

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012819\
 Data File : BM018668.D
 Acq On : 28 Jan 2019 17:32
 Operator : JU/SJ
 Sample : PB116655BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB116655BL

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_M\METHODS\8270-BM010919.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	4.869	298	303	313	rBV	183174	264302	1.82%	0.356%
2	5.316	372	379	395	rBV	4303549	6037117	41.53%	8.142%
3	6.898	641	648	666	rBV	3987674	6394938	44.00%	8.625%
4	7.257	702	709	725	rBV	4359600	6909432	47.53%	9.319%
5	7.722	782	788	798	rVB	759788	1217081	8.37%	1.641%
6	8.039	835	842	850	rBV	3345947	5367794	36.93%	7.239%
7	8.892	979	987	998	rBV	2436831	4005149	27.55%	5.402%
8	10.516	1255	1263	1276	rBV	937429	1593833	10.97%	2.150%
9	12.992	1676	1684	1698	rBV	6652984	9666369	66.50%	13.037%
10	14.374	1913	1919	1933	rVB2	1328528	2075198	14.28%	2.799%
11	15.868	2167	2173	2187	rBV	5448438	7574208	52.11%	10.215%
12	17.127	2380	2387	2395	rBV	1770666	2488609	17.12%	3.356%
13	19.762	2829	2835	2840	rBV	11316673	14535540	100.00%	19.604%
14	21.321	3095	3100	3105	rBV	2431014	3020279	20.78%	4.073%
15	23.586	3479	3485	3493	rVB	1592656	2997001	20.62%	4.042%

