

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF040218\
 Data File : BF104127.D
 Acq On : 2 Apr 2018 11:45
 Operator : JU/SJ
 Sample : PB107878BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB107878BL

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:\HPCHEM1\BNA_F\METHODS\8270-BF032818.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	5.198	526	529	540	rBV	97220	136022	4.04%	0.589%
2	5.592	593	596	623	rBV	1475695	2067919	61.48%	8.953%
3	6.592	762	766	786	rBV	1456262	2137456	63.54%	9.254%
4	6.728	786	789	820	rVB	1772480	2450783	72.86%	10.611%
5	6.957	825	828	845	rVV	293374	457440	13.60%	1.981%
6	7.110	850	854	884	rVB	1731604	1914027	56.90%	8.287%
7	7.522	920	924	962	rBV	888315	1447128	43.02%	6.266%
8	8.239	1042	1046	1071	rBV	509285	709581	21.10%	3.072%
9	9.310	1224	1228	1251	rBV	2644873	2986982	88.80%	12.933%
10	9.998	1341	1345	1366	rBV	733346	832443	24.75%	3.604%
11	10.792	1476	1480	1506	rBV	1603955	1907666	56.71%	8.259%
12	11.492	1595	1599	1623	rBV	713968	898931	26.72%	3.892%
13	13.080	1864	1869	1890	rBV	3028785	3363710	100.00%	14.564%
14	14.145	2046	2050	2062	rBV	807467	942277	28.01%	4.080%
15	15.686	2303	2312	2326	rBV	492509	844342	25.10%	3.656%

