

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF091219\
 Data File : BF116989.D
 Acq On : 12 Sep 2019 13:24
 Operator : HP/JU
 Sample : PB123013BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB123013BL

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.163	7	13	22	rVB	77355	114142	3.73%	0.481%
2	5.457	568	573	576	rBV	2438516	2363355	77.32%	9.953%
3	6.463	738	744	747	rBV	2343703	2529758	82.77%	10.653%
4	6.598	762	767	770	rBV	2756924	2641092	86.41%	11.122%
5	6.828	802	806	821	rVB	664610	634515	20.76%	2.672%
6	6.981	828	832	836	rBV	2125375	2020694	66.11%	8.510%
7	7.387	895	901	904	rBV	1639655	1568640	51.32%	6.606%
8	8.104	1018	1023	1039	rBV	861400	840584	27.50%	3.540%
9	9.181	1200	1206	1209	rBV	3179147	3056505	100.00%	12.872%
10	9.851	1316	1320	1335	rBV	974437	1004345	32.86%	4.230%
11	10.639	1450	1454	1467	rBV	1632022	1799833	58.89%	7.579%
12	11.333	1568	1572	1594	rBV	924784	977384	31.98%	4.116%
13	12.922	1837	1842	1845	rBV	3007608	2922010	95.60%	12.305%
14	13.969	2016	2020	2029	rBV	709501	711198	23.27%	2.995%
15	15.416	2260	2266	2277	rBV	377544	562037	18.39%	2.367%

