

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF021319\
 Data File : BF112530.D
 Acq On : 13 Feb 2019 11:06
 Operator : JU/SJ
 Sample : PB117061BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB117061BL

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-BF012519.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	5.069	503	507	518	rBV	84434	125220	2.94%	0.436%
2	5.481	573	577	601	rBV	2683205	2679616	62.89%	9.337%
3	6.487	744	748	766	rBV	2501366	2801600	65.75%	9.762%
4	6.616	766	770	773	rBV	3373036	2954025	69.33%	10.293%
5	6.834	804	807	819	rBV	727145	650579	15.27%	2.267%
6	6.992	830	834	853	rVB	2735248	2379857	55.85%	8.292%
7	7.404	899	904	924	rBV	1608124	1825340	42.84%	6.360%
8	8.116	1021	1025	1040	rBV	761797	834484	19.58%	2.908%
9	9.192	1202	1208	1211	rBV	3900657	3695274	86.72%	12.876%
10	9.875	1319	1324	1343	rBV	1009322	1003461	23.55%	3.496%
11	10.669	1455	1459	1482	rBV	2358670	2418280	56.75%	8.426%
12	11.363	1574	1577	1596	rVB	1067656	1078518	25.31%	3.758%
13	12.951	1842	1847	1850	rBV	4473403	4260969	100.00%	14.847%
14	14.016	2024	2028	2039	rBV	960343	1040930	24.43%	3.627%
15	15.092	2208	2211	2219	rVB2	41895	67669	1.59%	0.236%
16	15.492	2274	2279	2290	rVB	488788	834724	19.59%	2.908%
17	15.898	2345	2348	2352	rVB5	33888	49192	1.15%	0.171%

