

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111617\
 Data File : BF100623.D
 Acq On : 16 Nov 2017 11:46
 Operator : SJ/JU
 Sample : PB104287BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB104287BL

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	5.116	511	515	526	rBV	324085	404118	14.63%	1.858%
2	5.504	576	581	584	rBV	2128424	2232411	80.83%	10.266%
3	6.504	746	751	754	rBV	2064891	2190824	79.32%	10.075%
4	6.651	771	776	779	rBV	2363236	2242598	81.20%	10.313%
5	6.881	811	815	818	rBV	723185	572801	20.74%	2.634%
6	7.039	837	842	845	rBV	1937539	1802250	65.25%	8.288%
7	7.439	905	910	913	rBV	1399415	1423378	51.54%	6.545%
8	8.163	1028	1033	1036	rBV	852819	767315	27.78%	3.529%
9	9.233	1210	1215	1219	rBV	2590304	2712810	98.22%	12.475%
10	9.916	1326	1331	1335	rBV2	973923	862586	31.23%	3.967%
11	10.704	1460	1465	1468	rBV	1595201	1455911	52.71%	6.695%
12	11.398	1579	1583	1587	rBV	1050089	929577	33.66%	4.275%
13	12.986	1848	1853	1856	rBV	2822604	2761860	100.00%	12.700%
14	14.039	2028	2032	2035	rBV	910249	818581	29.64%	3.764%
15	15.509	2277	2282	2289	rBV	457895	569122	20.61%	2.617%