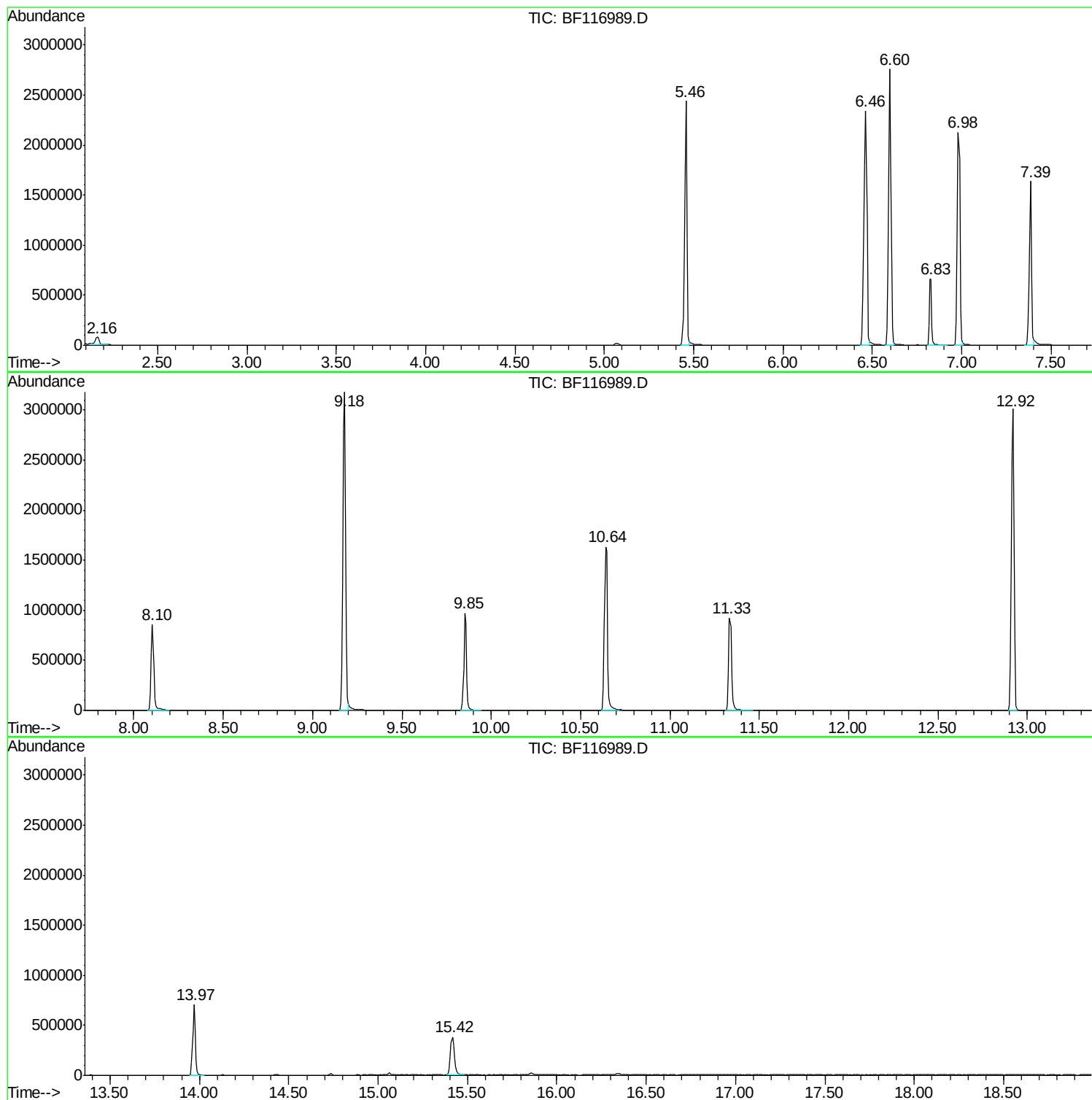
Sum of corrected areas: 23746092

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



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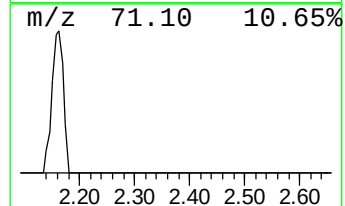
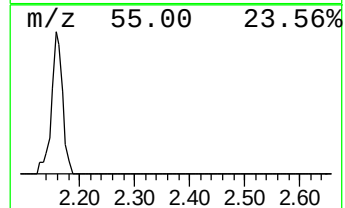
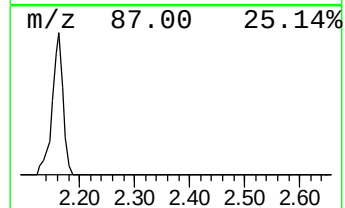
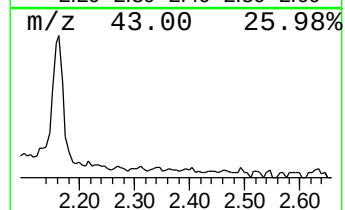
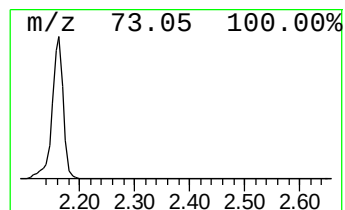
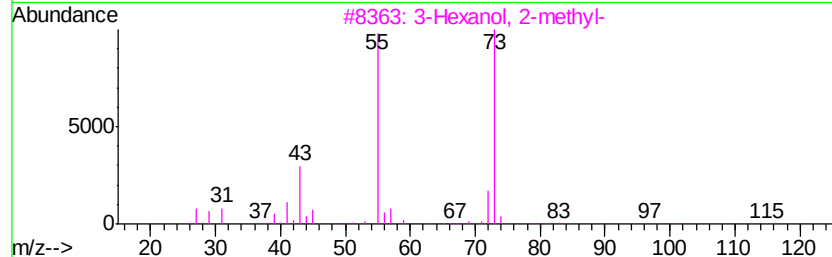
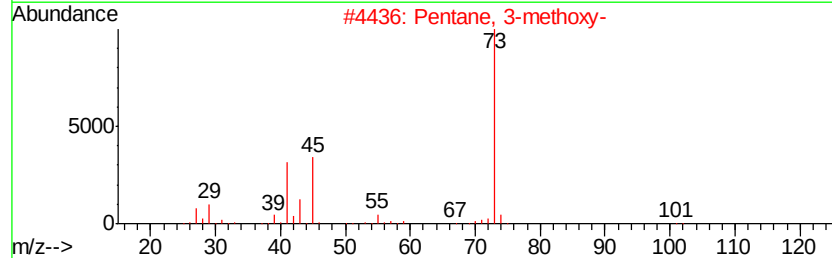
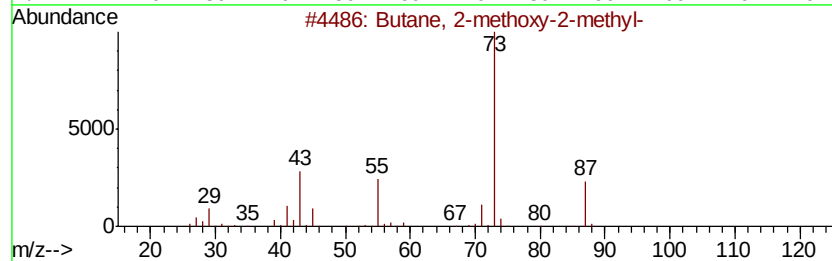
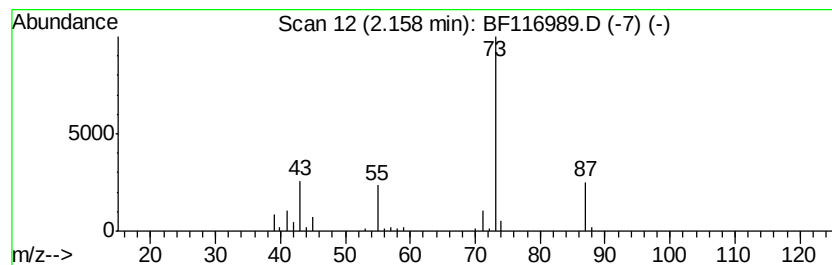
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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.16	3.60 ng	114142	1,4-Dichlorobenzene-d4	6.83

