

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM081718\
 Data File : BM016477.D
 Acq On : 18 Aug 2018 17:08
 Operator : SJ/JU
 Sample : PB112126BL
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB112126BL

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % 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\8270-BM081718.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	5.110	306	310	322	rBV	48571	78694	2.76%	0.870%
2	5.587	386	391	404	rBV	344440	553334	19.43%	6.116%
3	7.222	664	669	687	rBV	396947	631797	22.18%	6.983%
4	7.581	724	730	742	rBV	423470	692514	24.31%	7.654%
5	8.057	806	811	820	rBB	48270	81660	2.87%	0.903%
6	8.369	858	864	877	rBB	313279	495575	17.40%	5.477%
7	9.239	1006	1012	1027	rBV	206913	365650	12.84%	4.041%
8	10.863	1283	1288	1300	rBV	57087	98521	3.46%	1.089%
9	13.310	1698	1704	1720	rBB	736409	993828	34.89%	10.984%
10	14.692	1934	1939	1948	rBB2	94998	145702	5.11%	1.610%
11	16.186	2187	2193	2204	rBV	936699	1216374	42.70%	13.444%
12	17.433	2400	2405	2416	rBV	142467	209577	7.36%	2.316%
13	20.027	2840	2846	2851	rBV	2575427	2848547	100.00%	31.483%
14	21.580	3105	3110	3114	rBV	245842	314431	11.04%	3.475%
15	23.880	3496	3501	3509	rVB2	171734	321572	11.29%	3.554%

