

Data Path : Z:\HPCHEM1\BNA F\DATA\BF110916\
 Data File : BF091124.D
 Acq On : 9 Nov 2016 15:18
 Operator : UM/SJ
 Sample : PB94612BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB94612BL

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-BF110316.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	1.744	4	5	65	rVB	3848693	12931850	100.00%	20.681%
2	4.807	270	273	286	rBV	607608	1021887	7.90%	1.634%
3	5.253	309	312	314	rBV	4499409	5033915	38.93%	8.050%
4	6.316	402	405	407	rBV	3420397	5342280	41.31%	8.544%
5	6.419	412	414	417	rBV	3674841	5185389	40.10%	8.293%
6	6.647	432	434	437	rBV	932420	1142338	8.83%	1.827%
7	6.807	445	448	451	rBV	4048749	3962155	30.64%	6.336%
8	7.230	481	485	487	rBV	2246487	3450144	26.68%	5.518%
9	7.939	544	547	550	rBV	1720135	1639989	12.68%	2.623%
10	9.025	638	642	644	rBV	5449919	6300169	48.72%	10.075%
11	9.688	697	700	703	rBV	1963571	1826546	14.12%	2.921%
12	10.476	766	769	772	rBV2	2751417	3758267	29.06%	6.010%
13	11.173	827	830	832	rBV	1288928	1691015	13.08%	2.704%
14	12.762	966	969	971	rBV	5474345	6116261	47.30%	9.781%
15	13.802	1057	1060	1062	rBV	1516637	1702839	13.17%	2.723%
16	15.174	1177	1180	1183	rBV	1229359	1424771	11.02%	2.279%

