

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020916\
 Data File : BF084953.D
 Acq On : 9 Feb 2016 15:46
 Operator : SJ/IZ
 Sample : PB88264BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB88264BL

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-BF020116.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	4.784	233	236	245	rBV	828764	1275933	24.79%	3.119%
2	5.230	272	275	277	rBV	4128298	4270593	82.96%	10.440%
3	6.293	365	368	370	rBV	4005557	4616401	89.68%	11.286%
4	6.407	375	378	380	rBV	3741279	4521463	87.84%	11.053%
5	6.636	395	398	400	rBV	724861	853708	16.58%	2.087%
6	6.785	409	411	414	rBV	2866541	3224094	62.63%	7.882%
7	7.207	445	448	450	rBV	2662956	2988085	58.05%	7.305%
8	7.916	508	510	513	rBV	1168402	1151752	22.37%	2.816%
9	9.002	601	605	607	rBV	4920687	5044548	98.00%	12.332%
10	9.665	660	663	666	rVB	1391776	1314962	25.55%	3.215%
11	10.453	729	732	735	rBV	2925381	3026602	58.80%	7.399%
12	10.785	759	761	762	rBV	111403	99083	1.92%	0.242%
13	10.808	762	763	766	rVB	91718	75818	1.47%	0.185%
14	11.139	790	792	795	rBV	1212000	1295460	25.17%	3.167%
15	12.739	929	932	934	rBV	3781068	5147578	100.00%	12.584%
16	13.768	1020	1022	1025	rBV	1252386	1168869	22.71%	2.857%
17	15.128	1139	1141	1144	rBV	679267	830377	16.13%	2.030%

