

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031517\
 Data File : BF093697.D
 Acq On : 16 Mar 2017 7:24
 Operator : SJ/MA
 Sample : I2167-02
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 031117-DSDC-F-D-M

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-BF030717.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	4.836	256	258	271	rBV	482280	787167	19.23%	2.989%
2	5.316	298	300	322	rBV	635101	1538521	37.58%	5.842%
3	6.379	391	393	400	rBV3	478398	1030350	25.17%	3.912%
4	6.481	400	402	414	rVB	2266932	2657786	64.92%	10.092%
5	6.710	419	422	430	rBV	671914	954906	23.32%	3.626%
6	6.859	433	435	438	rBV	3458775	3331488	81.37%	12.650%
7	7.282	470	472	486	rBV	2498202	2856374	69.77%	10.846%
8	8.002	532	535	537	rBV	1043739	1101805	26.91%	4.184%
9	9.076	627	629	632	rBV	3273327	4094081	100.00%	15.546%
10	9.488	663	665	668	rBV	93584	89567	2.19%	0.340%
11	9.750	686	688	691	rBV	758019	1025541	25.05%	3.894%
12	10.276	732	734	739	rVB	46536	52397	1.28%	0.199%
13	10.550	756	758	766	rBV	1064300	1177030	28.75%	4.469%
14	10.665	766	768	770	rVB	30419	42825	1.05%	0.163%
15	11.248	817	819	821	rBV	677113	842461	20.58%	3.199%
16	12.642	939	941	943	rBV	64172	56569	1.38%	0.215%
17	12.825	954	957	959	rBV	3138762	2977727	72.73%	11.307%
18	13.351	1000	1003	1005	rBV	34594	42462	1.04%	0.161%
19	13.877	1047	1049	1052	rBV	638584	864073	21.11%	3.281%
20	15.282	1170	1172	1182	rVB	549814	812223	19.84%	3.084%

