

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111617\
 Data File : BF100647.D
 Acq On : 16 Nov 2017 23:38
 Operator : SJ/JU
 Sample : PB104306BL
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB104306BL

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-BF110817.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	5.110	510	514	524	rBV	373555	419754	17.06%	2.500%
2	5.504	576	581	584	rBV	1537896	1371733	55.76%	8.169%
3	6.504	746	751	754	rBV	1474718	1359543	55.26%	8.096%
4	6.651	771	776	779	rBV	1555812	1417243	57.61%	8.440%
5	6.881	812	815	819	rBB	744564	599052	24.35%	3.567%
6	7.039	837	842	845	rBV	1900336	1664267	67.65%	9.911%
7	7.445	905	911	914	rBV	1326510	1296789	52.71%	7.722%
8	8.163	1028	1033	1037	rBV	896575	781794	31.78%	4.656%
9	9.239	1210	1216	1219	rBV	2689200	2460076	100.00%	14.650%
10	9.922	1327	1332	1335	rBV	829187	828021	33.66%	4.931%
11	10.704	1461	1465	1479	rBB	822194	722354	29.36%	4.302%
12	11.404	1580	1584	1588	rBV	930193	767349	31.19%	4.570%
13	12.986	1849	1853	1857	rBV	2047094	1937963	78.78%	11.541%
14	14.039	2028	2032	2036	rBV	688849	606260	24.64%	3.610%
15	15.515	2277	2283	2290	rBV	410440	534180	21.71%	3.181%
16	16.480	2443	2447	2453	rVB2	16430	26152	1.06%	0.156%