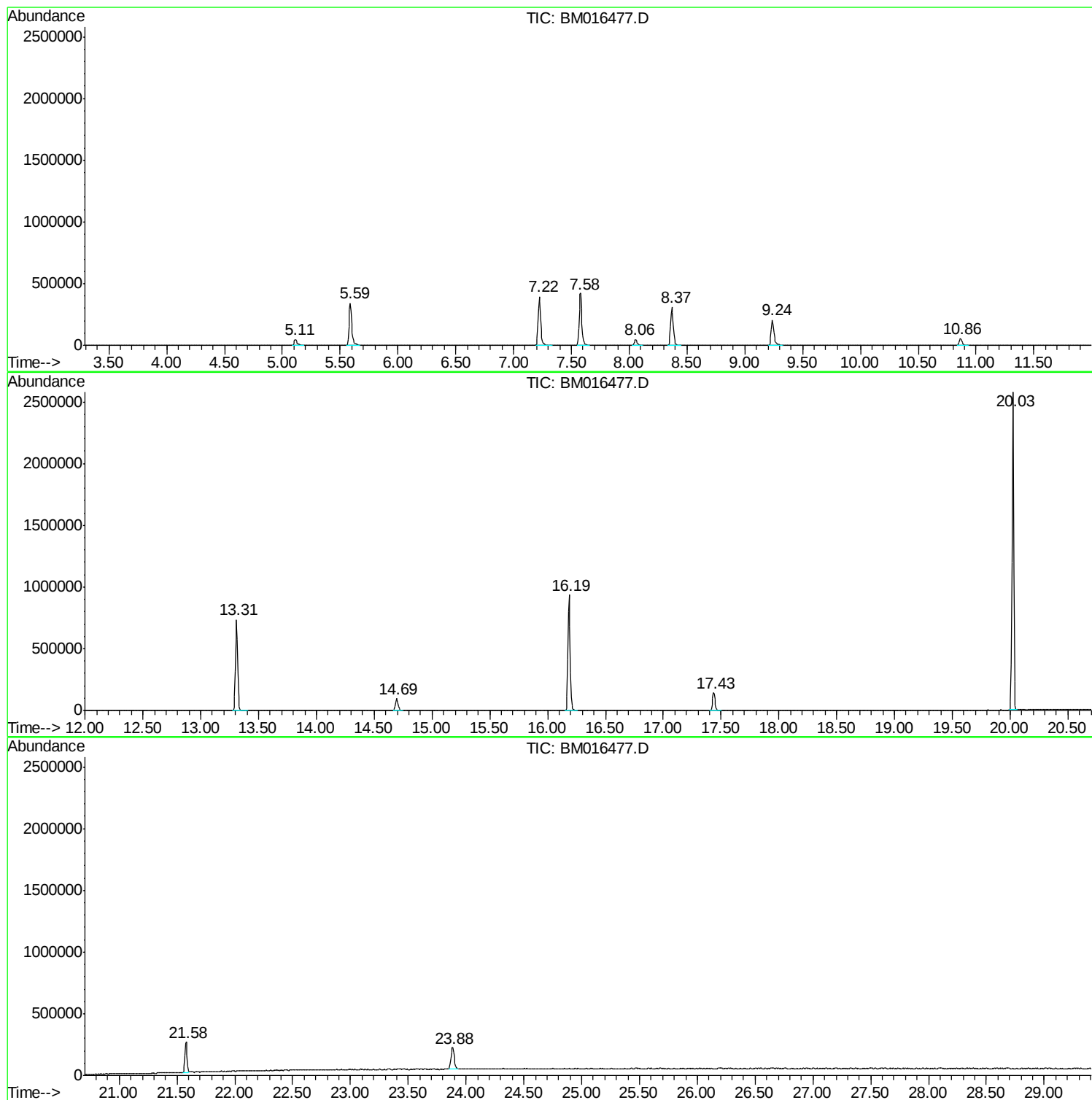
Sum of corrected areas: 9047776

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



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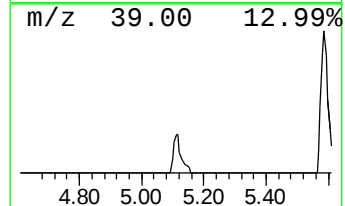
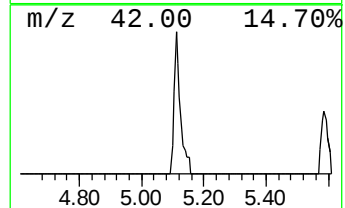
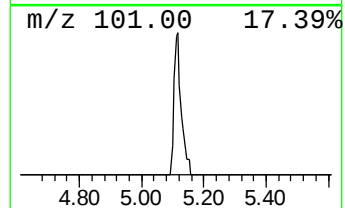
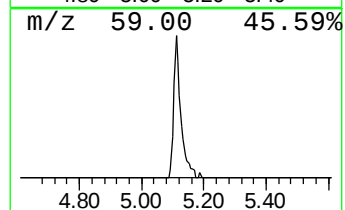
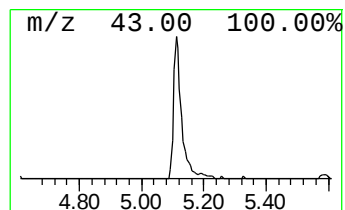
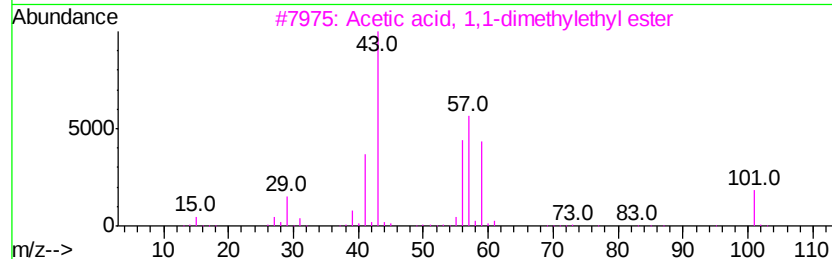
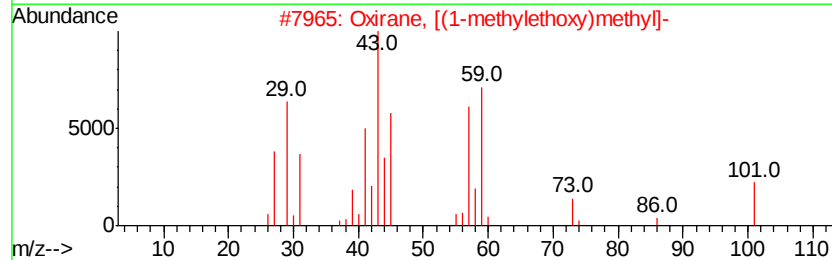
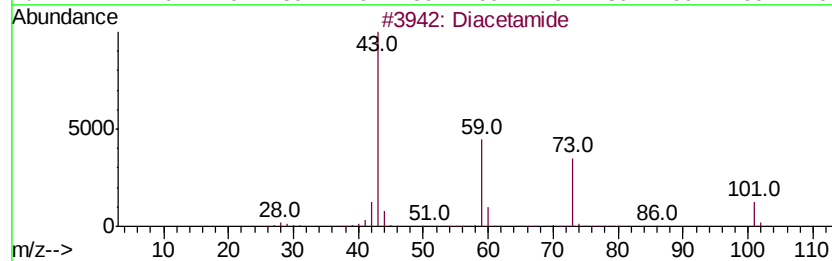
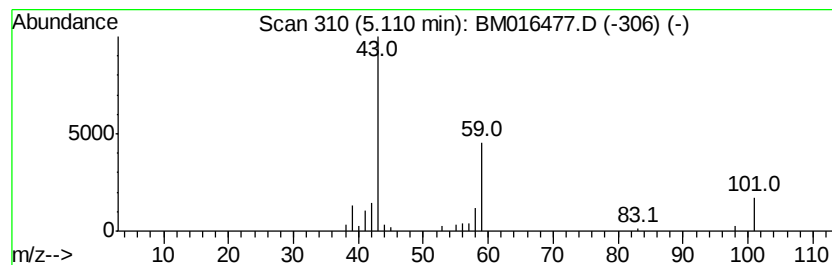
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown5.11 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.11	19.27 ng	78694	1,4-Dichlorobenzene-d4	8.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Diacetamide	101	C4H7NO2	000625-77-4	45
2		Oxirane, [(1-methylethoxy)methyl]-	116	C6H12O2	004016-14-2	39
3		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	36
4		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	33
5		Butane	58	C4H10	000106-97-8	27