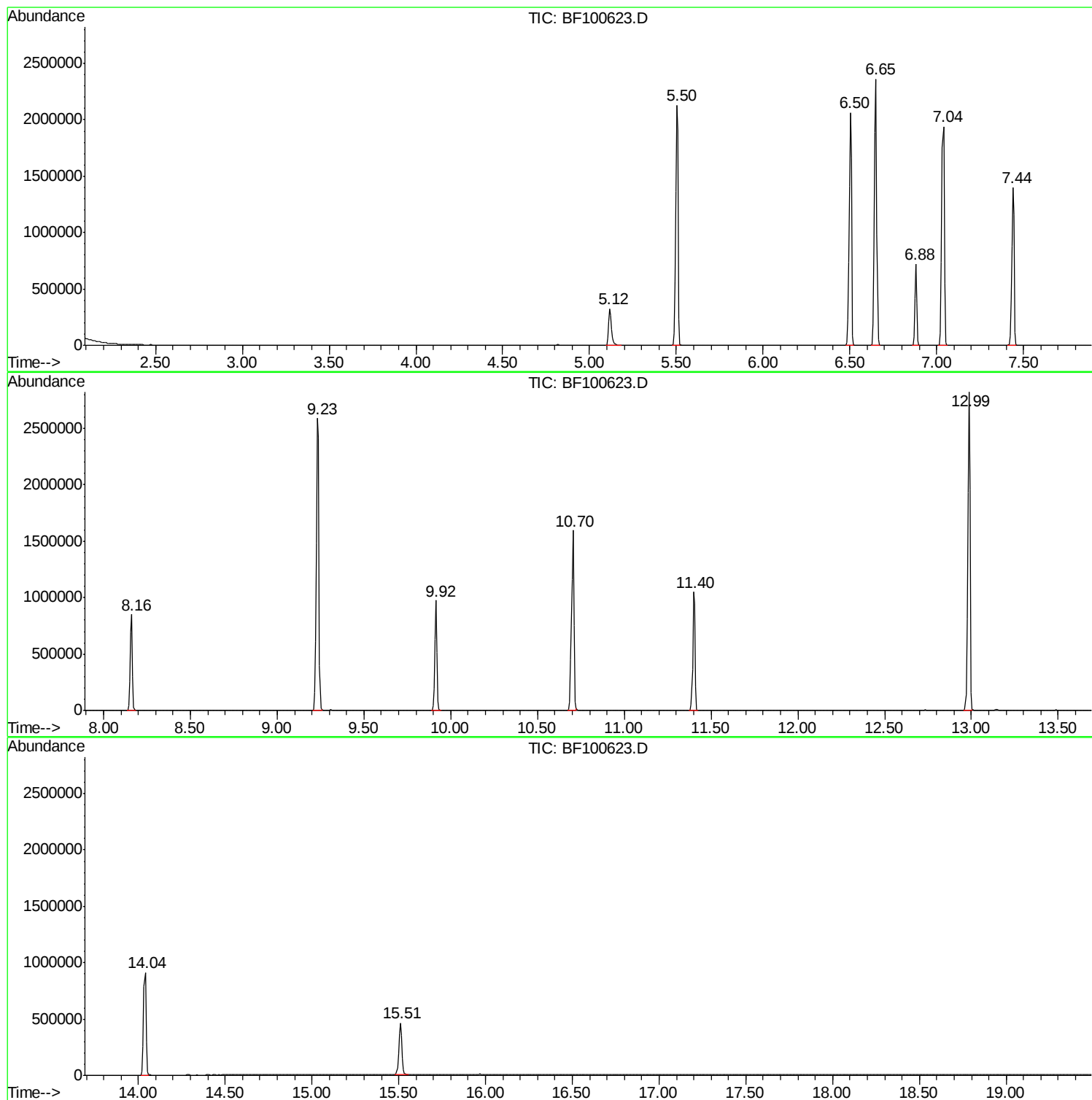
Sum of corrected areas: 21746142

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



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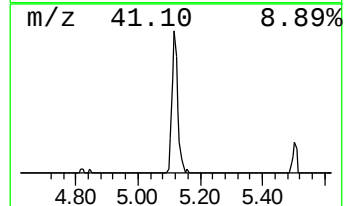
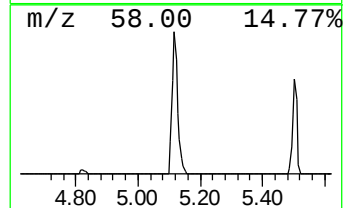
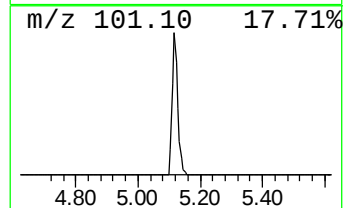
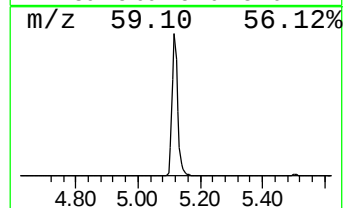
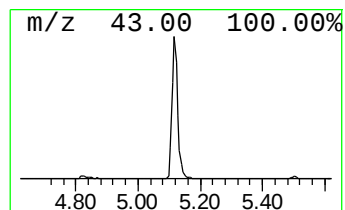
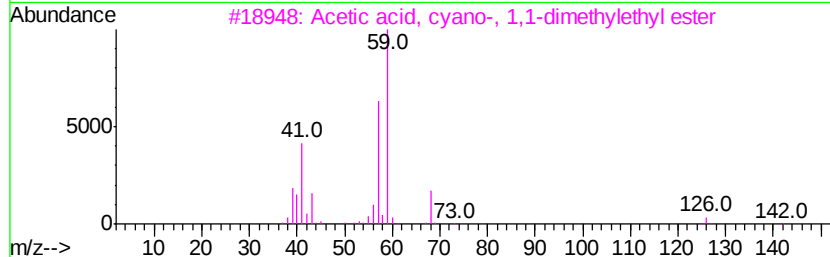
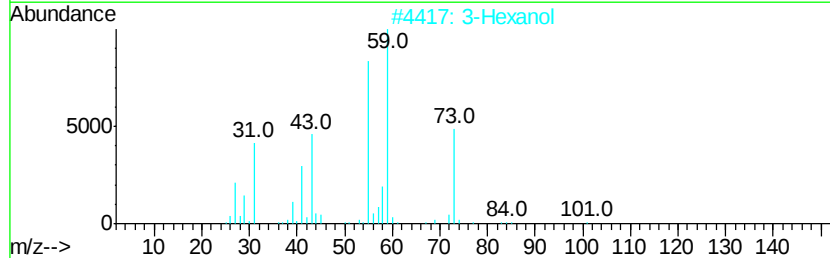
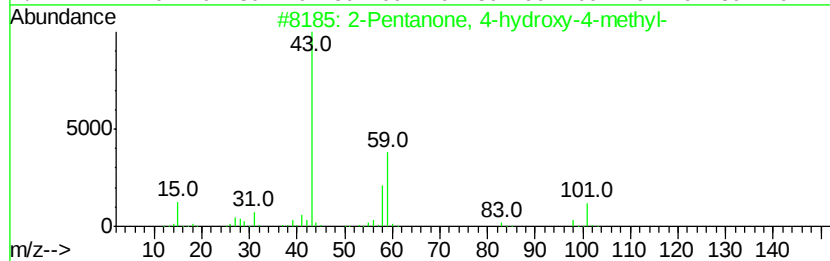
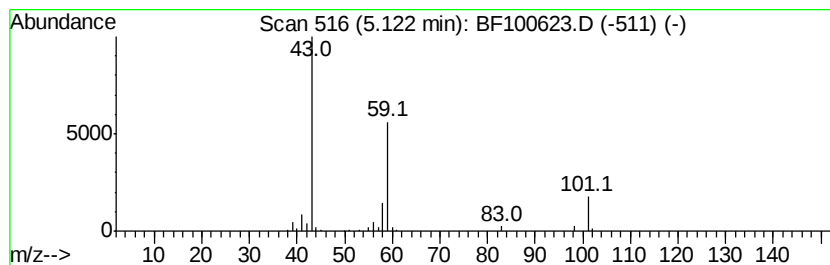
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.12	14.11 ng	404118	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	64
2			3-Hexanol	102	C6H14O	000623-37-0	33
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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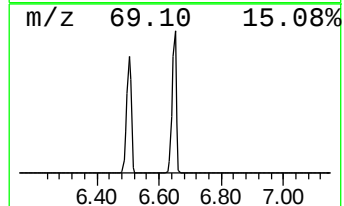