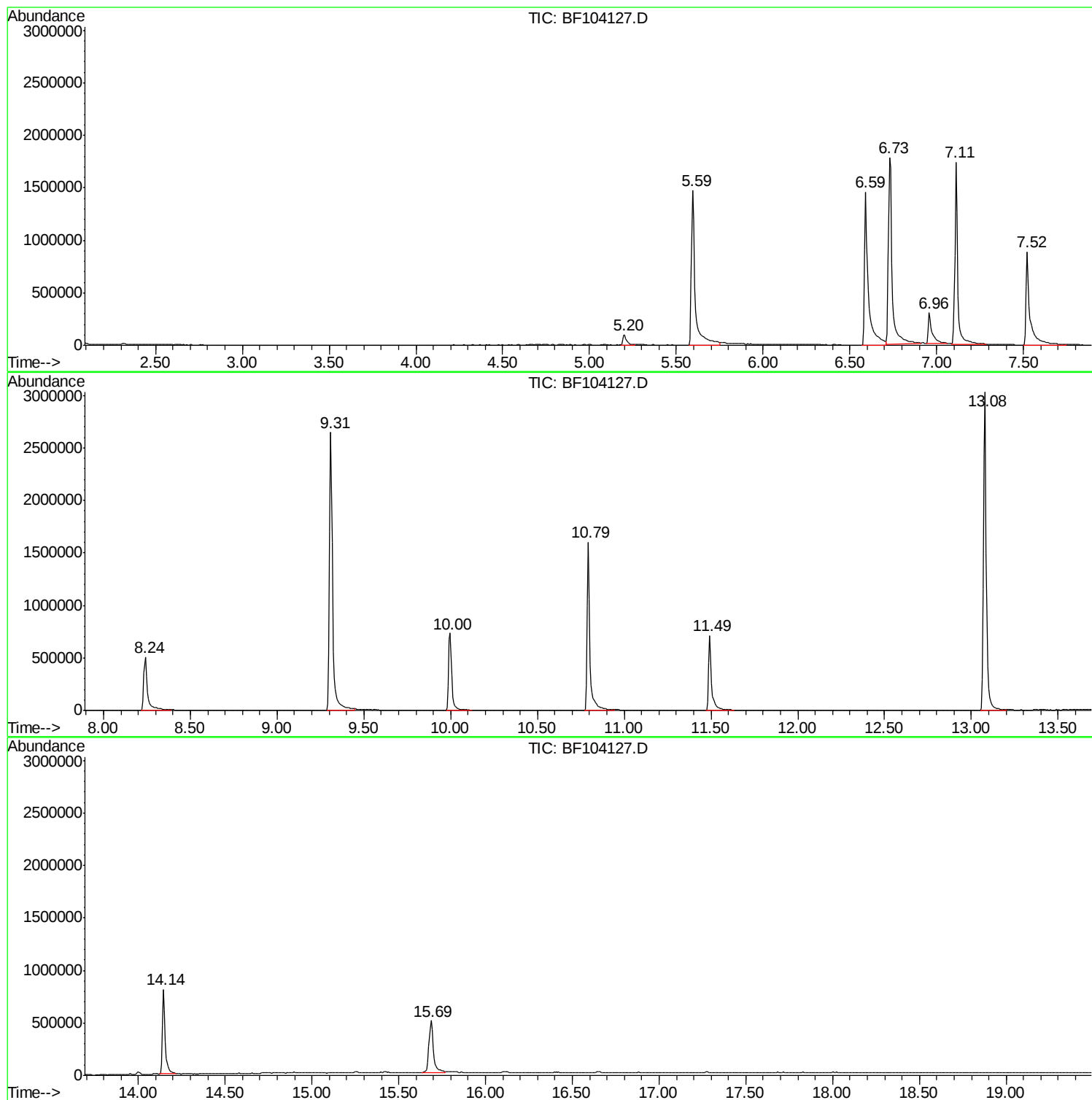
Sum of corrected areas: 23096707

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



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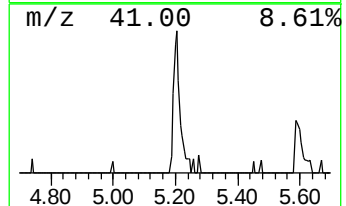
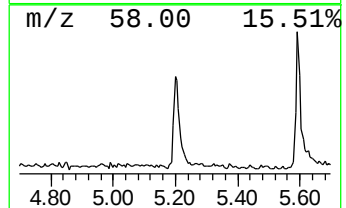
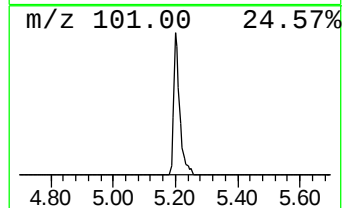
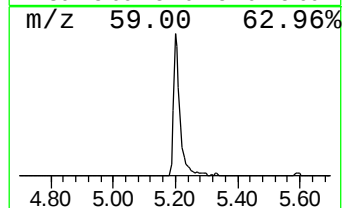
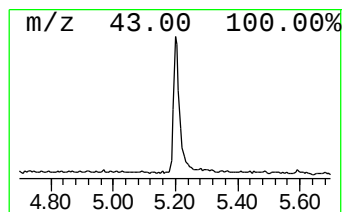
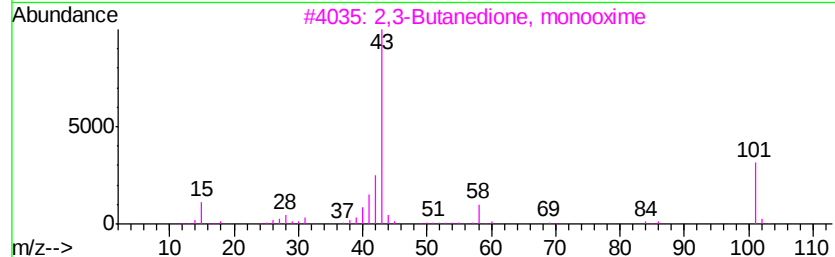
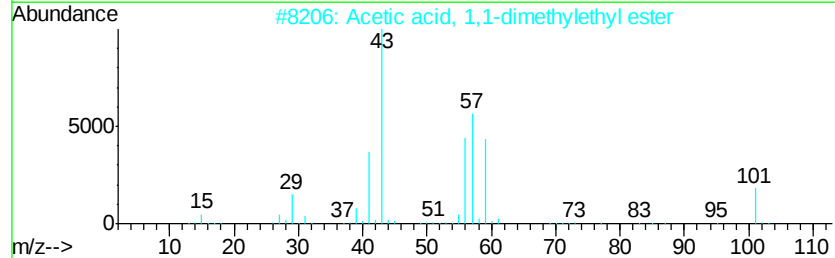
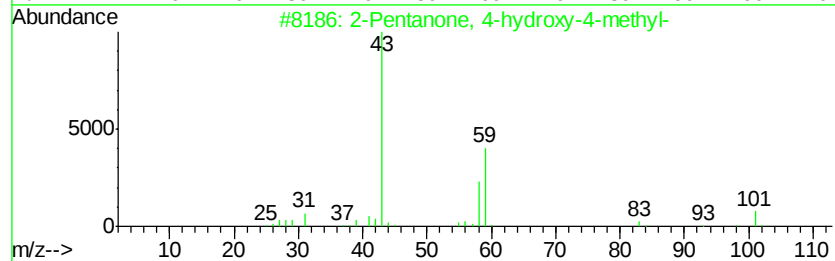
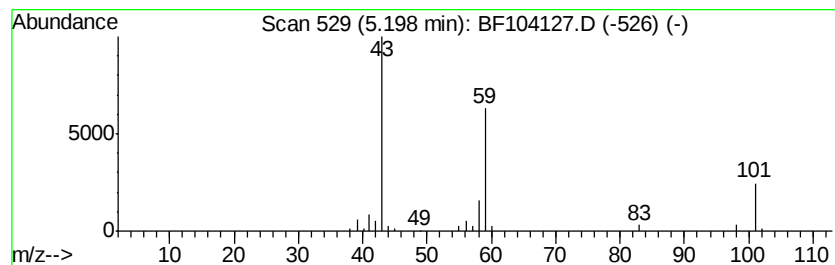
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	5.95 ng	136022	1,4-Dichlorobenzene-d4	6.96

