

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN072518\
 Data File : BN002152.D
 Acq On : 26 Jul 2018 00:47
 Operator : JU/SJ
 Sample : PB111180BL
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB111180BL

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_N\METHODS\8270-BN072518.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	4.858	313	318	332	rVB	747972	1070862	10.74%	1.505%
2	5.311	389	395	410	rBV	4799302	6922494	69.43%	9.731%
3	6.881	654	662	676	rBV	4447896	7200035	72.21%	10.121%
4	7.234	714	722	737	rBV	4688128	7339764	73.61%	10.318%
5	7.693	793	800	809	rVB	936376	1497145	15.02%	2.105%
6	8.011	846	854	865	rVB	3532443	5643657	56.60%	7.933%
7	8.840	988	995	1007	rBV	2668966	4518524	45.32%	6.352%
8	10.463	1264	1271	1280	rBV	1182452	1948981	19.55%	2.740%
9	12.940	1685	1692	1702	rBV	6418100	9502625	95.31%	13.358%
10	14.322	1920	1927	1937	rBV2	1571835	2367211	23.74%	3.328%
11	15.816	2174	2181	2199	rBV	4483934	6356357	63.75%	8.935%
12	17.069	2388	2394	2407	rBV	1802783	2544301	25.52%	3.577%
13	19.710	2837	2843	2849	rBV	7926599	9970536	100.00%	14.016%
14	21.269	3102	3108	3116	rBV	1835507	2273693	22.80%	3.196%
15	23.492	3479	3486	3499	rVB	1048998	1982615	19.88%	2.787%