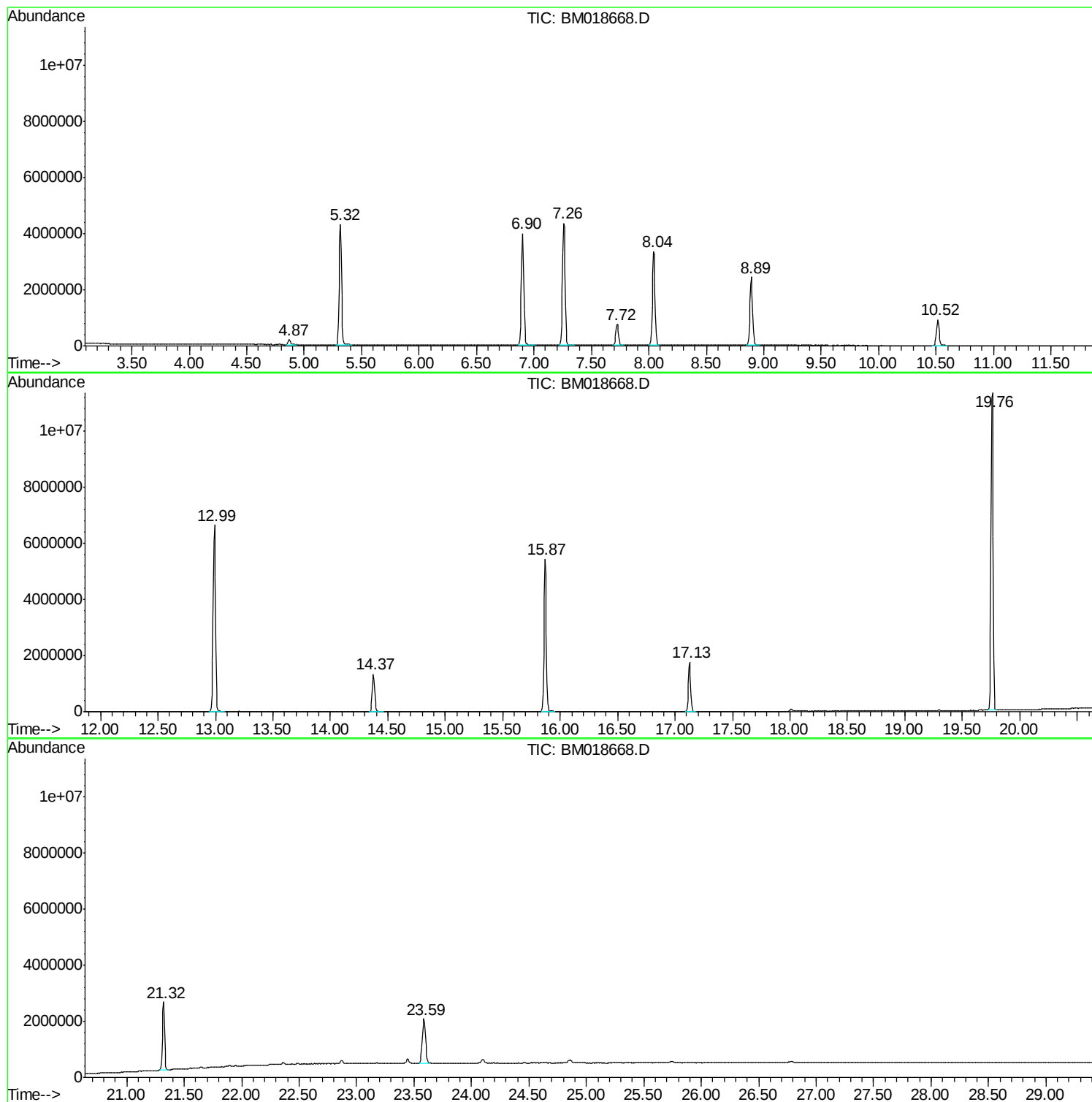
Sum of corrected areas: 74146850

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



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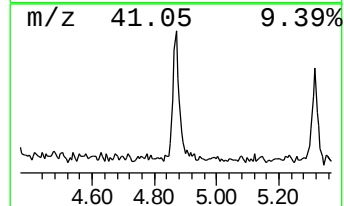
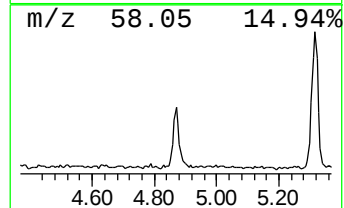
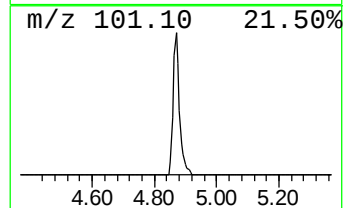
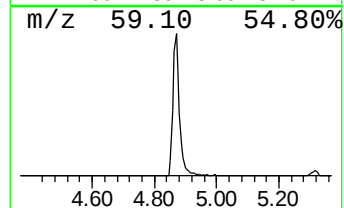
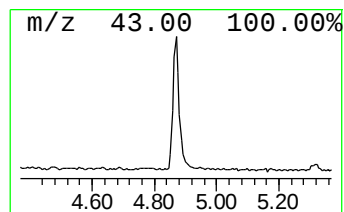
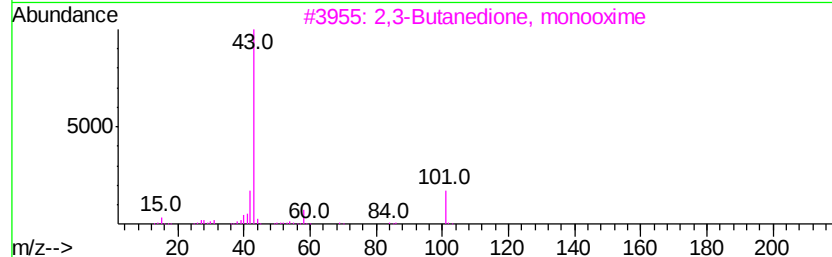
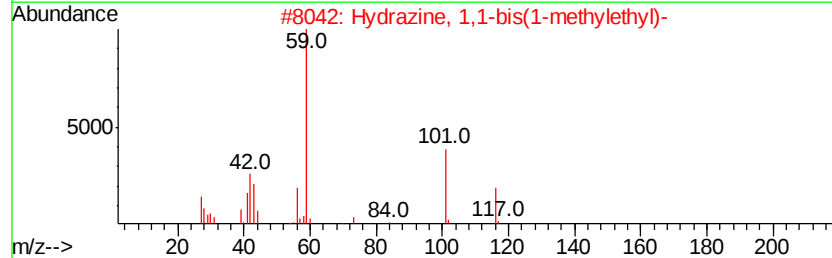
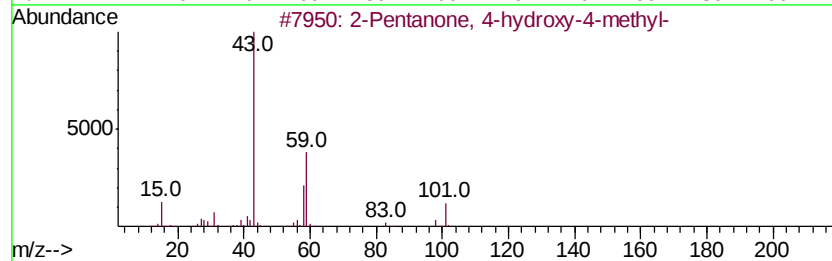
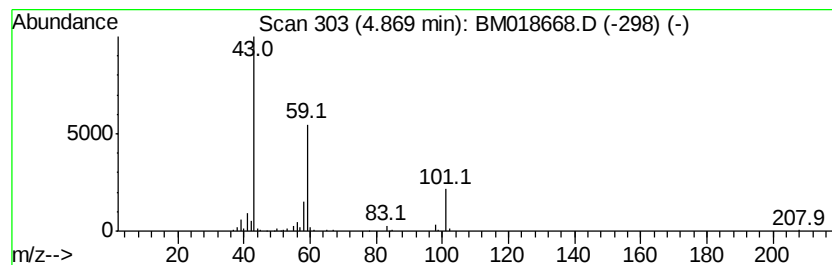
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.87	4.34 ng	264302	1,4-Dichlorobenzene-d4	7.73