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
5			Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	17



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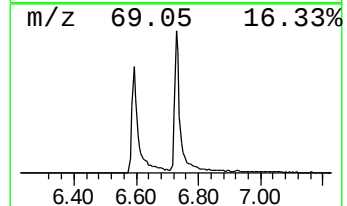
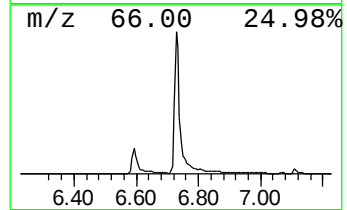
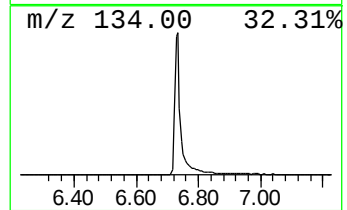
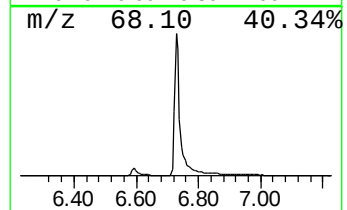
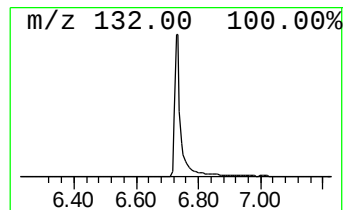
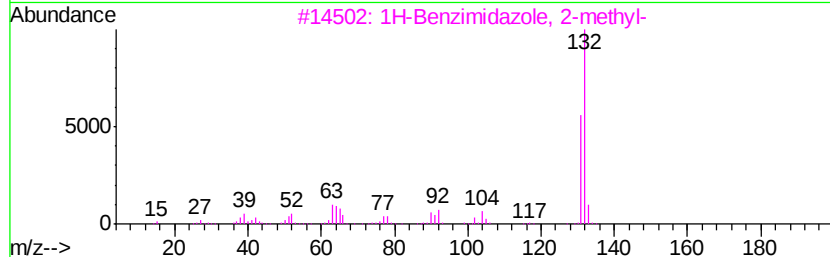
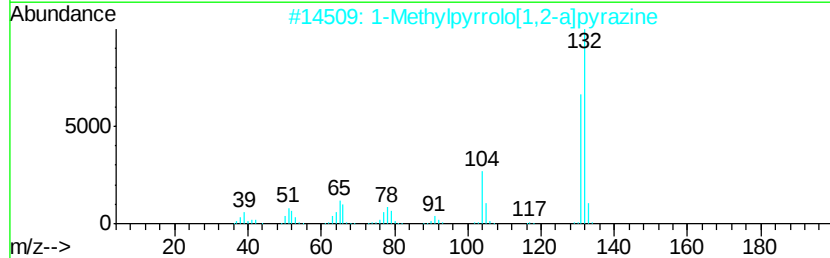
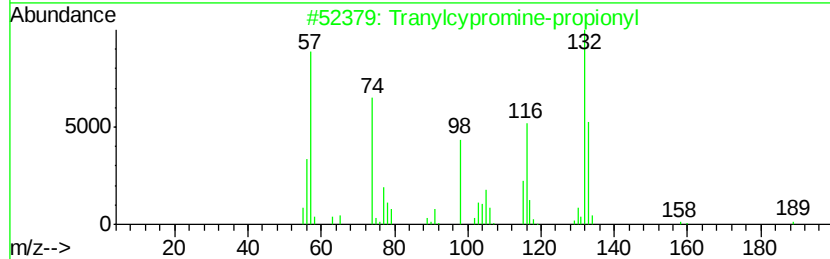
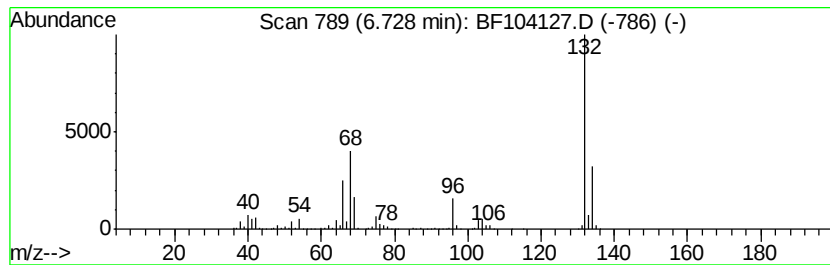
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Peak Number 2 unknown6.73 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.73	107.15 ng	2450780	1,4-Dichlorobenzene-d4	6.96

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
2			1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
3			1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	9
4			2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5			(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.20	6.0	ng	136022	1	6.96	457440	20.0
unknown6.73	6.73	107.2	ng	2450780	1	6.96	457440	20.0