

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM081518\
 Data File : BM016388.D
 Acq On : 15 Aug 2018 19:49
 Operator : SJ/JU
 Sample : PB112066BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB112066BL

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-BM072318.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	5.163	314	319	333	rBV	123988	182004	7.10%	1.348%
2	5.646	395	401	414	rBV	858894	1309161	51.04%	9.695%
3	7.287	673	680	694	rBV	822072	1297332	50.58%	9.608%
4	7.640	734	740	750	rBV	984679	1501039	58.52%	11.116%
5	8.110	815	820	829	rBB	135609	211726	8.25%	1.568%
6	8.428	868	874	887	rBB	721764	1143457	44.58%	8.468%
7	9.298	1015	1022	1037	rBV	493712	798954	31.15%	5.917%
8	10.922	1292	1298	1309	rBV	138980	233912	9.12%	1.732%
9	13.369	1707	1714	1723	rBV	1203224	1668214	65.04%	12.354%
10	14.745	1943	1948	1956	rBB2	169389	252998	9.86%	1.874%
11	16.239	2196	2202	2213	rBV	1008066	1397890	54.50%	10.352%
12	17.486	2407	2414	2420	rBV	201708	280355	10.93%	2.076%
13	20.062	2847	2852	2858	rBV	2188109	2564883	100.00%	18.995%
14	21.615	3111	3116	3121	rBV	274642	332618	12.97%	2.463%
15	23.944	3506	3512	3520	rBV	171207	328597	12.81%	2.433%