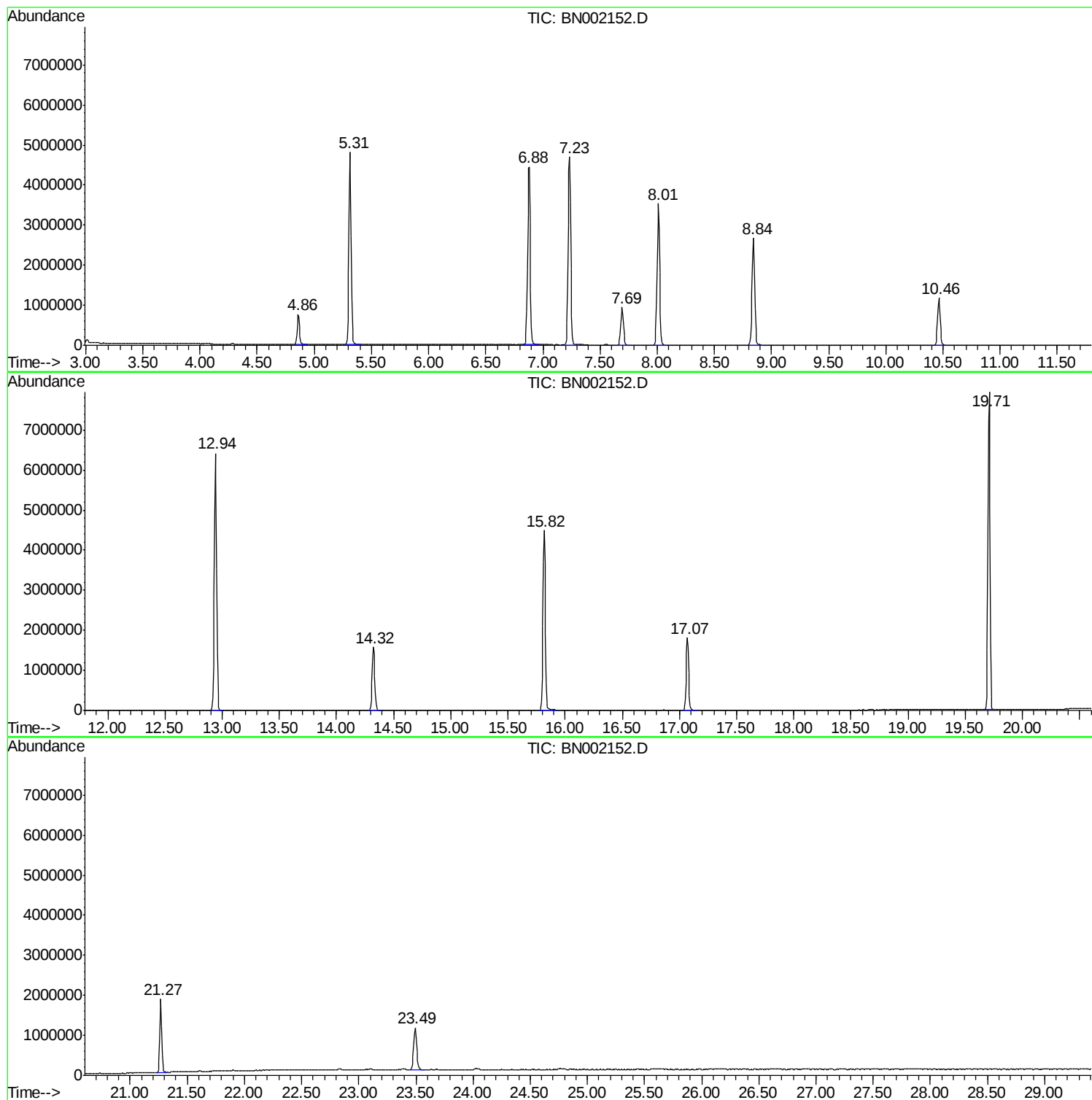
Sum of corrected areas: 71138800

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



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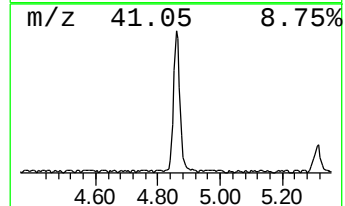
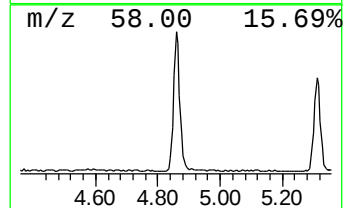
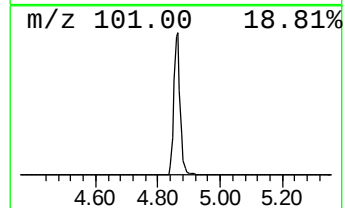
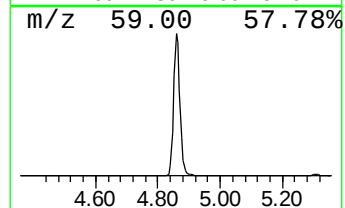
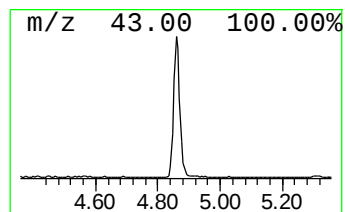
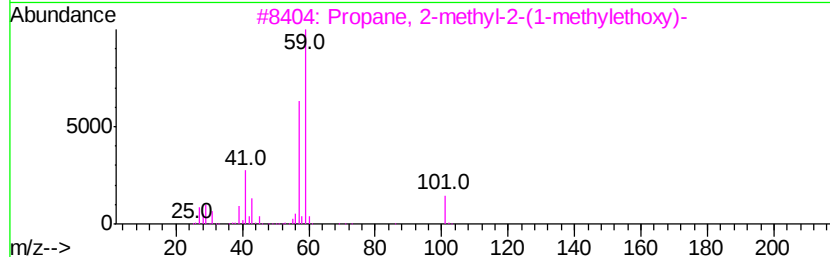
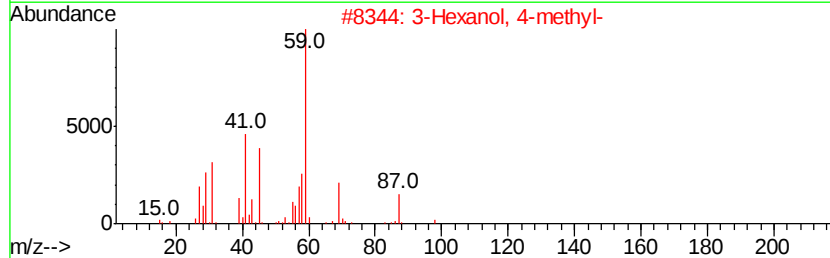
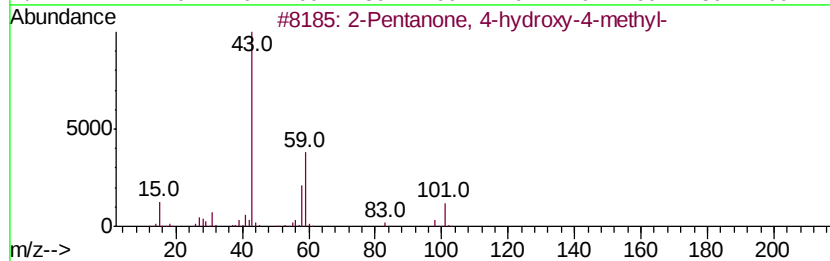
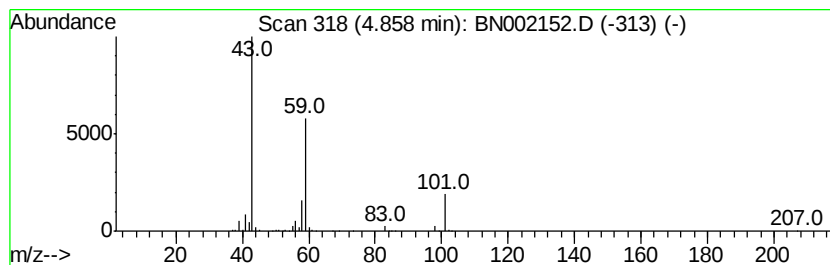
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.86	14.31 ng	1070860	1,4-Dichlorobenzene-d4	7.69