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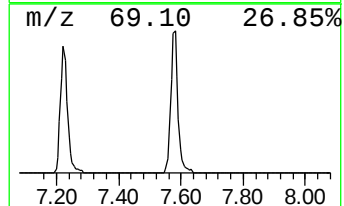
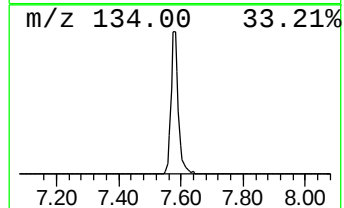
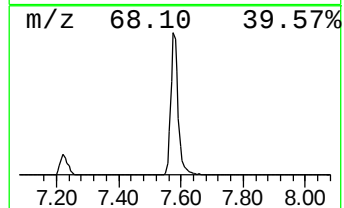
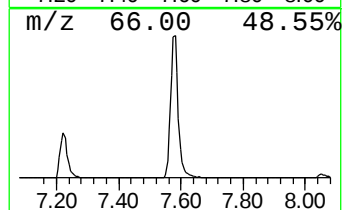
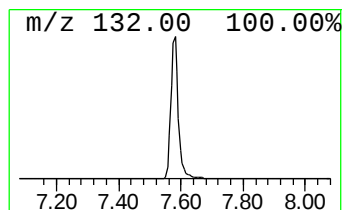
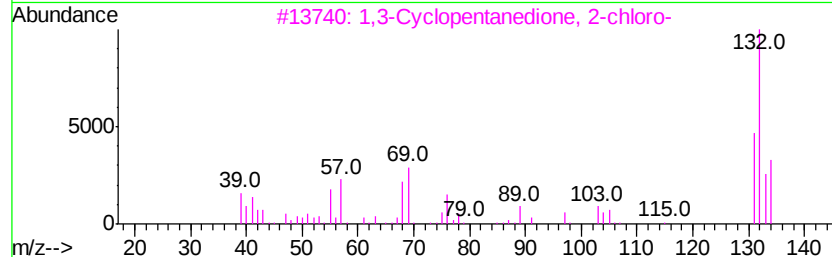
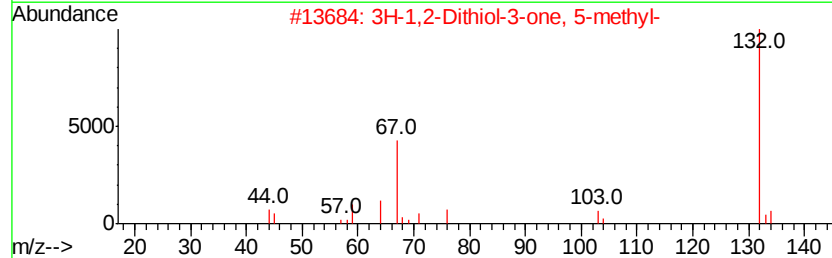
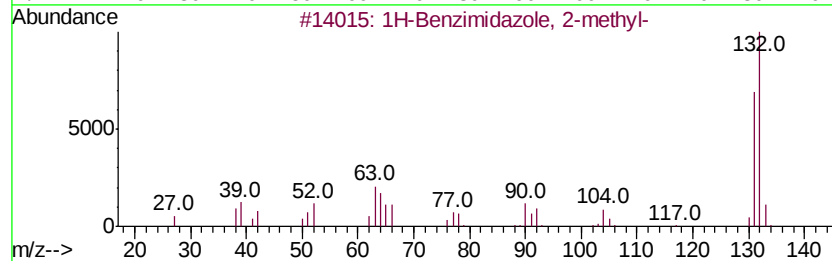
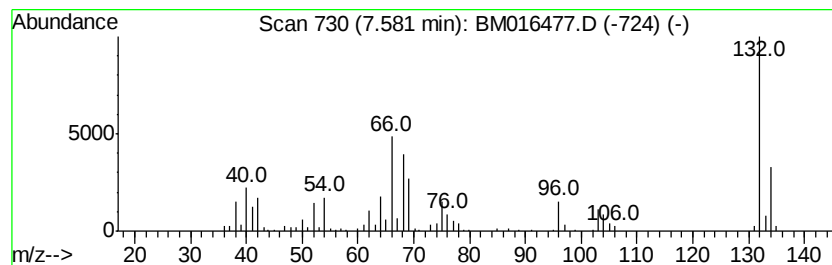
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Peak Number 2 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	169.61 ng	692514	1,4-Dichlorobenzene-d4	8.05

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
2			3H-1,2-Dithiol-3-one, 5-methyl-	132	C4H4OS2	003620-08-4	14
3			1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	10
4			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	10
5			(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown5.11	5.11	19.3	ng	78694	1	8.05	81660	20.0
unknown7.58	7.58	169.6	ng	692514	1	8.05	81660	20.0