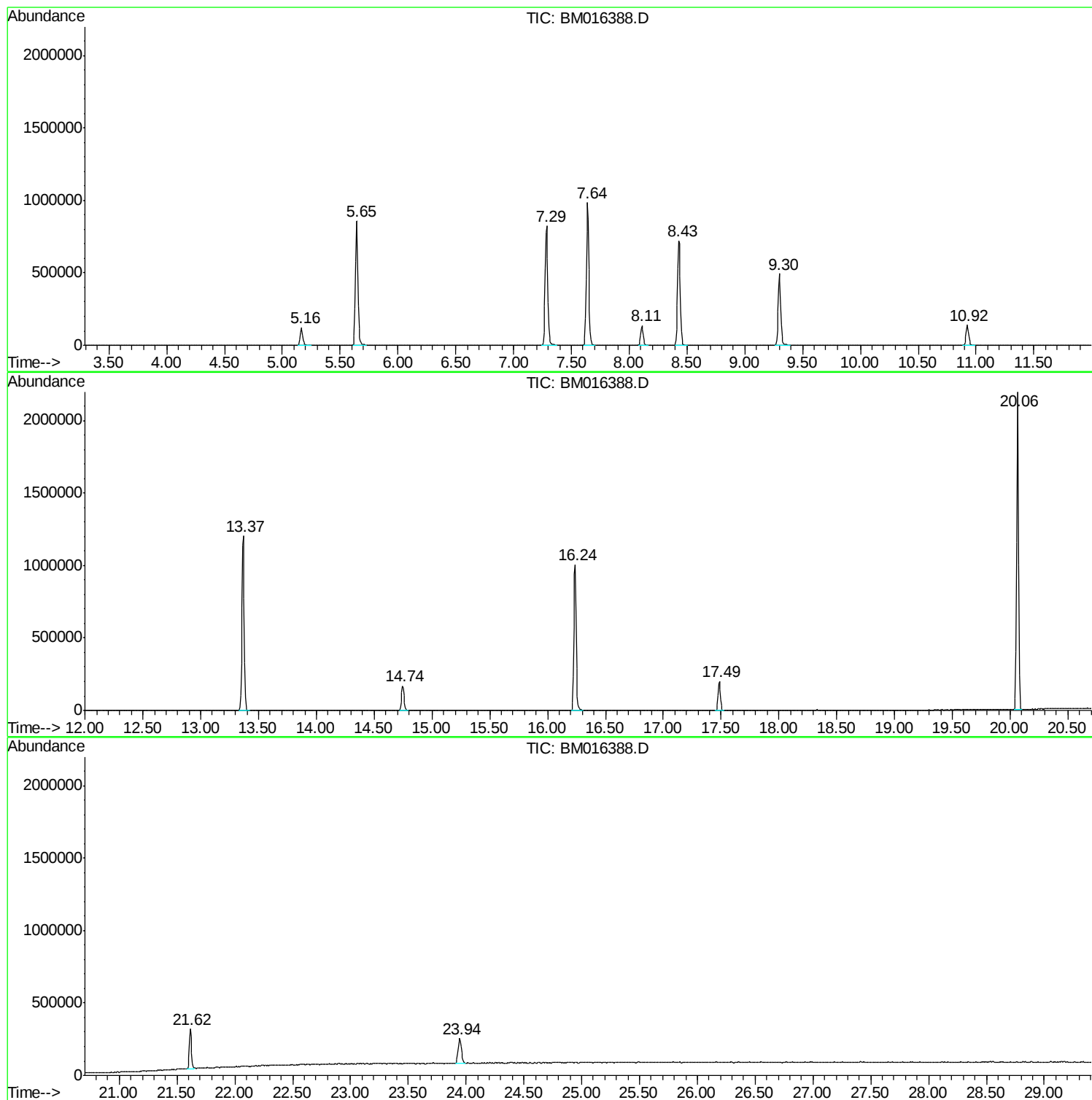
Sum of corrected areas: 13503140

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



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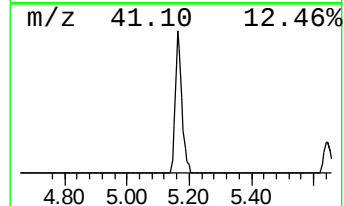
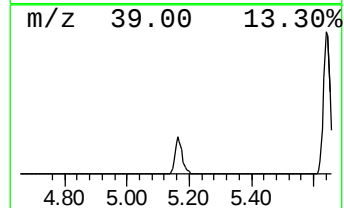
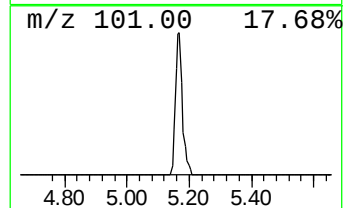
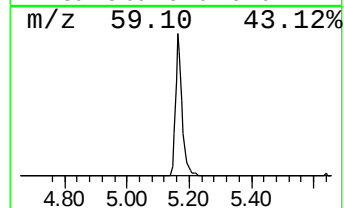
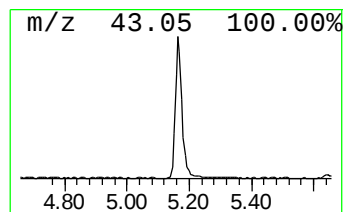
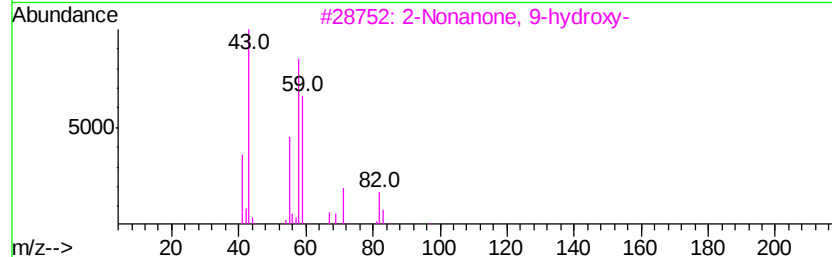
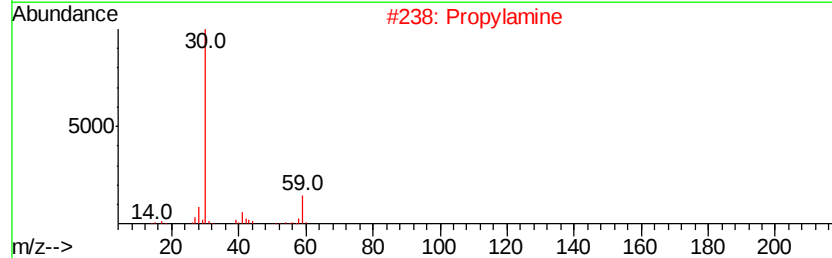
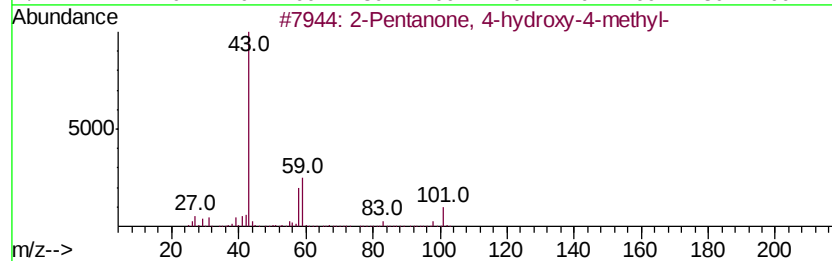
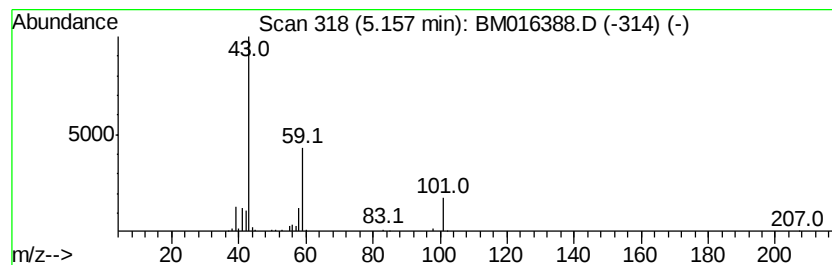
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	17.19 ng	182004	1,4-Dichlorobenzene-d4	8.11