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3			Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	33
4			Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	25
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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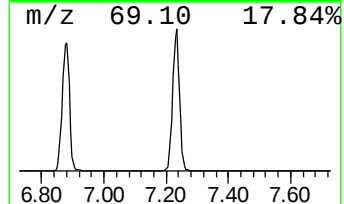
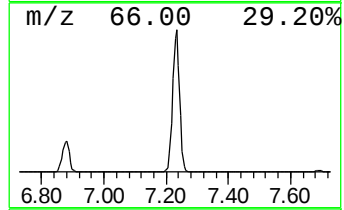
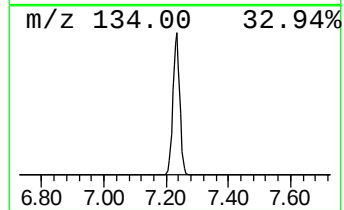
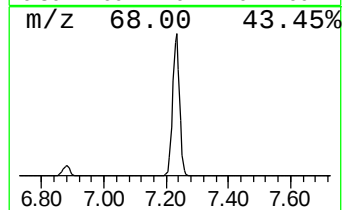
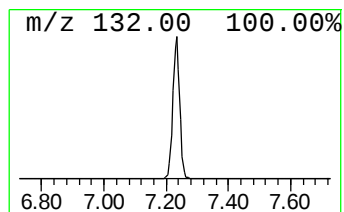
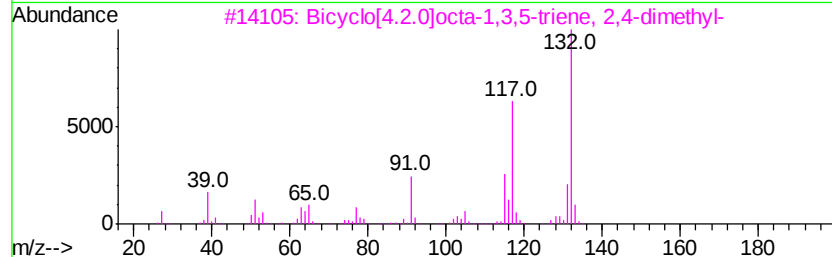
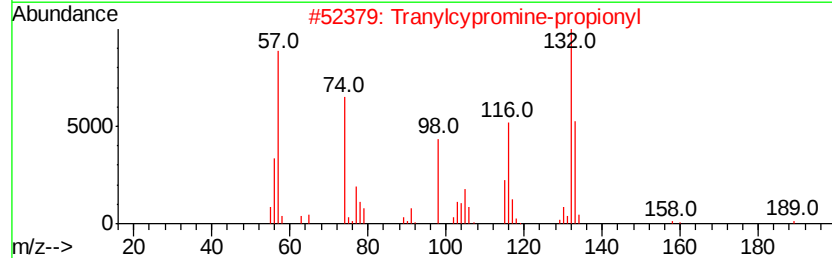
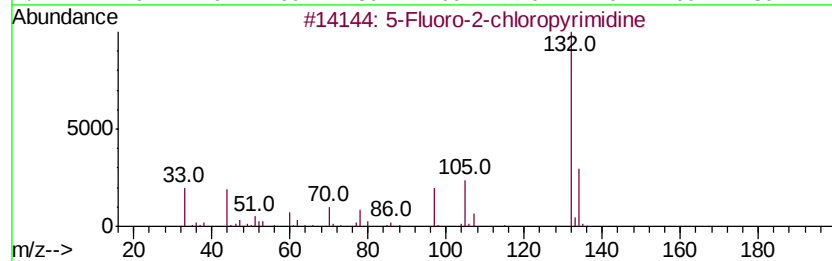
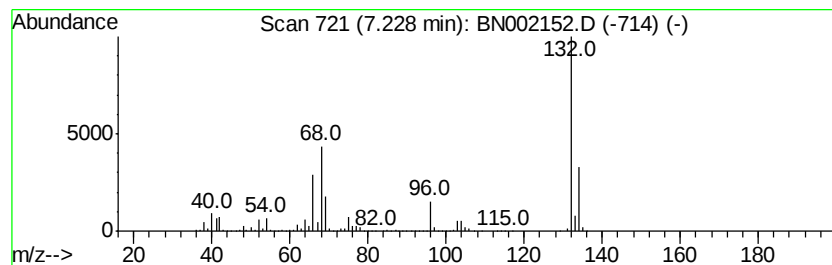
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Peak Number 2 unknown7.23 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.23	98.05 ng	7339760	1,4-Dichlorobenzene-d4	7.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
5		Xenon	132	Xe	007440-63-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.86	14.3	ng	1070860	1	7.69	1497150	20.0
unknown7.23	7.23	98.0	ng	7339760	1	7.69	1497150	20.0