

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062819\
 Data File : BM021250.D
 Acq On : 29 Jun 2019 02:33
 Operator : HP/JU
 Sample : K3381-12DL 20X
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BFCM6DL

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:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.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	3.016	3	5	11	rVB	162037	169944	7.31%	1.303%
2	3.310	51	55	67	rVB	179354	236999	10.19%	1.817%
3	7.104	696	700	707	rVB	25193	41005	1.76%	0.314%
4	7.293	728	732	738	rVB2	31084	48638	2.09%	0.373%
5	7.757	805	811	821	rVB	602361	970115	41.71%	7.439%
6	8.475	928	933	940	rBV3	16320	26160	1.12%	0.201%
7	8.928	1005	1010	1017	rBV2	34623	60181	2.59%	0.462%
8	9.287	1069	1071	1074	rVB3	3451	3626	0.16%	0.028%
9	9.645	1126	1132	1138	rBV2	22179	41303	1.78%	0.317%
10	9.898	1171	1175	1177	rBV3	2398	2431	0.10%	0.019%
11	10.192	1219	1225	1235	rBV2	23519	58474	2.51%	0.448%
12	10.545	1275	1285	1296	rBV	852878	1436614	61.77%	11.017%
13	11.181	1391	1393	1396	rVB3	2093	2473	0.11%	0.019%
14	11.204	1396	1397	1401	rBV2	1973	2666	0.11%	0.020%
15	11.775	1490	1494	1500	rVB2	3804	5314	0.23%	0.041%
16	12.163	1558	1560	1567	rBV2	960	2447	0.11%	0.019%
17	13.769	1829	1833	1834	rBV	2650	2344	0.10%	0.018%
18	13.816	1837	1841	1855	rVB	93215	144984	6.23%	1.112%
19	14.092	1882	1888	1896	rBV	105409	162966	7.01%	1.250%
20	14.398	1934	1940	1954	rBV2	1369419	2084889	89.65%	15.988%
21	15.351	2099	2102	2104	rBV2	2112	2576	0.11%	0.020%
22	15.398	2104	2110	2119	rVV	149183	229323	9.86%	1.759%
23	15.539	2130	2134	2136	rBV	3760	3661	0.16%	0.028%
24	16.127	2231	2234	2238	rBV4	1591	3101	0.13%	0.024%
25	16.239	2250	2253	2254	rBV2	3008	3107	0.13%	0.024%
26	16.357	2270	2273	2277	rVB4	3899	3893	0.17%	0.030%
27	16.686	2328	2329	2334	rVB2	2494	2588	0.11%	0.020%
28	16.727	2334	2336	2339	rBV4	2226	2680	0.12%	0.021%
29	16.851	2355	2357	2360	rVB3	2440	2352	0.10%	0.018%
30	16.892	2360	2364	2366	rBV3	1871	2813	0.12%	0.022%
31	16.939	2368	2372	2375	rVV2	4204	7207	0.31%	0.055%
32	17.151	2401	2408	2420	rBV	1591707	2325623	100.00%	17.834%
33	17.251	2420	2425	2433	rVB	131973	205416	8.83%	1.575%
34	17.598	2481	2484	2488	rVB4	1936	2490	0.11%	0.019%

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Title : SVOA CALIBRATION

35	17.909	2535	2537	2539	rBV3	2258	2509	0.11%	0.019%
36	19.545	2810	2815	2823	rBV	168381	243134	10.45%	1.864%
37	21.333	3114	3119	3124	rBV	1753481	2262216	97.27%	17.348%
38	23.597	3498	3504	3513	rVB	1187616	2232005	95.97%	17.116%