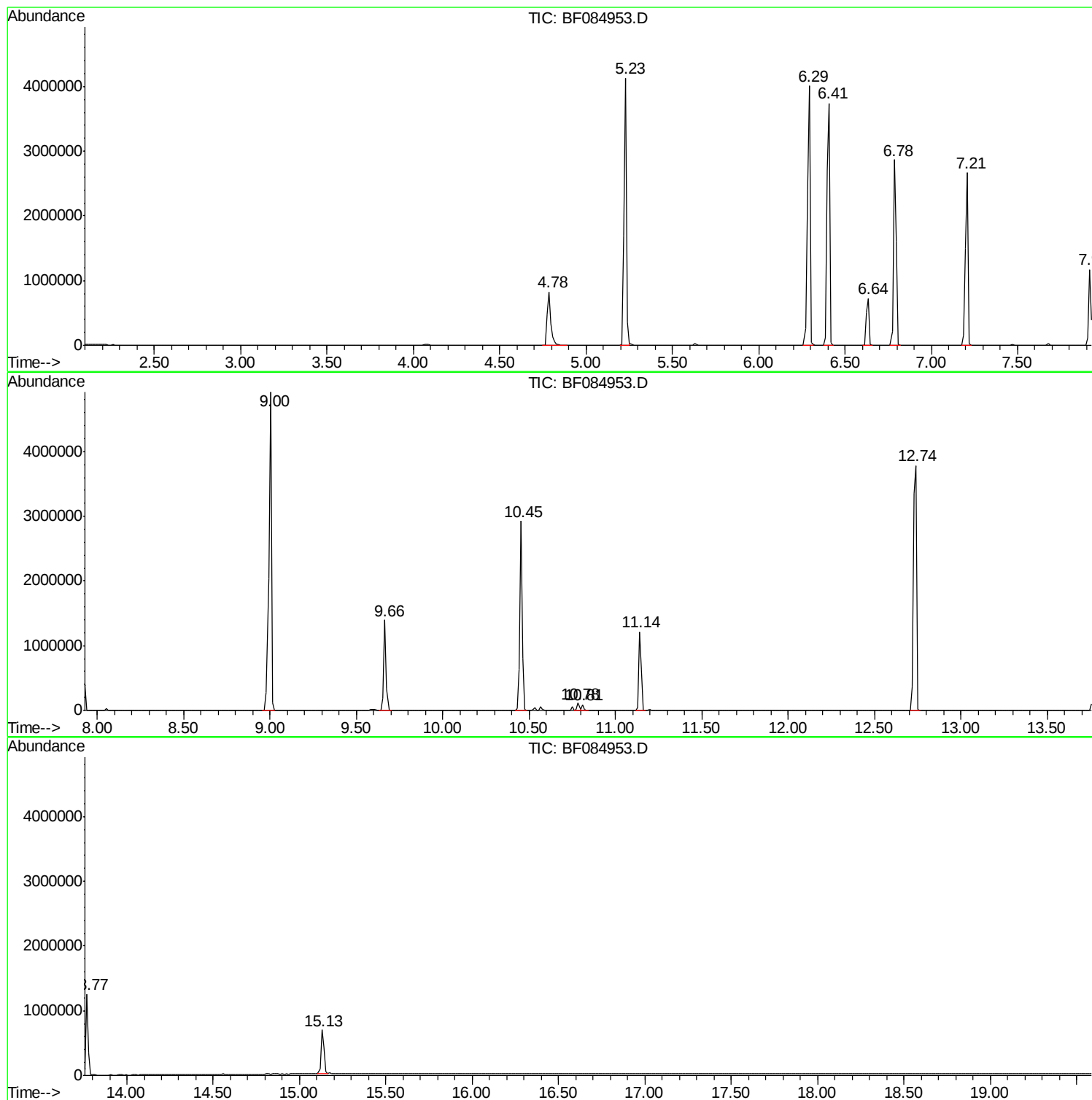
Sum of corrected areas: 40905326

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020116.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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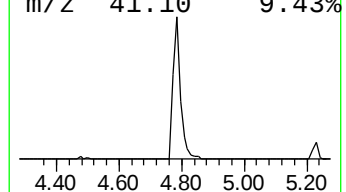
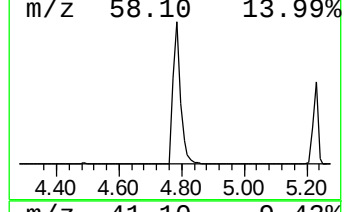
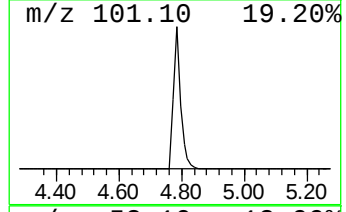
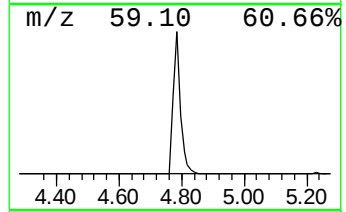
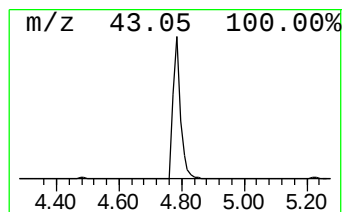
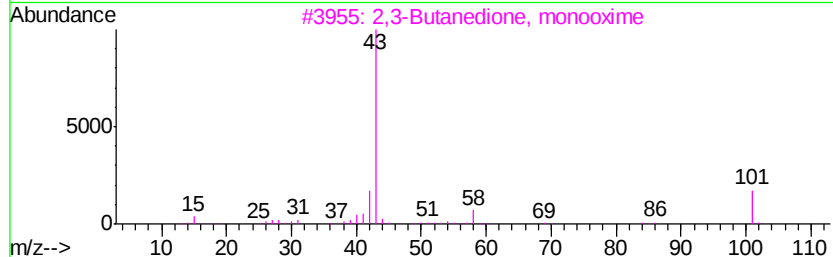
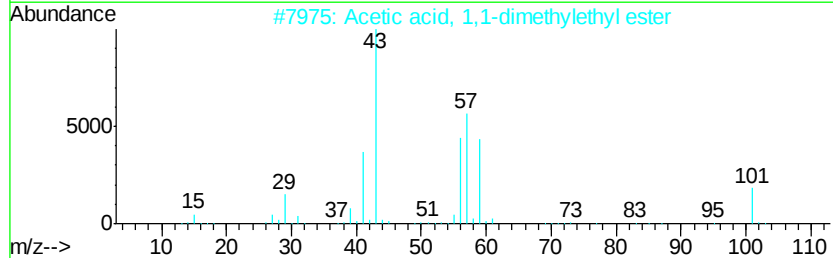
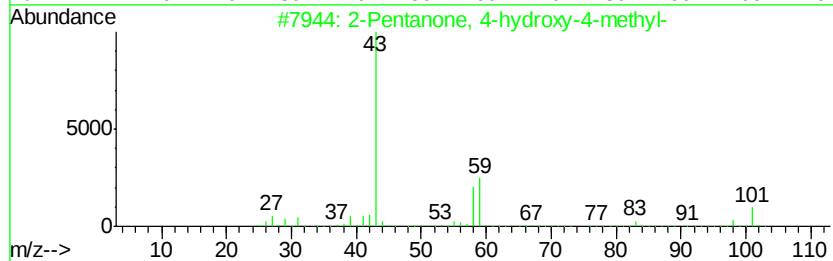
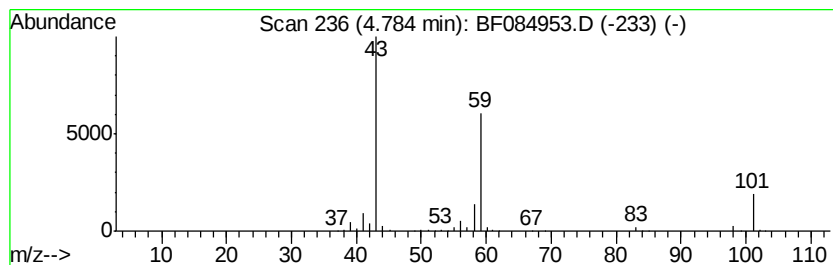
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.78	29.89 ng	1275930	1,4-Dichlorobenzene-d4	6.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	9
5		tert-Butyl carbamate	117	C5H11NO2	004248-19-5	9