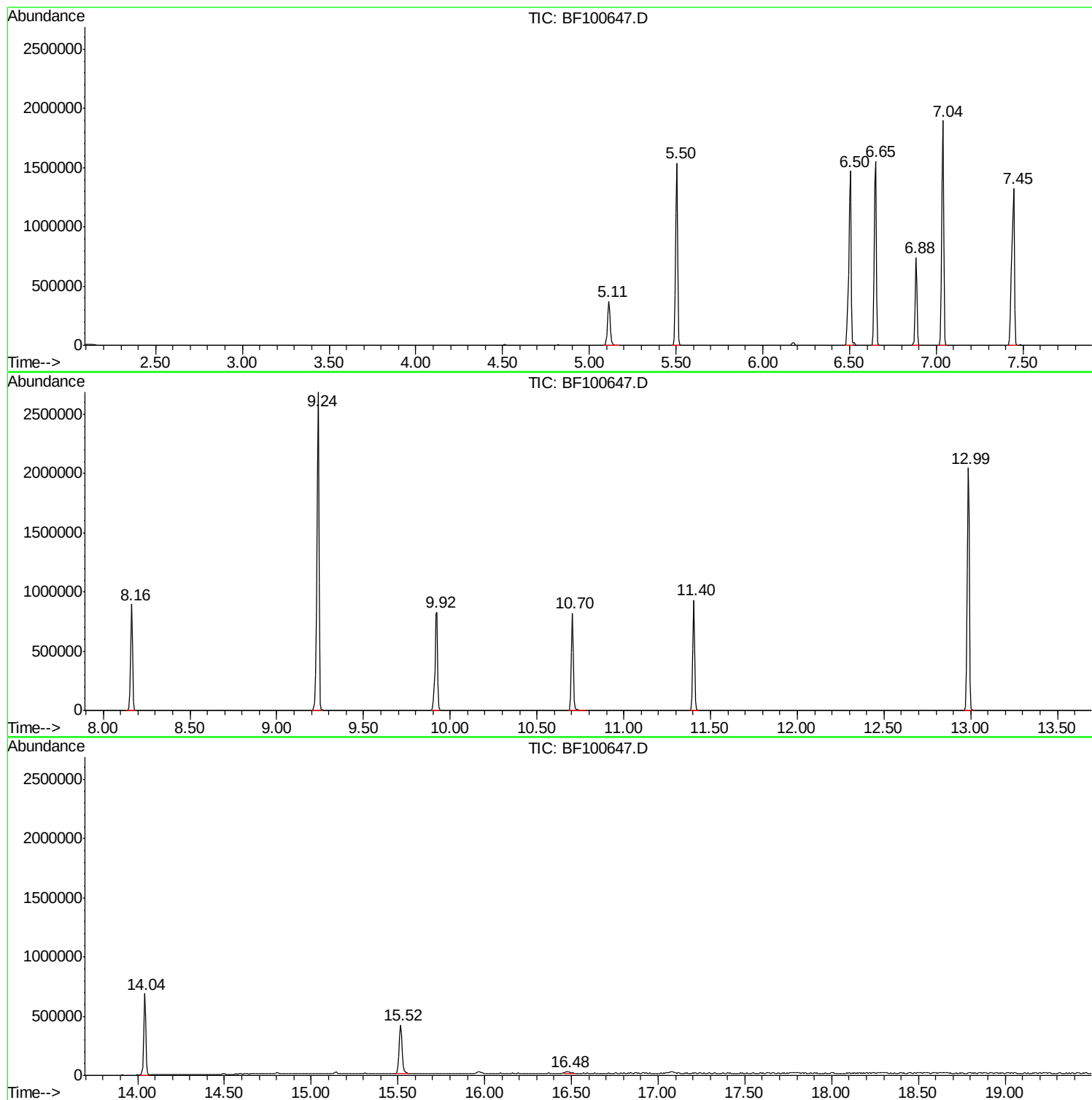
Sum of corrected areas: 16792530

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



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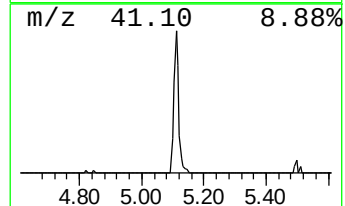
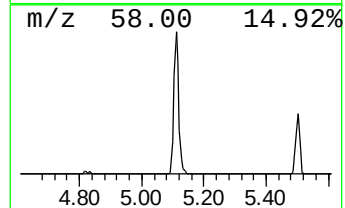
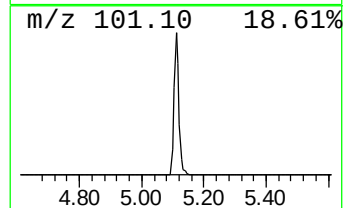
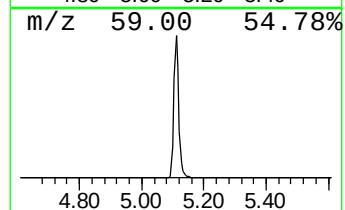
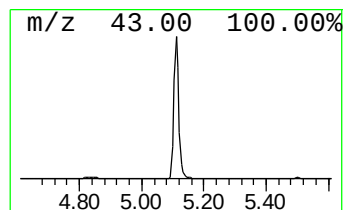
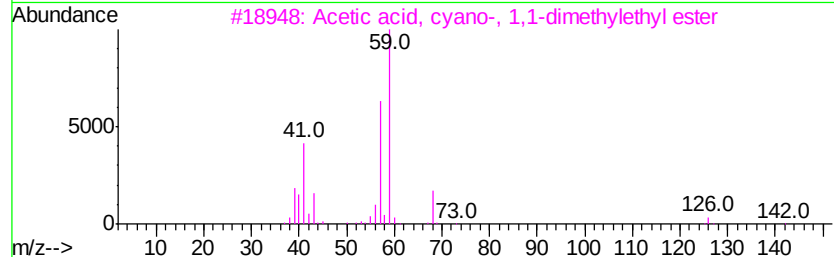
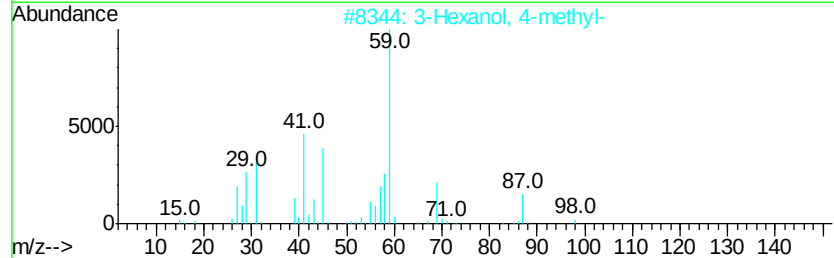
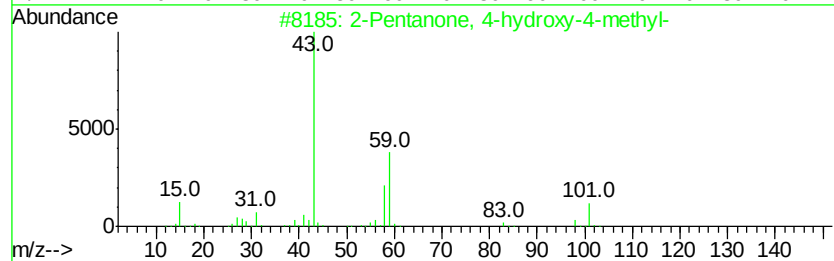
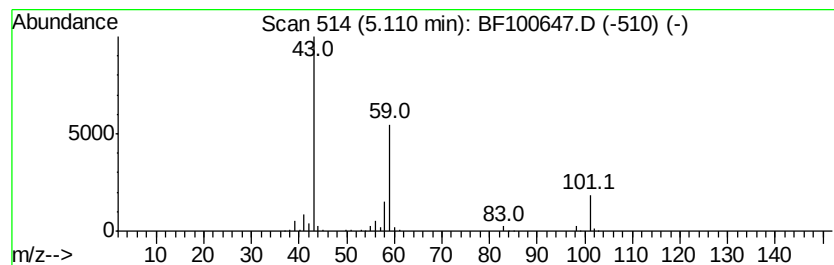
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.11	14.01 ng	419754	1,4-Dichlorobenzene-d4	6.88

