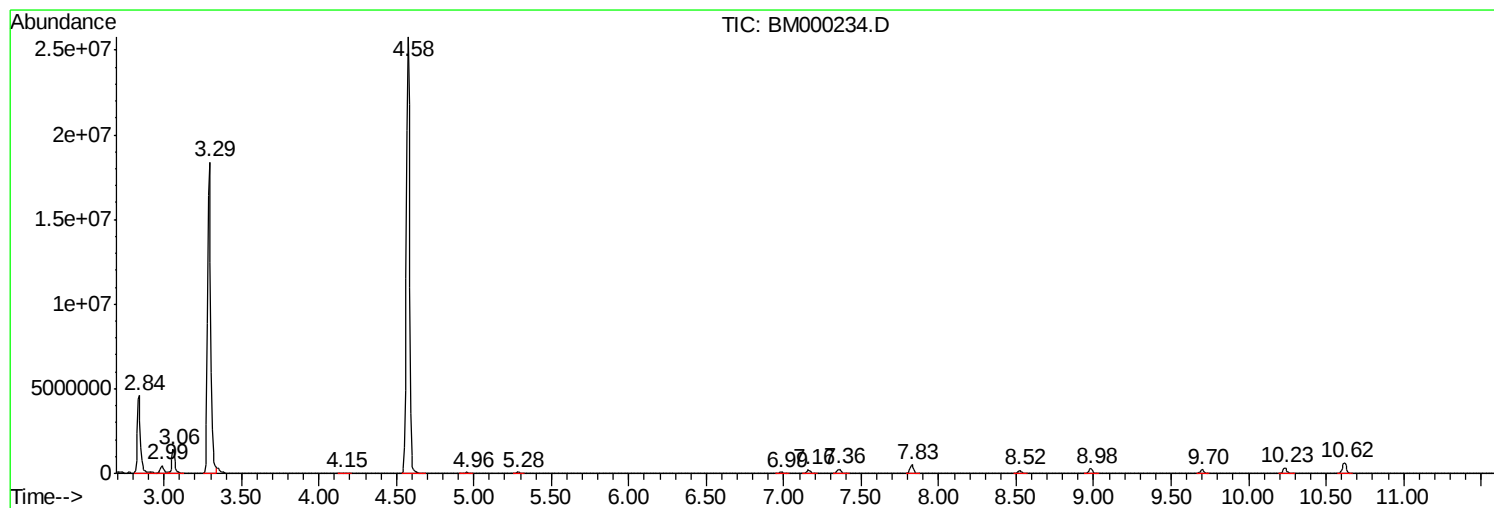
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM01.2-EPA-BM122914.M
Quant Title : SVOA CALIBRATION

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



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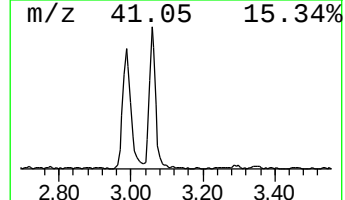
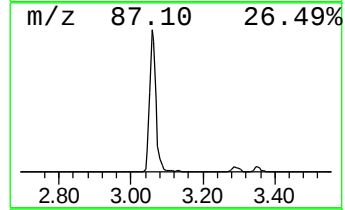
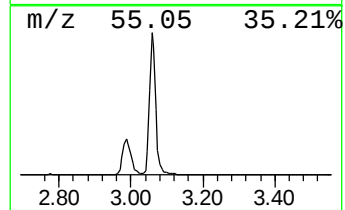
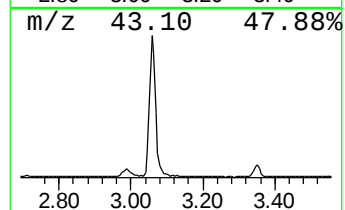
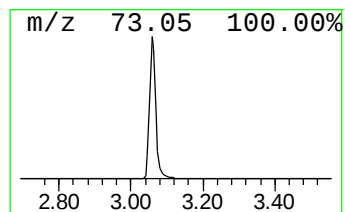
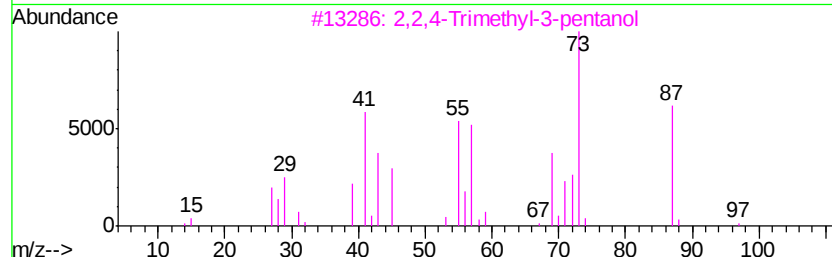
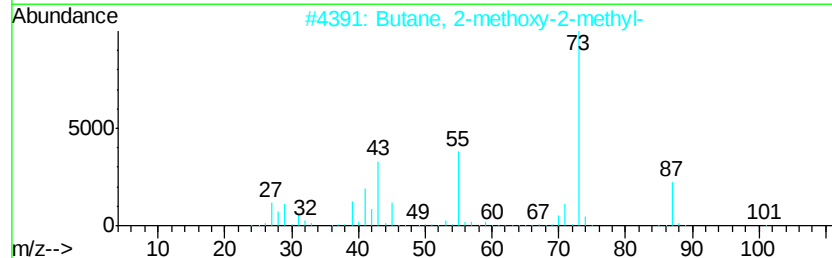
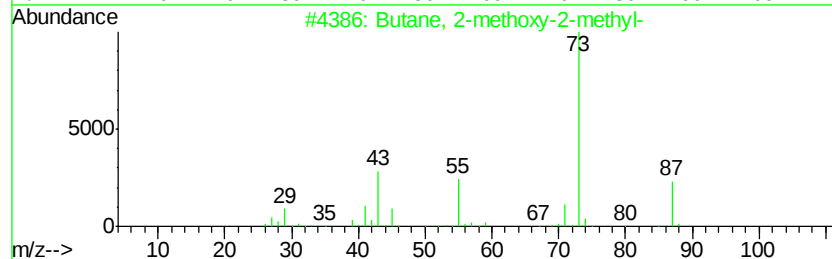
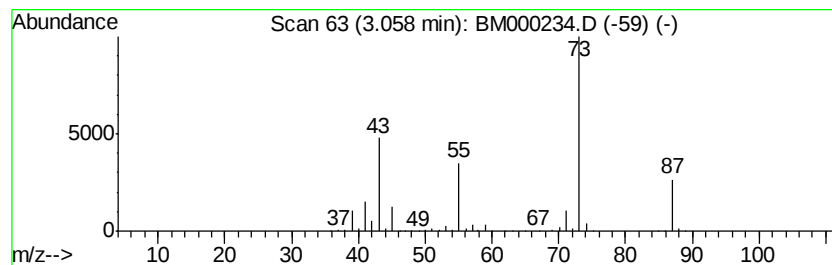