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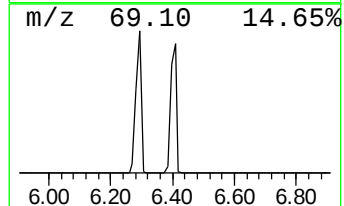
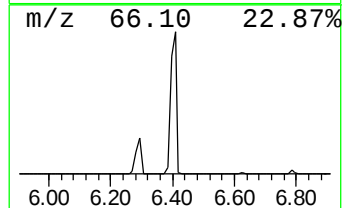
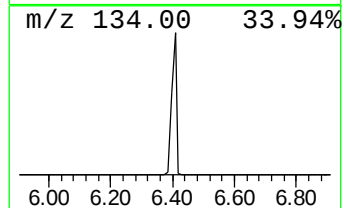
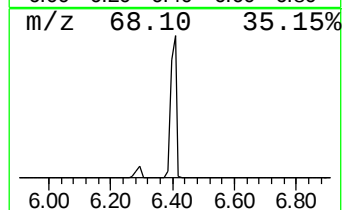
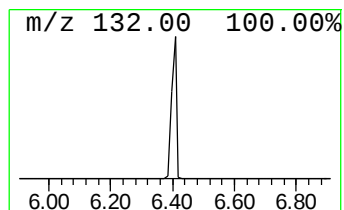
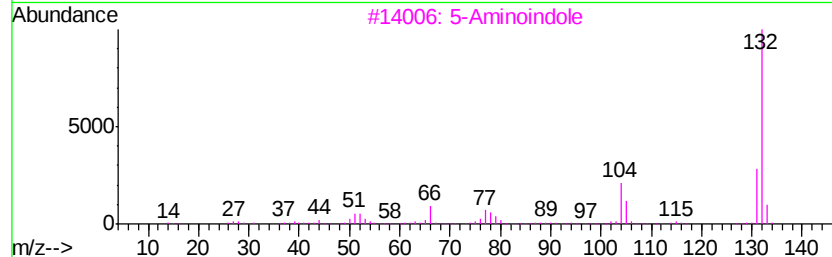
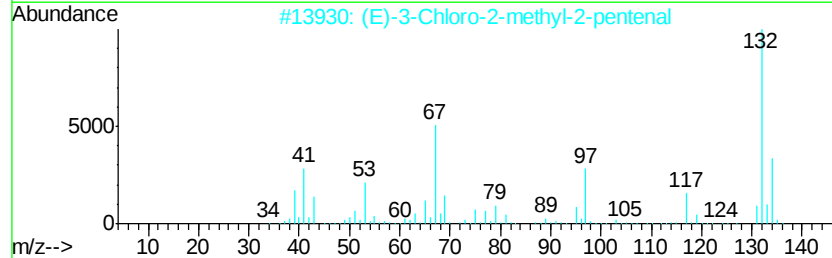
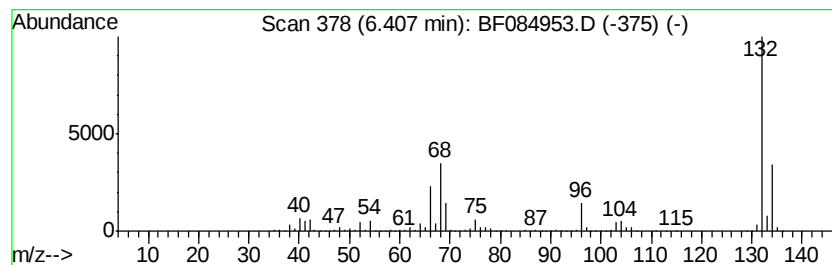
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Peak Number 2 unknown6.41 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.41	105.93 ng	4521460	1,4-Dichlorobenzene-d4	6.64

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		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.78	29.9	ng	1275930	1	6.64	853708	20.0
unknown6.41	6.41	105.9	ng	4521460	1	6.64	853708	20.0