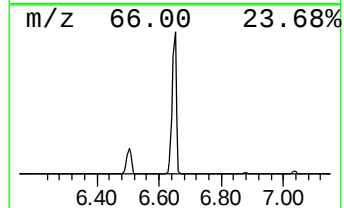
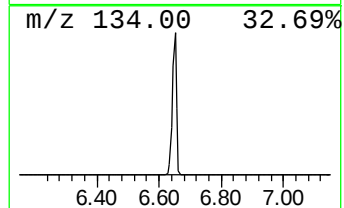
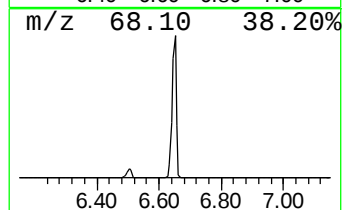
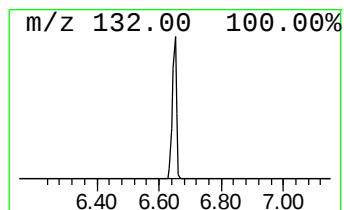
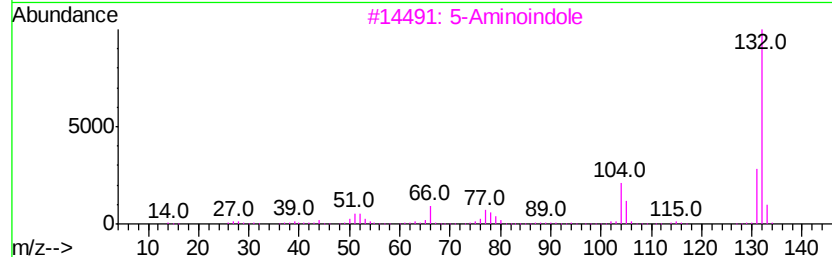
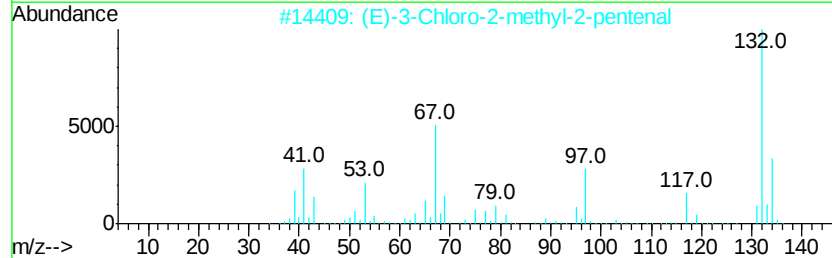
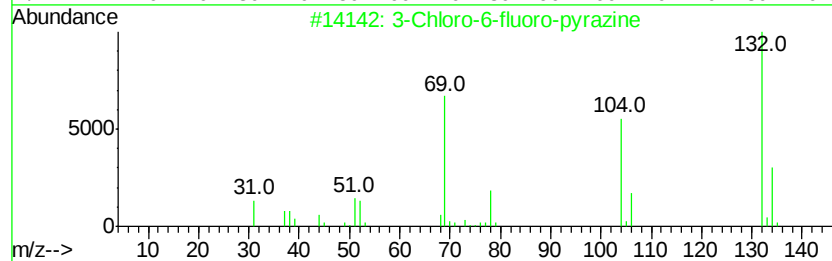
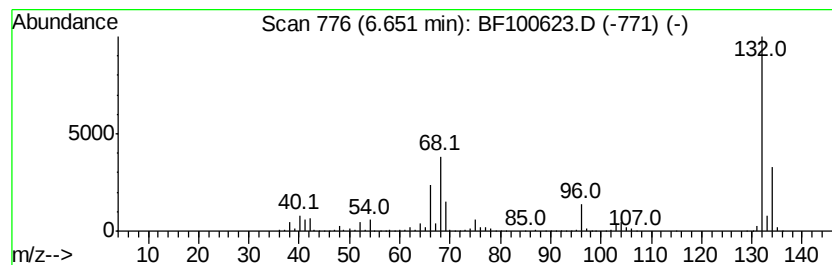
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Peak Number 2 unknown6.65 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.65	78.30 ng	2242600	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	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Amphetaminil	250	C17H18N2	017590-01-1	10
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.12	14.1	ng	404118	1	6.88	572801	20.0
unknown6.65	6.65	78.3	ng	2242600	1	6.88	572801	20.0