

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM070615\
 Data File : BM001833.D
 Acq On : 06 Jul 2015 12:58
 Operator : TP/UM
 Sample : PB84338BL
 Misc : SBLK82
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK82

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\SOM02.2-EPA-BM061515.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	2.805	16	20	27	rVB	21826	32455	1.96%	0.178%
2	2.881	29	33	44	rVB	105587	127528	7.71%	0.701%
3	3.140	74	77	86	rVB	16333	19510	1.18%	0.107%
4	3.693	168	171	173	rBV	4888	5146	0.31%	0.028%
5	6.828	698	704	712	rBV	264575	395458	23.90%	2.175%
6	6.999	727	733	741	rBB	196312	292746	17.69%	1.610%
7	7.193	760	766	774	rBB	271757	410107	24.78%	2.255%
8	7.663	839	846	853	rBB	273961	424258	25.64%	2.333%
9	8.363	959	965	973	rBV	315356	500353	30.24%	2.752%
10	8.822	1036	1043	1051	rBB	237938	381378	23.05%	2.097%
11	9.540	1158	1165	1172	rVB	203549	323547	19.55%	1.779%
12	10.069	1248	1255	1266	rVB	328383	528902	31.96%	2.909%
13	10.451	1313	1320	1328	rBB	419741	685777	41.44%	3.771%
14	10.593	1337	1344	1352	rBV	320862	517284	31.26%	2.845%
15	13.728	1871	1877	1883	rBV	545139	723240	43.71%	3.977%
16	14.010	1918	1925	1932	rBV	619897	920082	55.60%	5.060%
17	14.316	1970	1977	1984	rBV2	686522	1035334	62.57%	5.694%
18	14.516	2006	2011	2021	rBV	237905	335335	20.27%	1.844%
19	15.310	2140	2146	2154	rVB	855148	1155535	69.83%	6.355%
20	15.433	2161	2167	2174	rBV	219620	290573	17.56%	1.598%
21	16.845	2404	2407	2411	rBV	2595	3335	0.20%	0.018%
22	17.063	2438	2444	2450	rBV2	943839	1303635	78.78%	7.169%
23	17.163	2454	2461	2468	rBV	916401	1289490	77.93%	7.091%
24	19.351	2829	2833	2838	rBV	5606	7626	0.46%	0.042%
25	19.463	2846	2852	2861	rVB	1160961	1526248	92.24%	8.393%
26	20.486	3025	3026	3029	rBV3	1787	2025	0.12%	0.011%
27	20.827	3083	3084	3088	rBV4	2234	2676	0.16%	0.015%
28	20.927	3099	3101	3103	rBV2	1858	1794	0.11%	0.010%
29	21.051	3119	3122	3126	rVB2	12340	12701	0.77%	0.070%
30	21.257	3152	3157	3163	rBV	1352119	1654686	100.00%	9.100%
31	23.339	3504	3511	3521	rVB	943109	1631808	98.62%	8.974%
32	23.486	3528	3536	3545	rVB	884765	1643574	99.33%	9.038%

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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