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	32
4			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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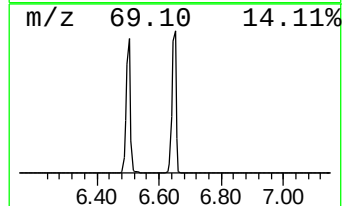
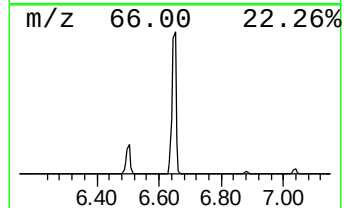
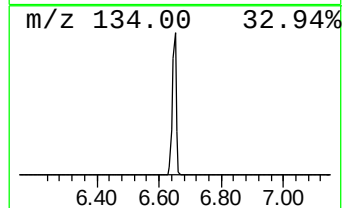
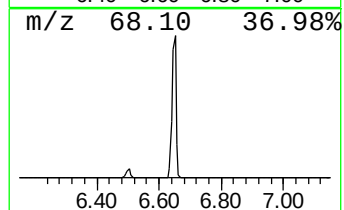
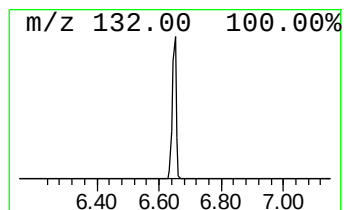
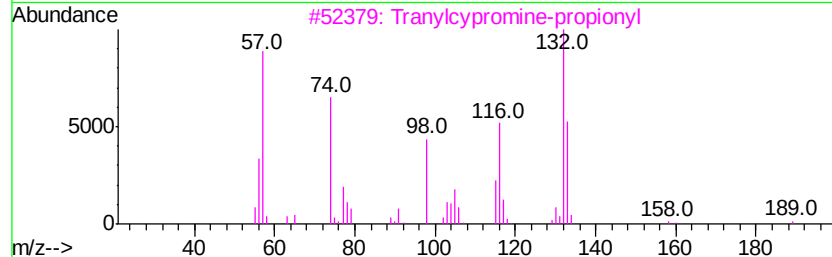
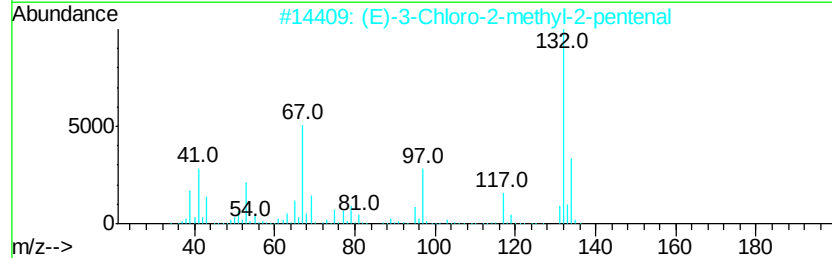
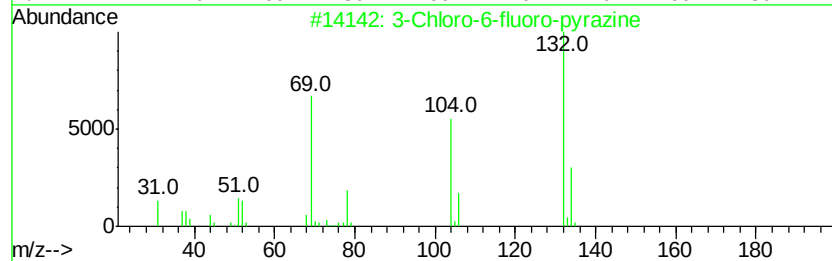
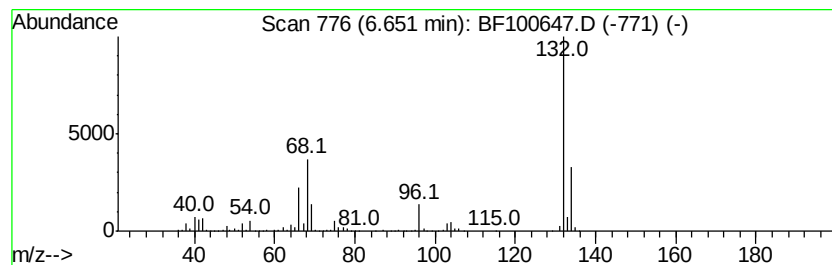
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Peak Number 2 unknown6.65 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.65	47.32 ng	1417240	1,4-Dichlorobenzene-d4	6.88

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	17
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Amphetaminil	250	C17H18N2	017590-01-1	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.11	14.0	ng	419754	1	6.88	599052	20.0
unknown6.65	6.65	47.3	ng	1417240	1	6.88	599052	20.0