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Propylamine	59	C3H9N	000107-10-8	42
3			2-Nonanone, 9-hydroxy-	158	C9H18O2	025368-56-3	36
4			Butane	58	C4H10	000106-97-8	22
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16



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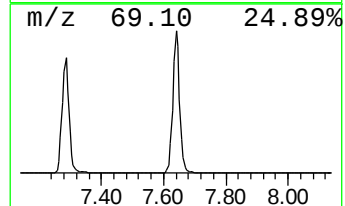
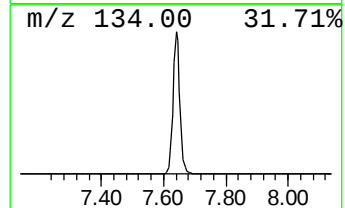
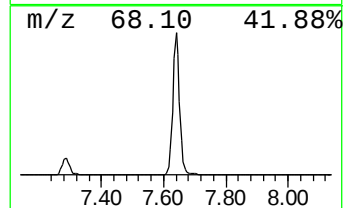
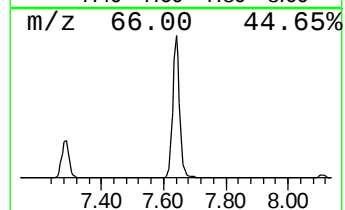
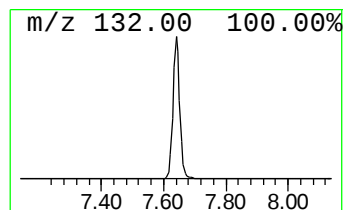
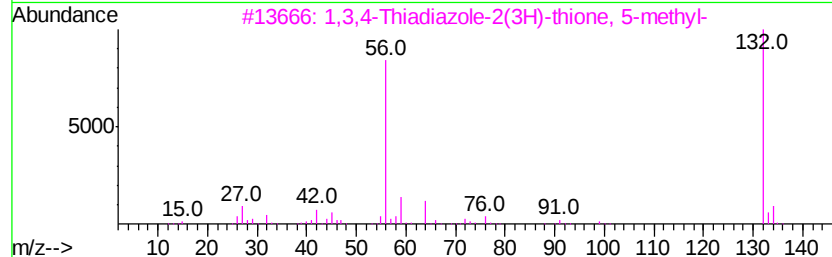
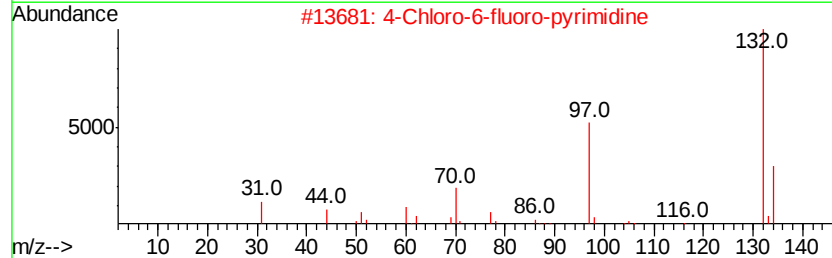
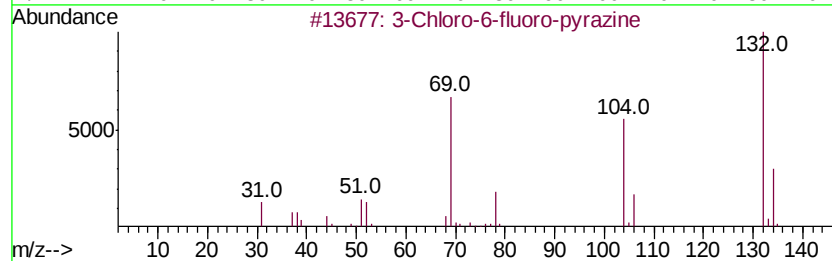
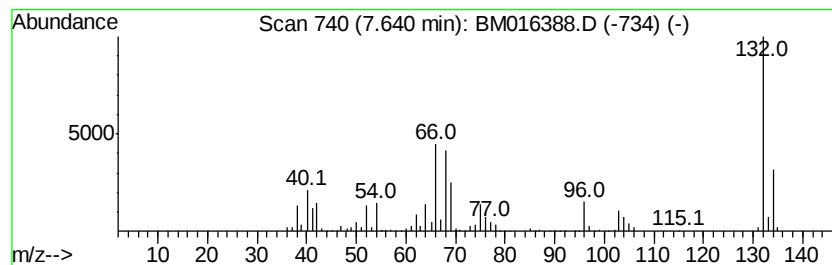
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Peak Number 2 unknown7.64 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.64	141.79 ng	1501040	1,4-Dichlorobenzene-d4	8.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	14
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	14
3		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	14
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	14
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.16	17.2	ng	182004	1	8.11	211726	20.0
unknown7.64	7.64	141.8	ng	1501040	1	8.11	211726	20.0