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Peak separation: 5

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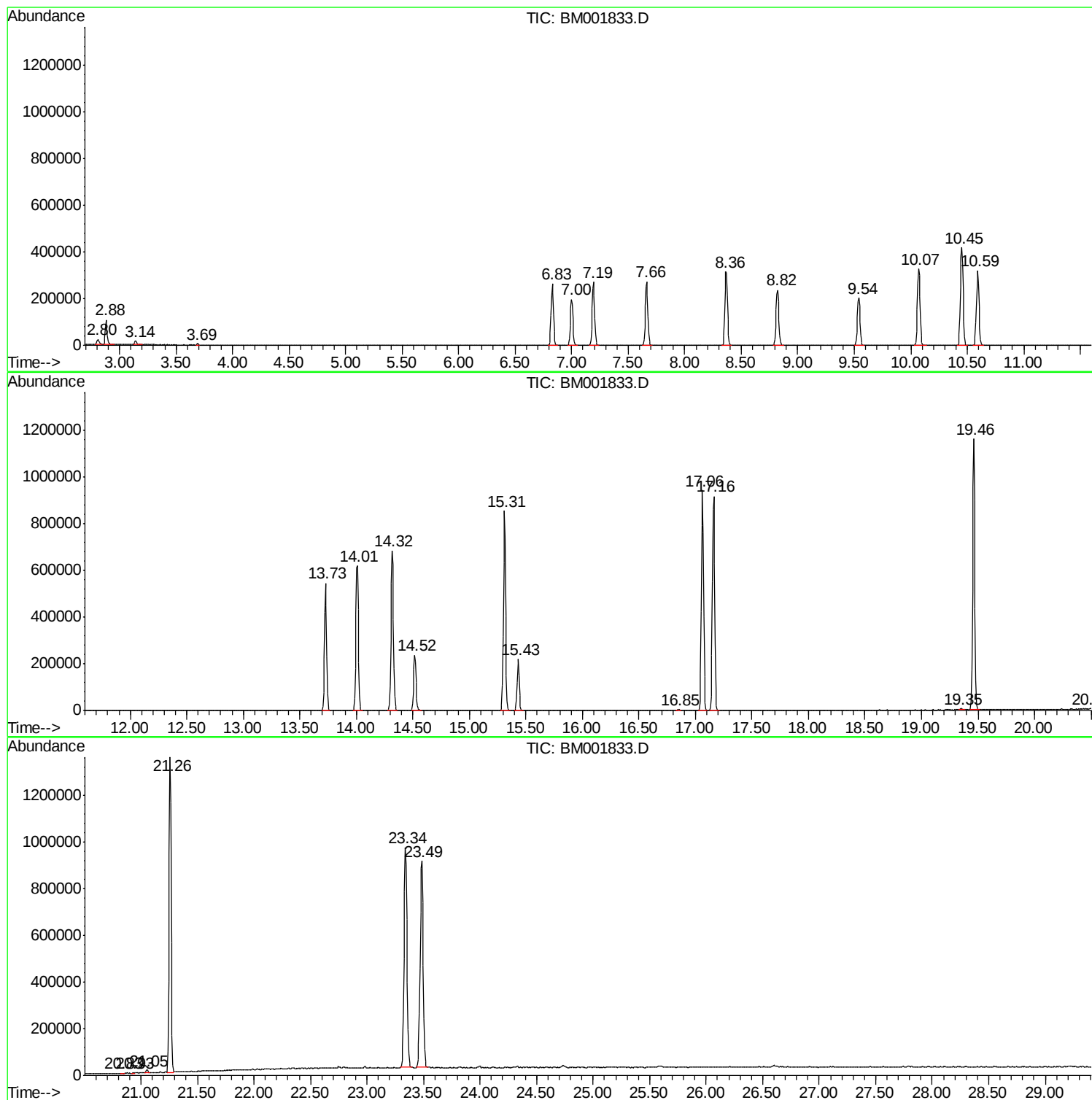
Sum of corrected areas: 18184146

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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061515.M
Quant Title : SVOA CALIBRATION