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	86
2			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
3			3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	17
4			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
5			N-Ethylformamide	73	C3H7NO	000627-45-2	9



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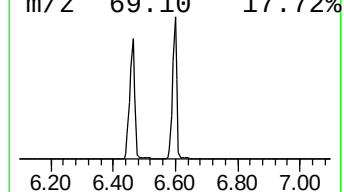
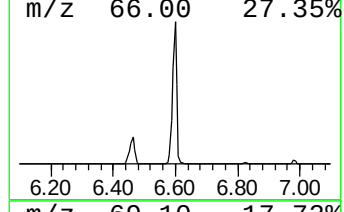
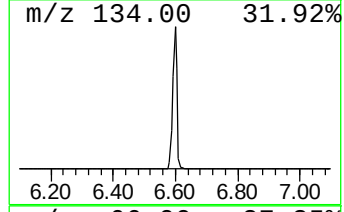
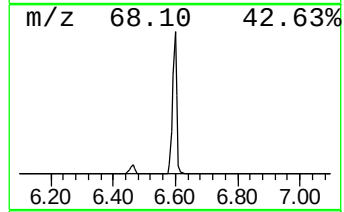
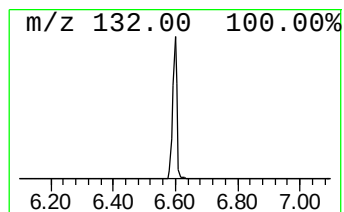
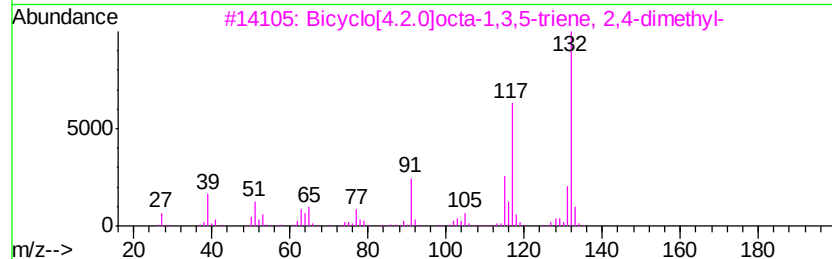
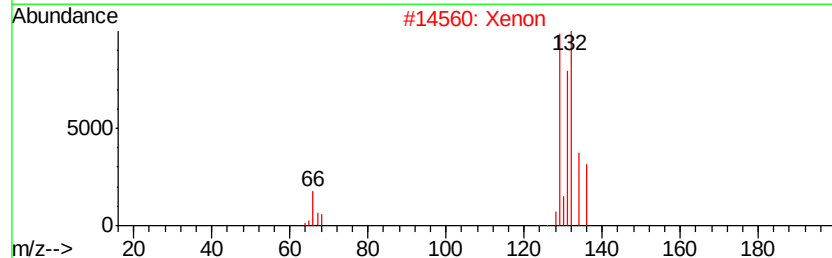
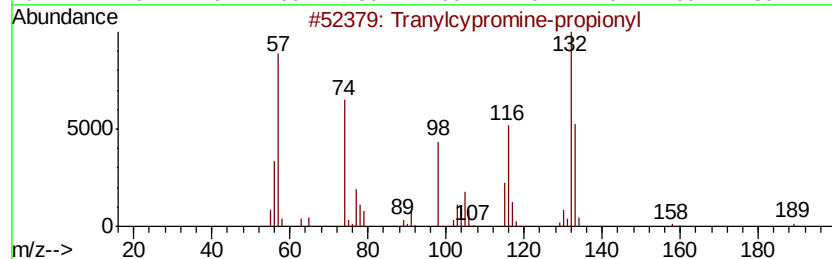
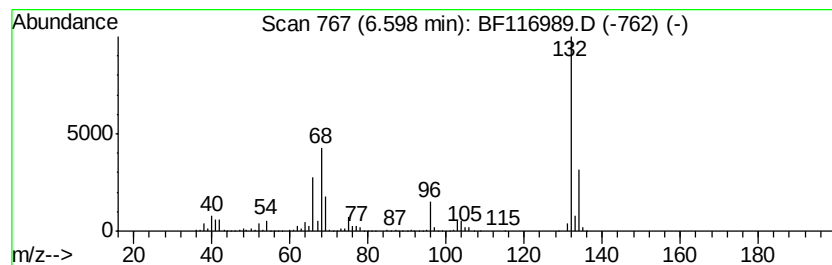
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Peak Number 2 unknown6.60 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.60	83.25 ng	2641090	1,4-Dichlorobenzene-d4	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tranlylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
2		Xenon	132	Xe	007440-63-3	9
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.16	3.6	ng	114142	1	6.83	634515	20.0
unknown6.60	6.60	83.3	ng	2641090	1	6.83	634515	20.0