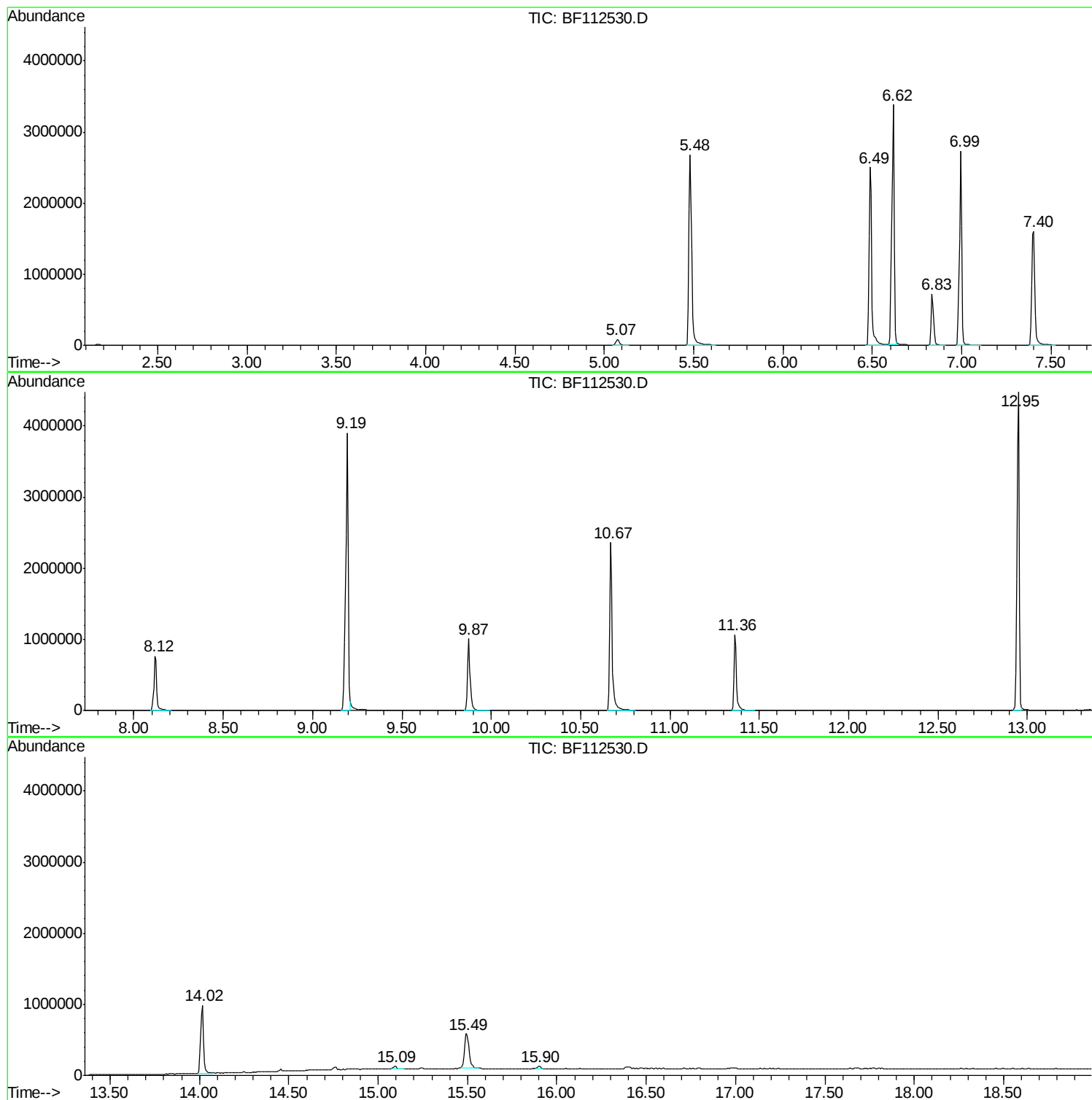
Sum of corrected areas: 28699738

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



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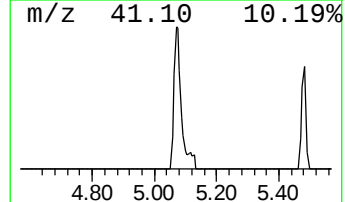
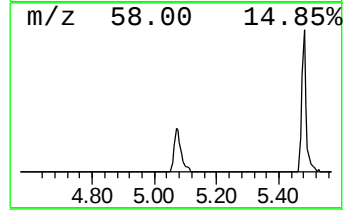
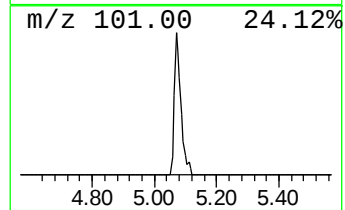
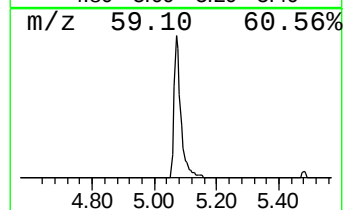
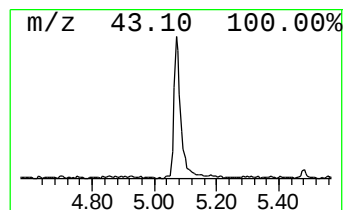
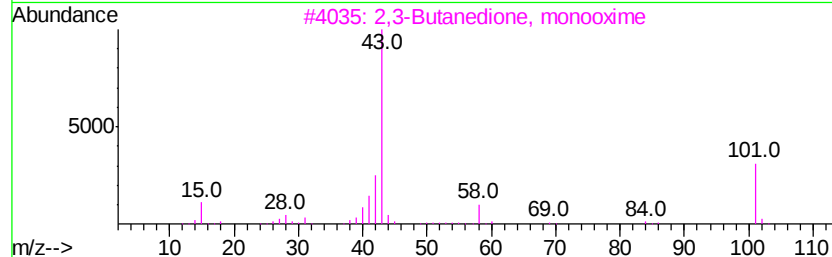
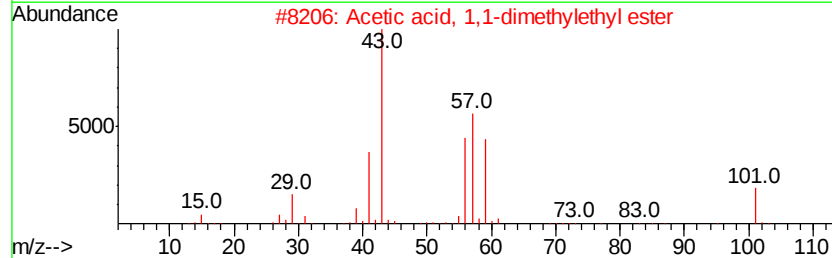
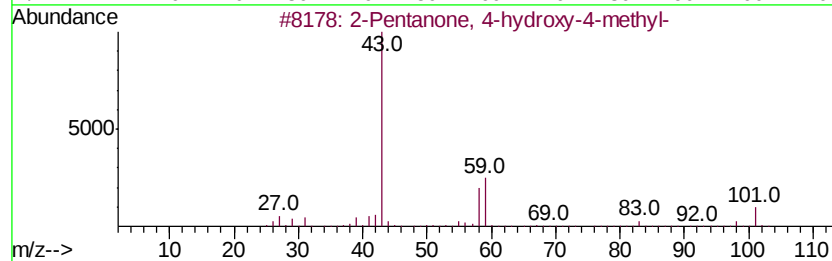
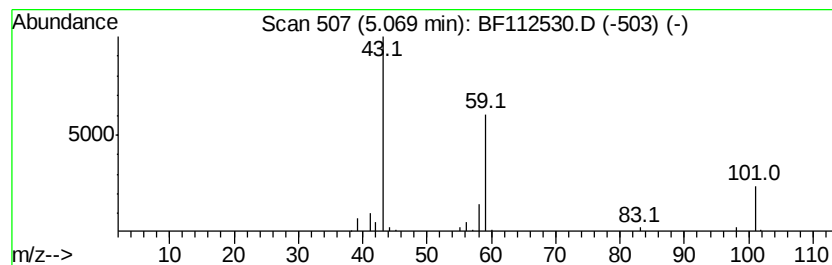
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.07	3.85 ng	125220	1,4-Dichlorobenzene-d4	6.83