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



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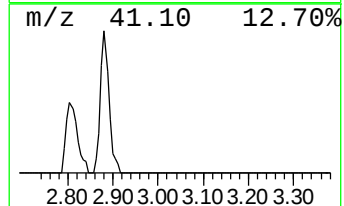
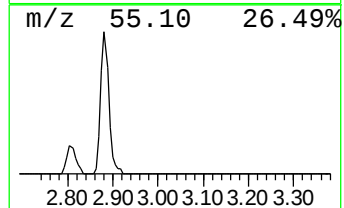
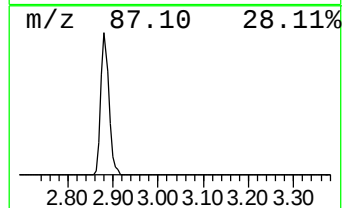
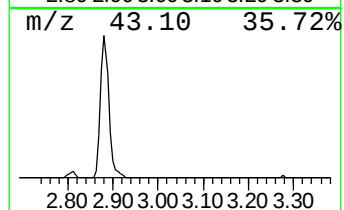
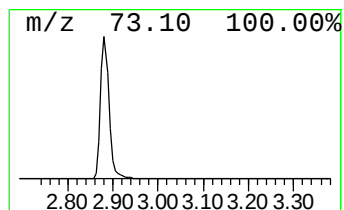
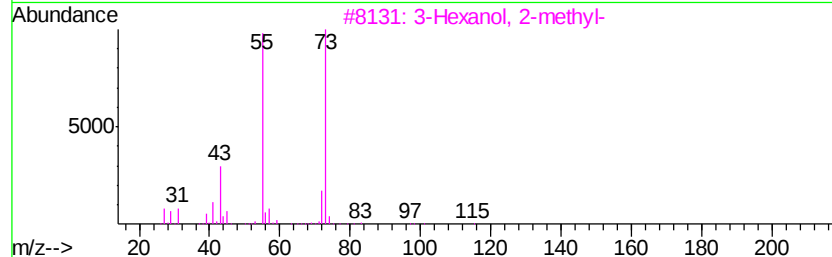
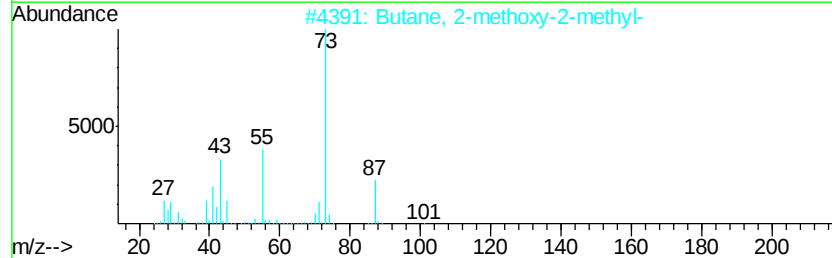
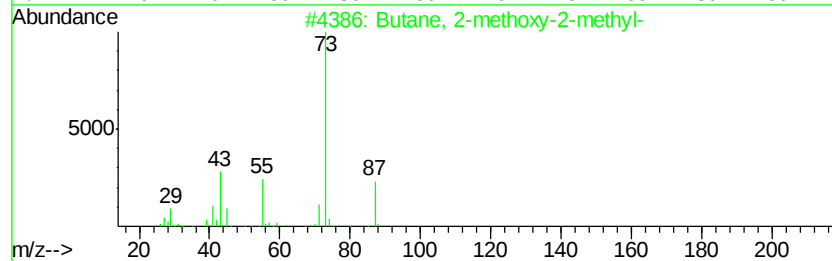
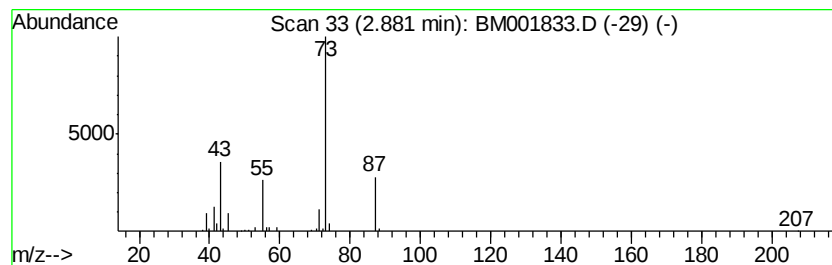
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.88	6.01 ng/ul	127528	1,4-Dichlorobenzene-d4	7.66

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
3			3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	12
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5			N-Ethylformamide	73	C3H7NO	000627-45-2	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...	2.88	6.0	ng/ul	127528	1	7.66	424258	20.0