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM01.2-EPA-BM122914.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.06	41.88 ng/ul	1686840	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	64
3		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4		Silane, tetramethyl-	88	C4H12Si	000075-76-3	17
5		Boronic acid, ethyl-, dimethyl e...	102	C4H11BO2	007318-82-3	10



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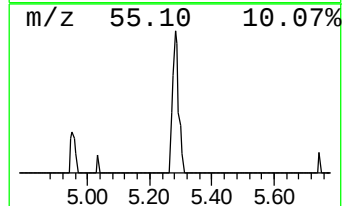
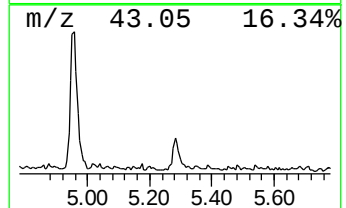
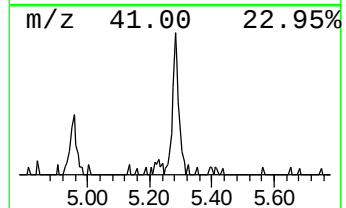
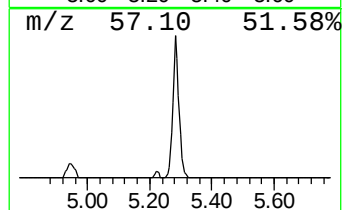
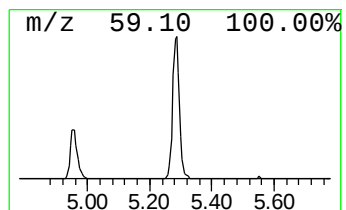
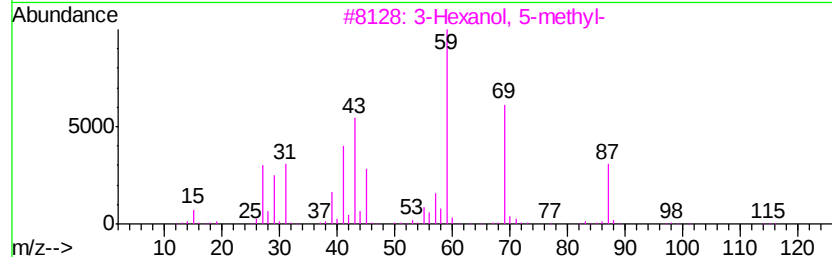
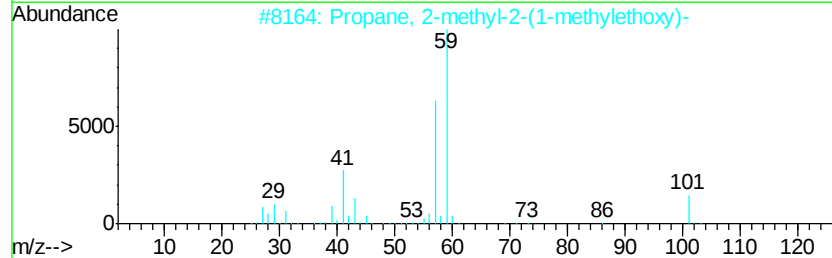
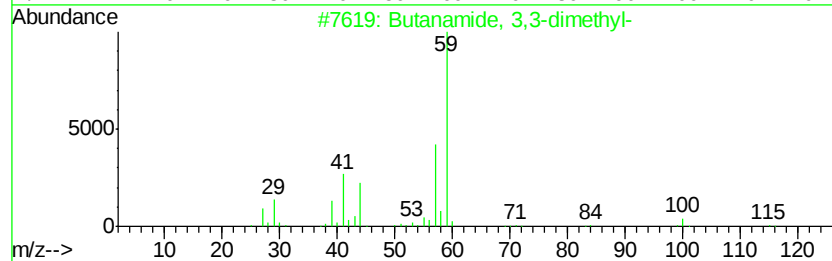
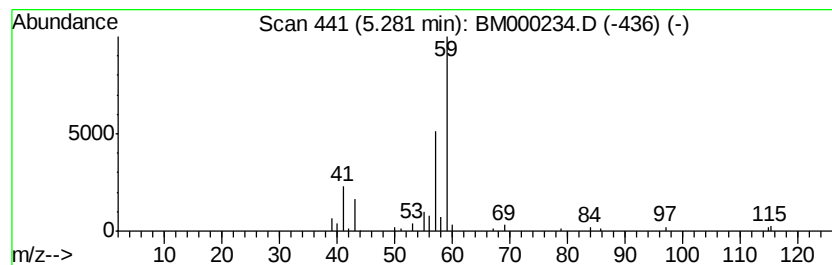
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Peak Number 5 Butanamide, 3,3-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.28	2.45 ng/ul	98704	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	56
2		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	56
3		3-Hexanol, 5-methyl-	116	C7H16O	000623-55-2	40
4		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	39
5		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	39



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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	41.9	ng/ul	1686840	1	7.83	805579	20.0
Butanamide, 3,3-d...	5.28	2.5	ng/ul	98704	1	7.83	805579	20.0