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72	
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39	
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	32	
4	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9	
5	5-Hexen-2-one	98	C6H10O	000109-49-9	9	



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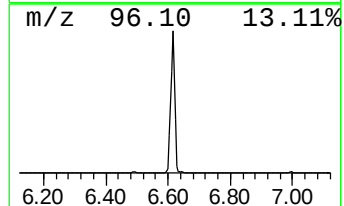
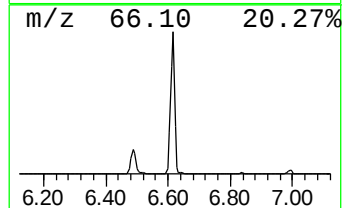
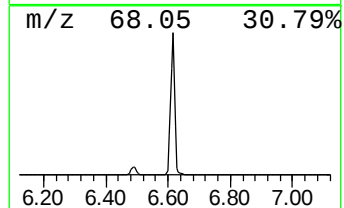
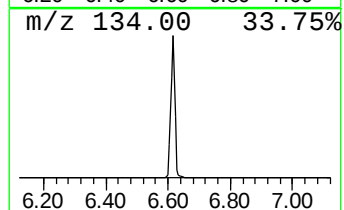
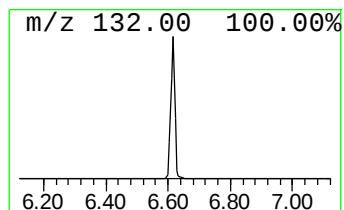
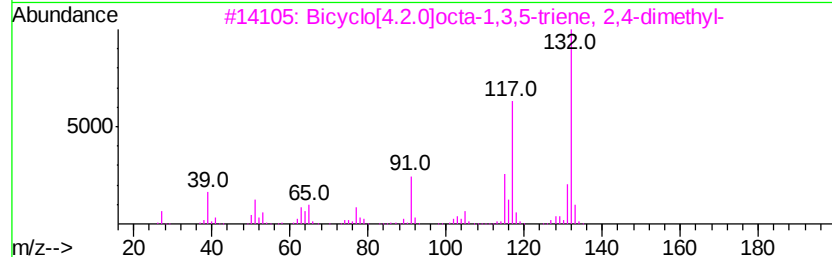
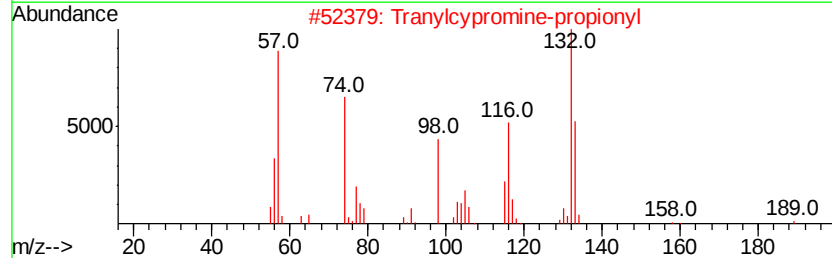
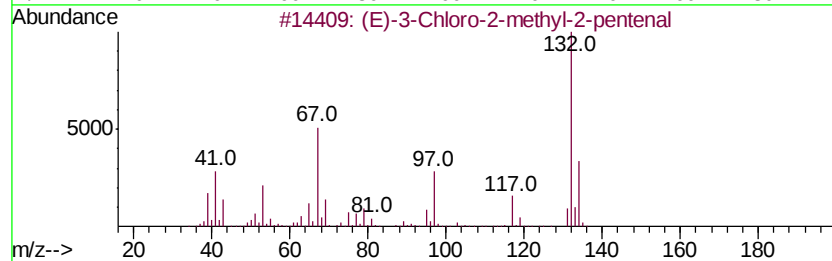
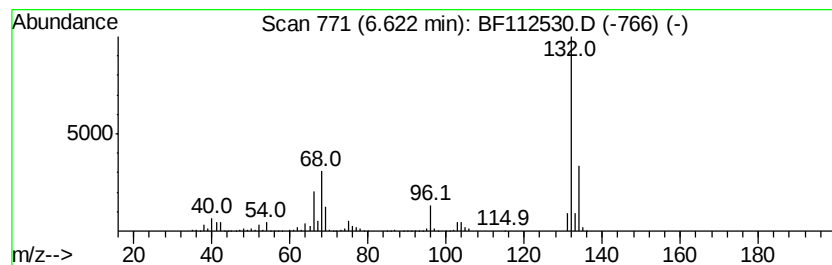
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Peak Number 2 unknown6.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	90.81 ng	2954030	1,4-Dichlorobenzene-d4	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.07	3.9	ng	125220	1	6.83	650579	20.0
unknown6.62	6.62	90.8	ng	2954030	1	6.83	650579	20.0