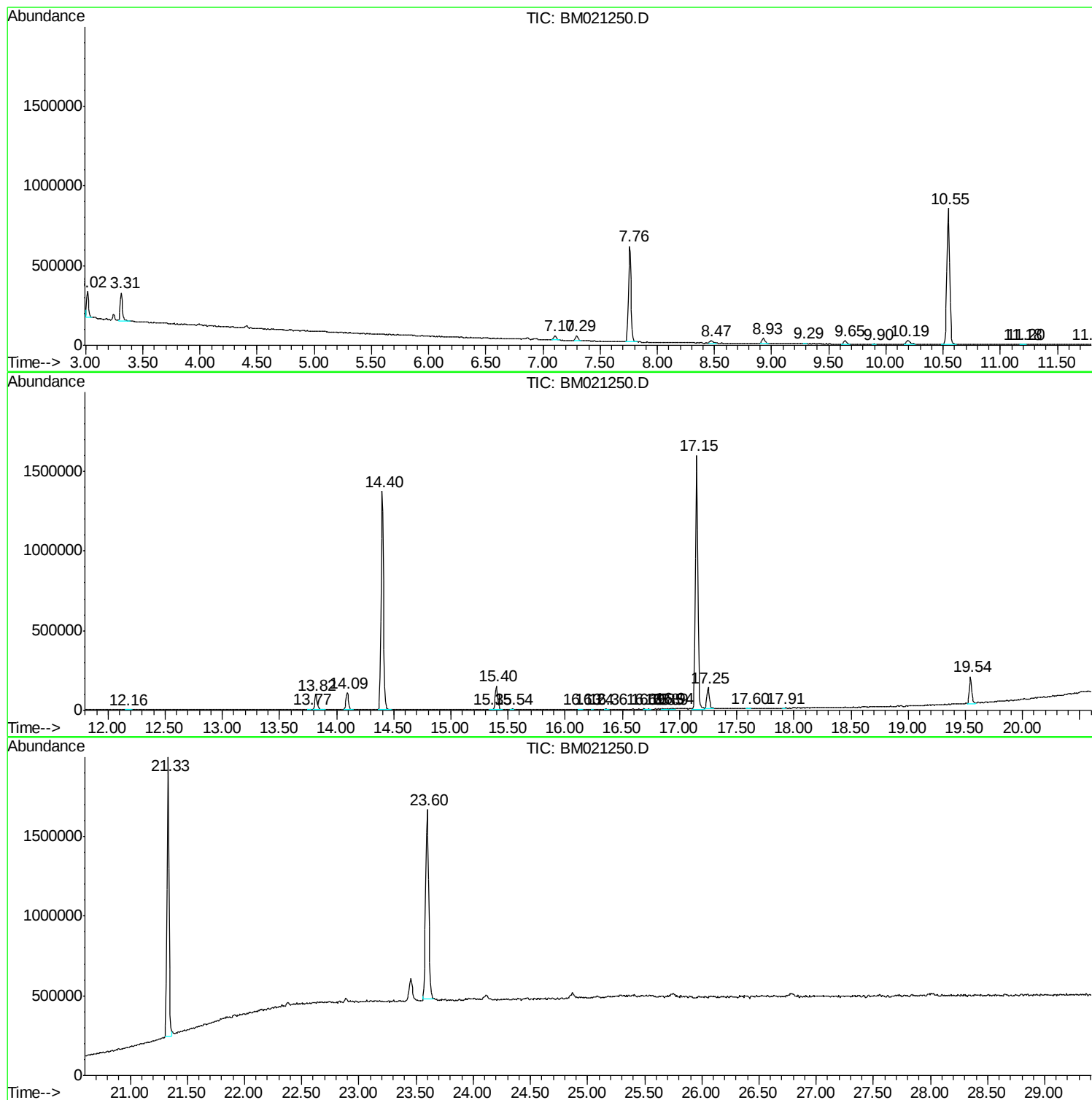
Sum of corrected areas: 13040267

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
Quant Title : SVOA CALIBRATION