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2			Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	33
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
4			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
5			Propanenitrile, 3-ethoxy-	99	C5H9NO	002141-62-0	9



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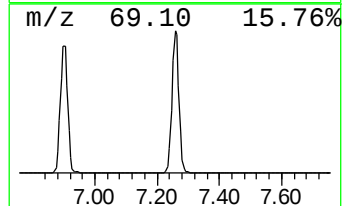
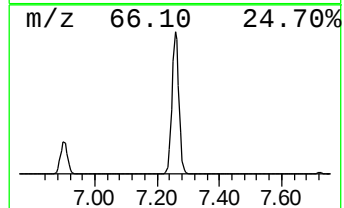
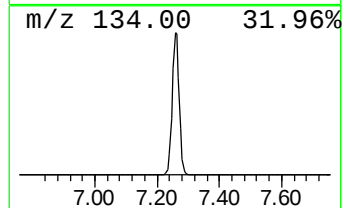
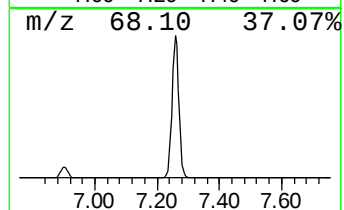
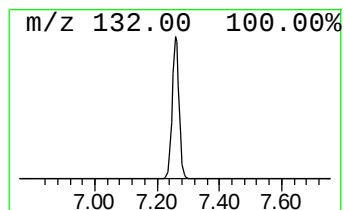
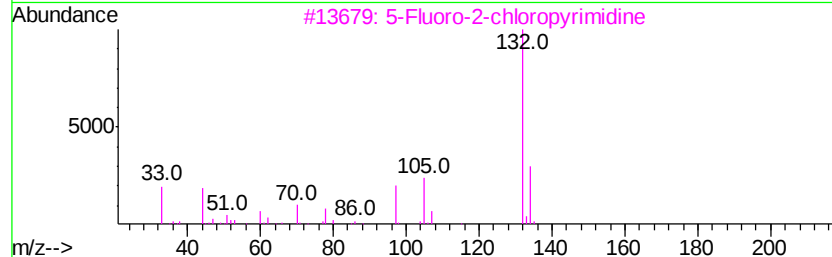
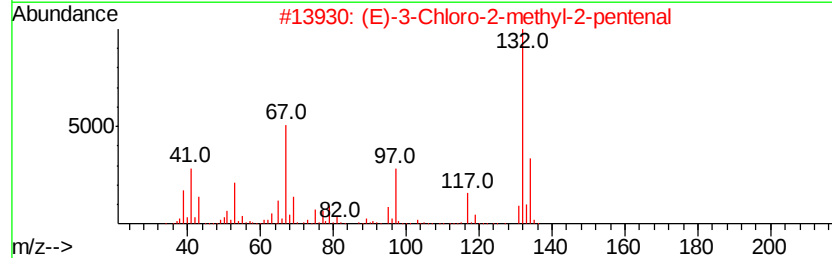
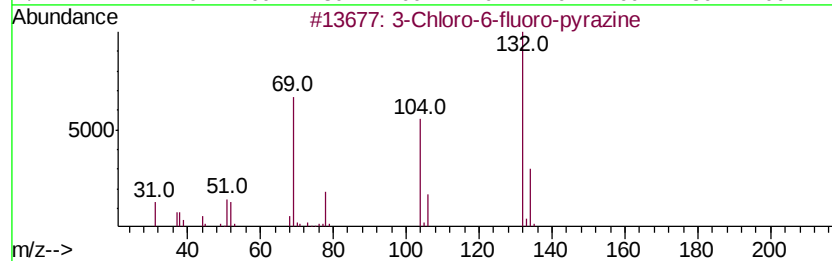
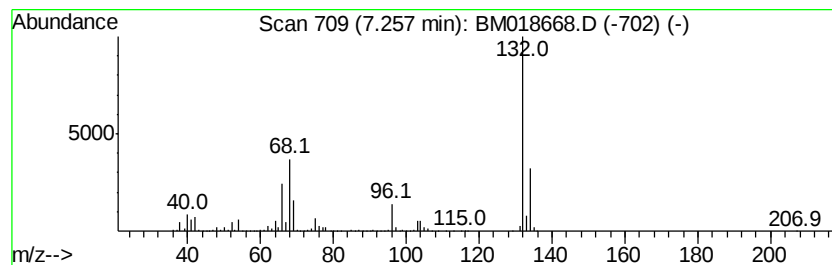
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Peak Number 2 unknown7.26 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.26	113.54 ng	6909430	1,4-Dichlorobenzene-d4	7.73

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	25
4		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
5		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.87	4.3	ng	264302	1	7.73	1217080	20.0
unknown7.26	7.26	113.5	ng	6909430	1	7.73	1217080	20.0