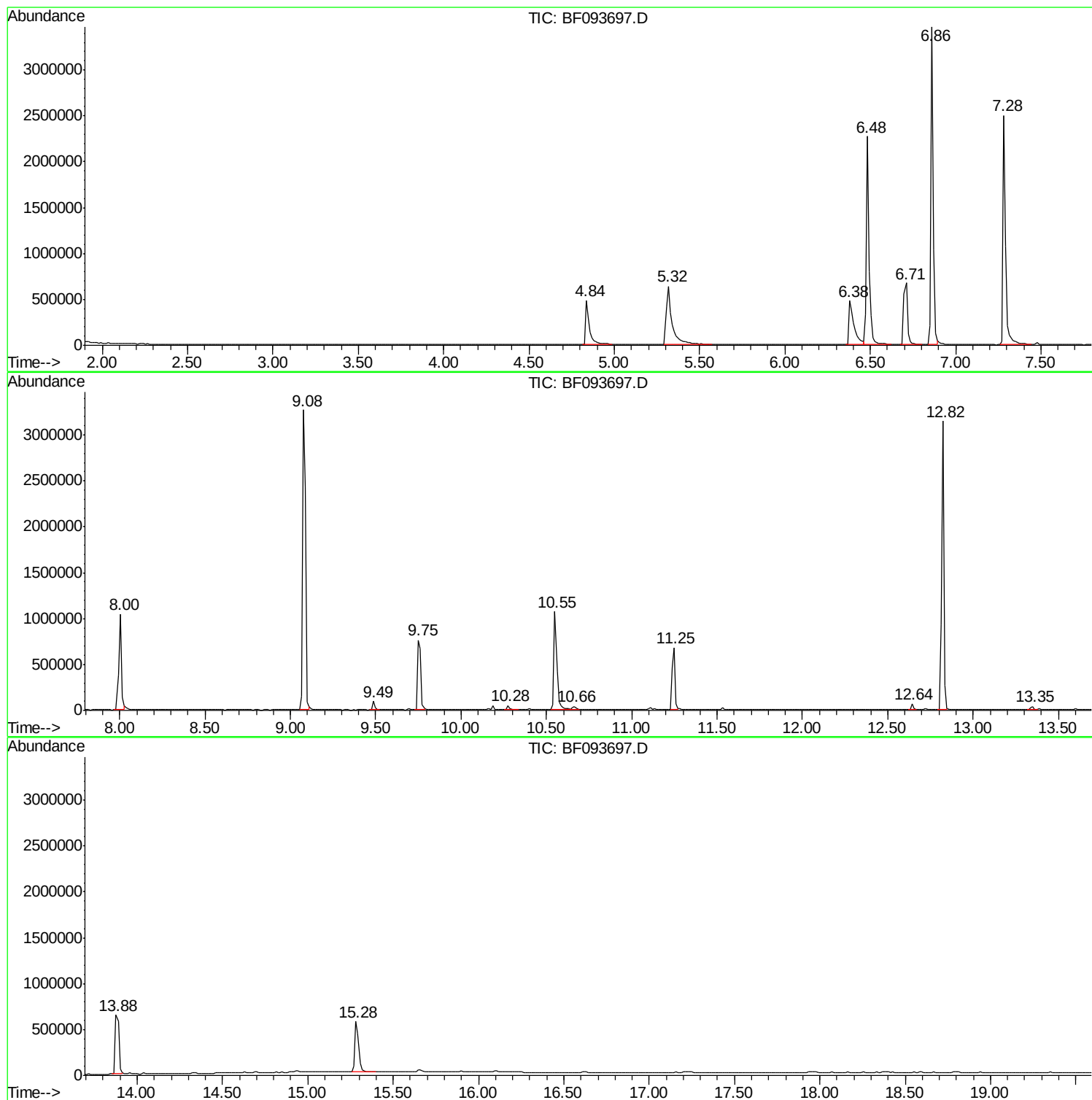
Sum of corrected areas: 26335353

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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF030717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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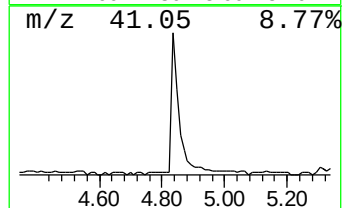
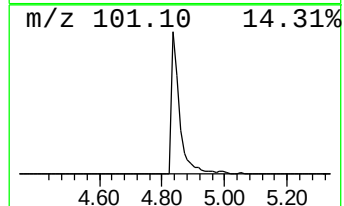
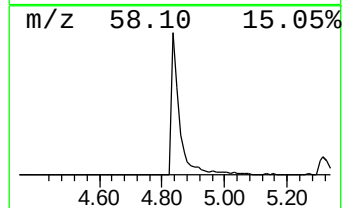
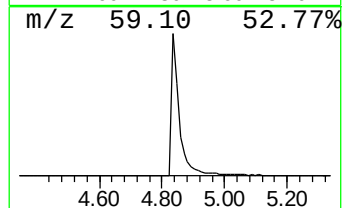
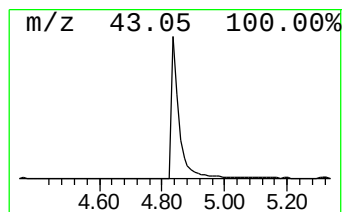
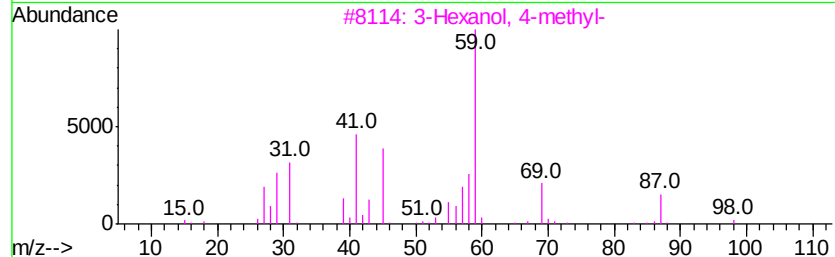
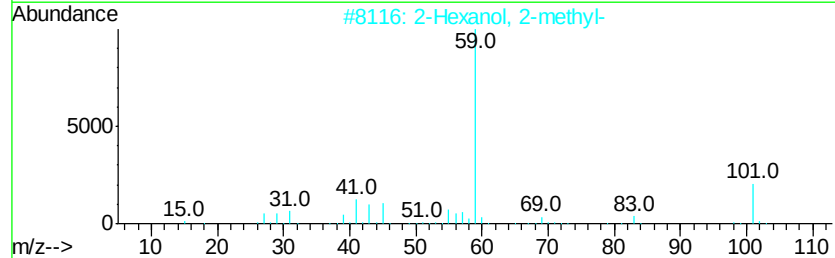
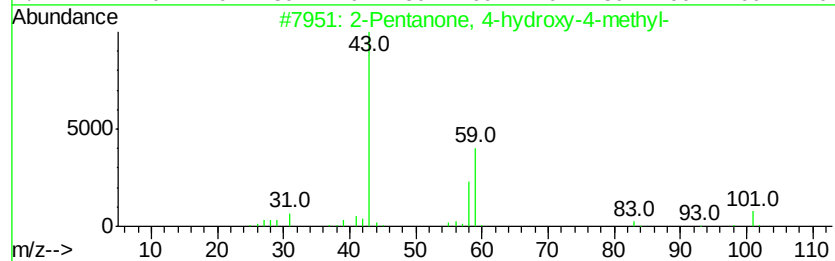
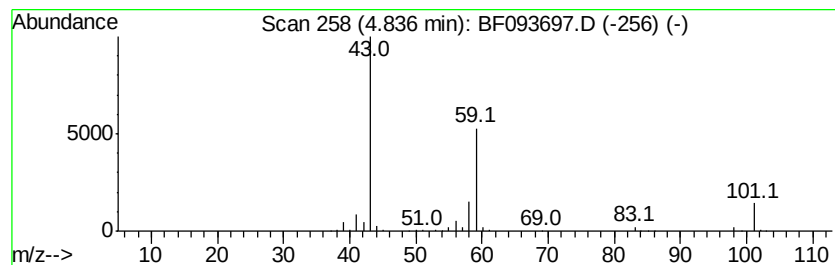
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.84	16.49 ng	787167	1,4-Dichlorobenzene-d4	6.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
5		2-Propanone, 1-(1-methylethoxy)-	116	C6H12O2	042781-12-4	9