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



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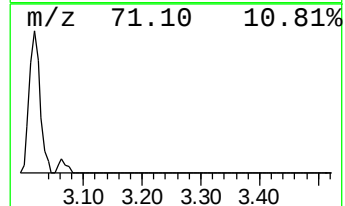
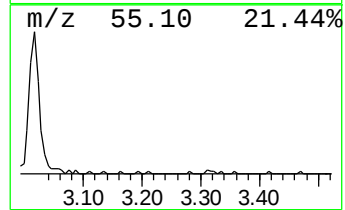
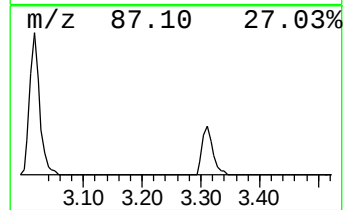
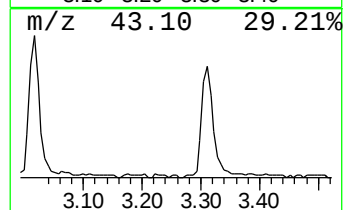
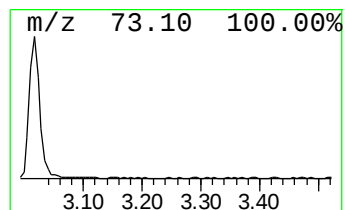
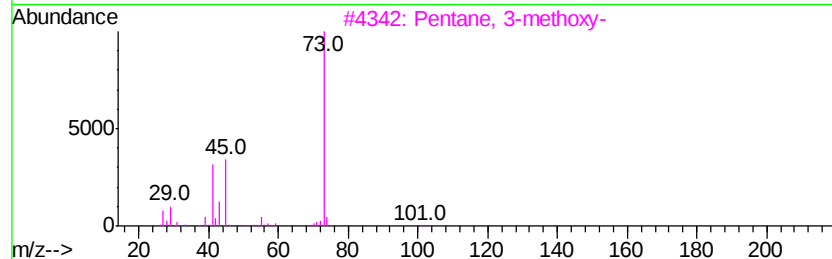
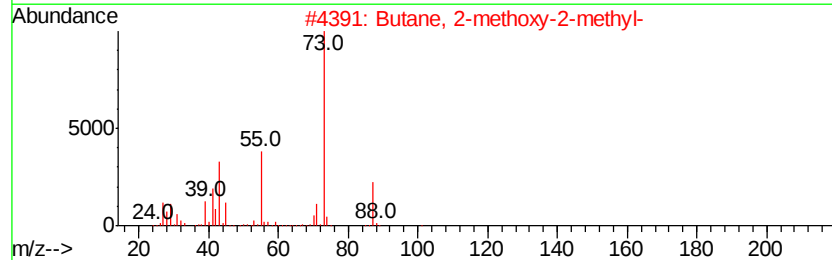
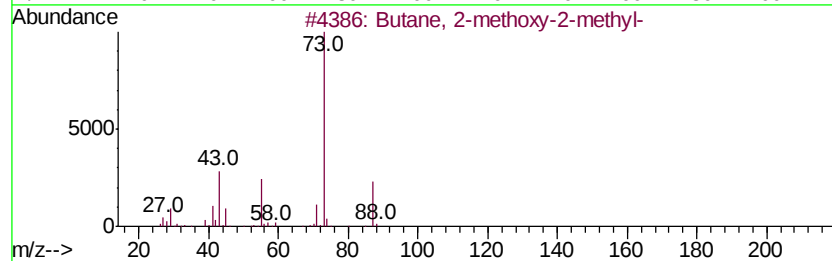
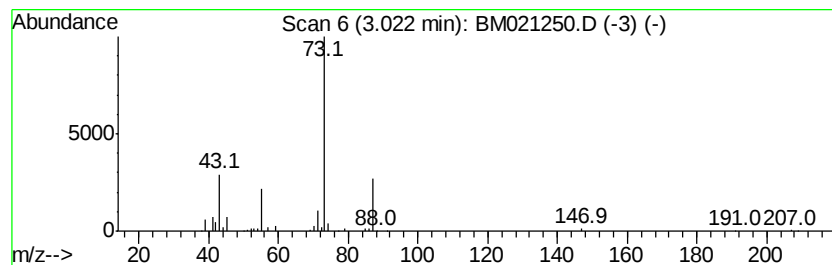
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061419MA.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.02	3.50 ng/ul	169944	1,4-Dichlorobenzene-d4	7.76

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4			N-Ethylformamide	73	C3H7NO	000627-45-2	9
5			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9



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Quant Title : SVOA CALIBRATION

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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...	3.02	3.5	ng/ul	169944	1	7.76	970115	20.0