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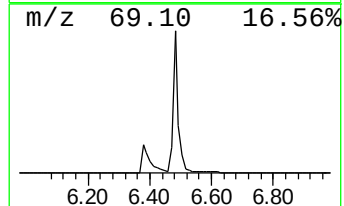
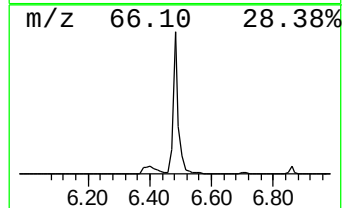
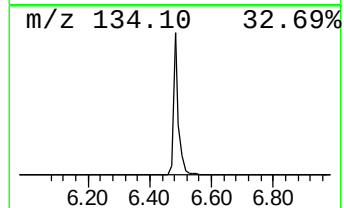
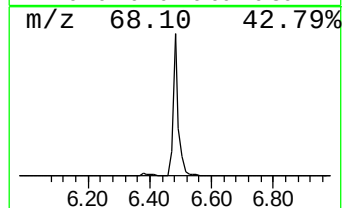
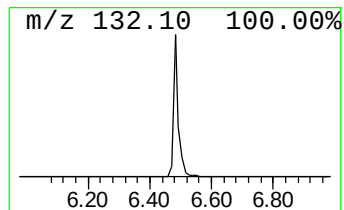
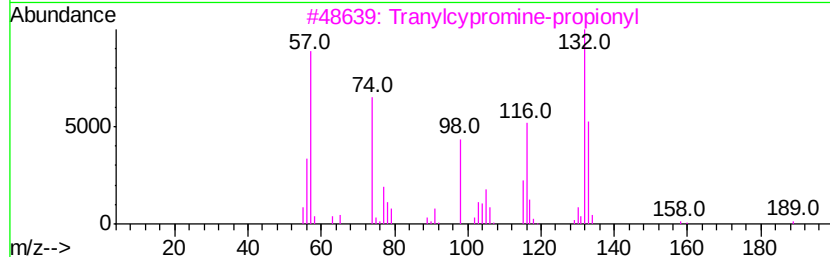
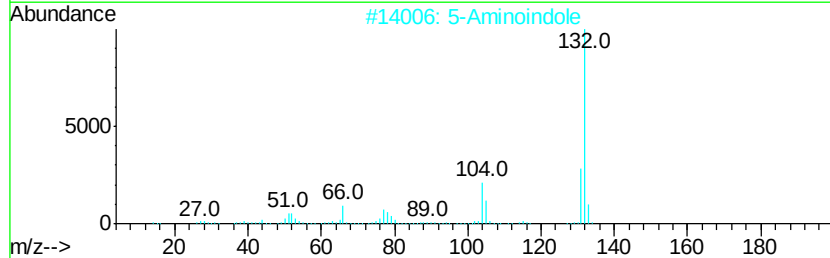
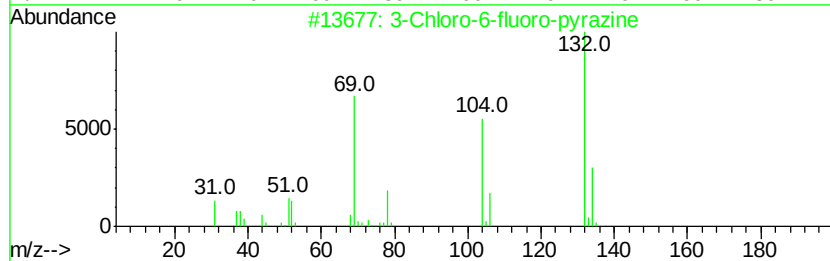
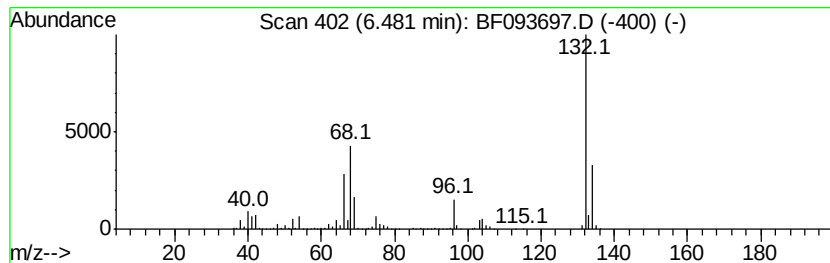
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Peak Number 2 unknown6.48 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.48	55.67 ng	2657790	1,4-Dichlorobenzene-d4	6.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4			Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.84	16.5	ng	787167	1	6.71	954906	20.0
unknown6.48	6.48	55.7	ng	2657790	1	6.71	954906	20.0