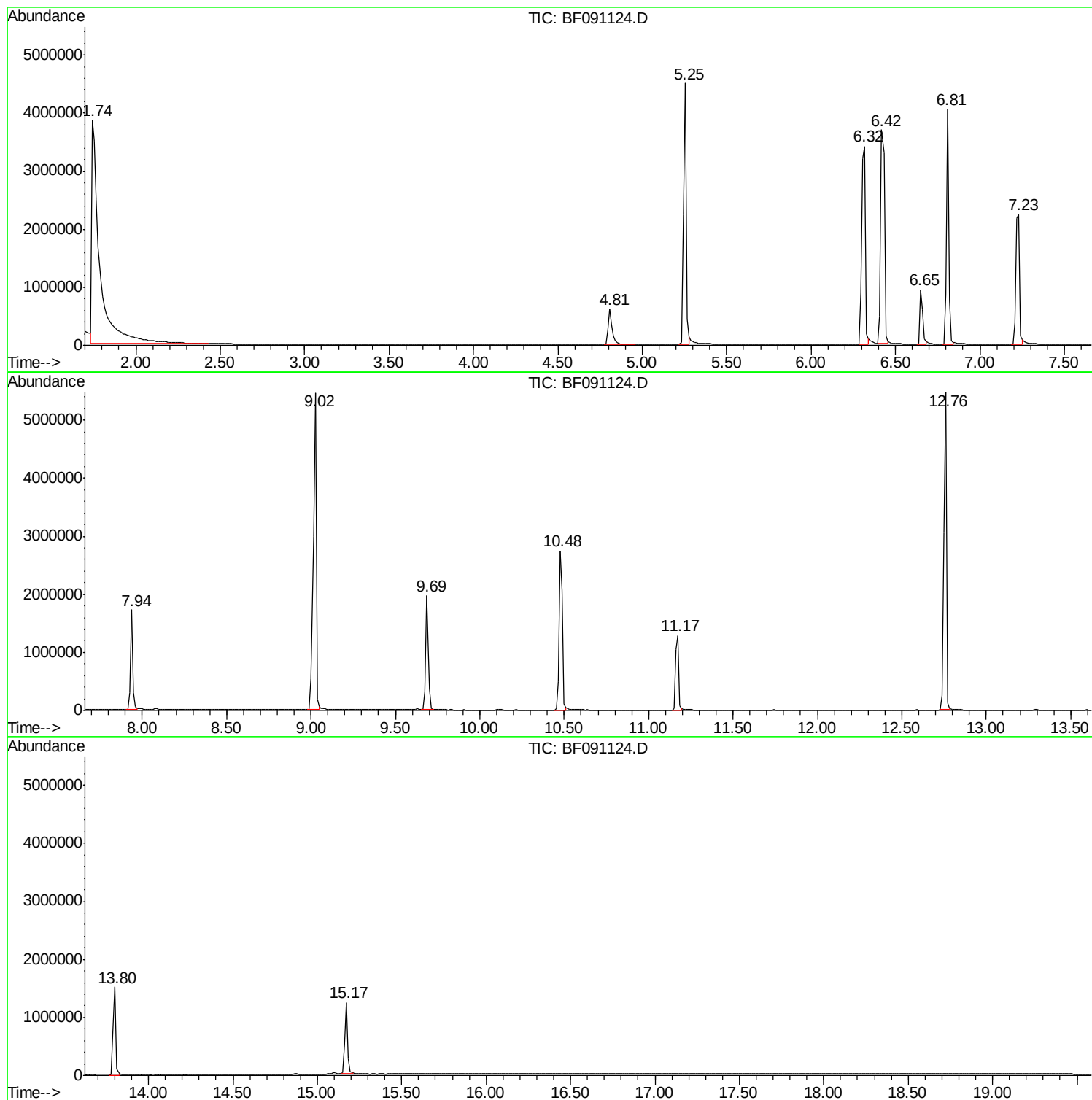
Sum of corrected areas: 62529815

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110316.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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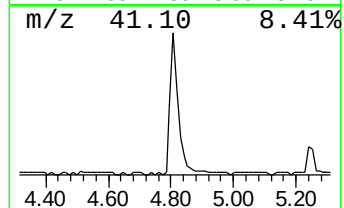
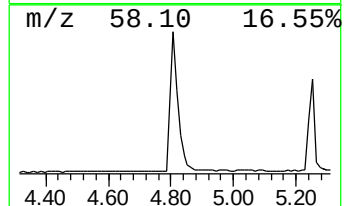
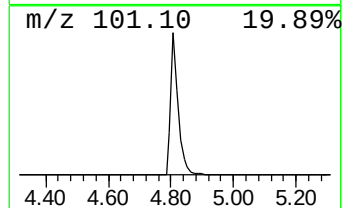
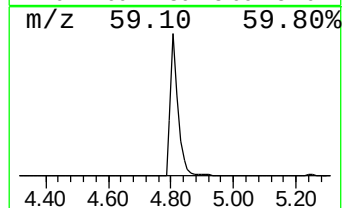
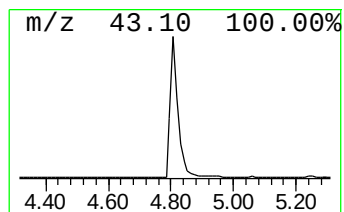
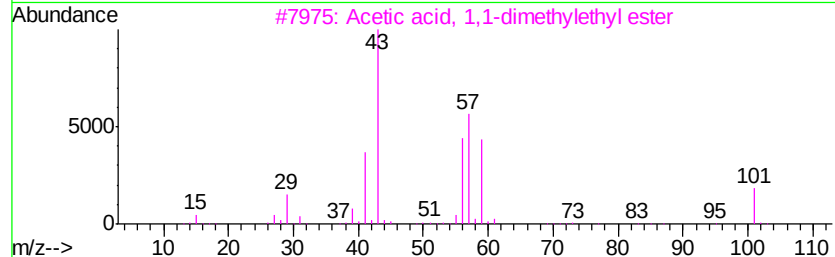
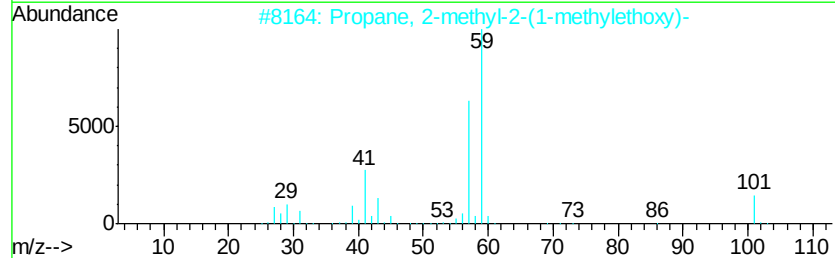
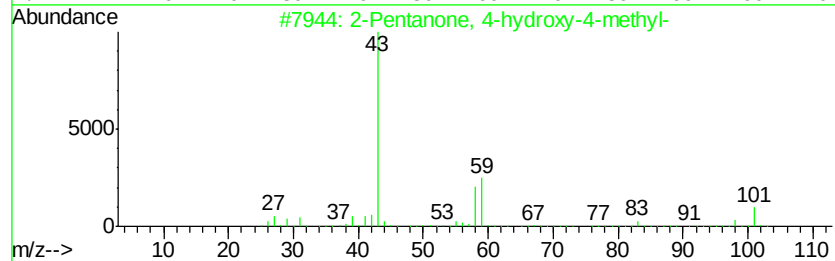
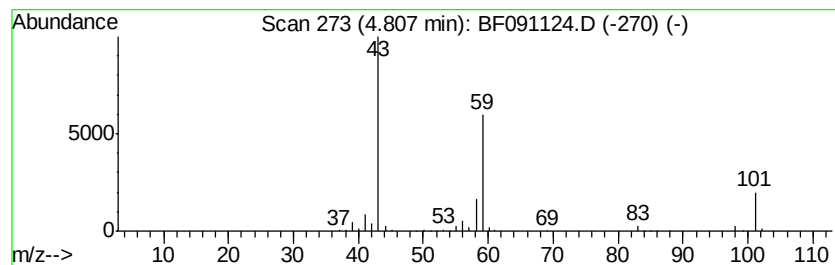
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.81	17.89 ng	1021890	1,4-Dichlorobenzene-d4	6.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	42
3		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
4		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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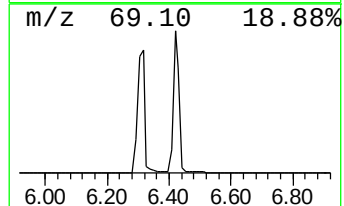
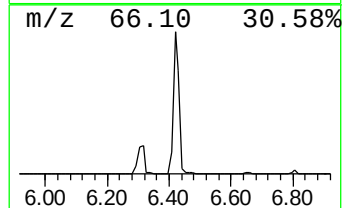
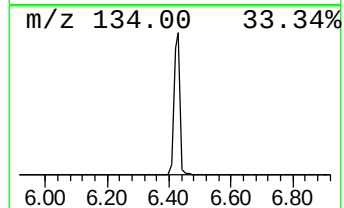
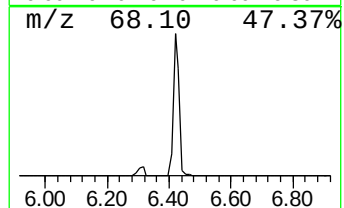
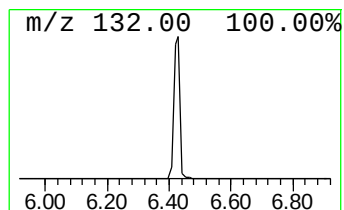
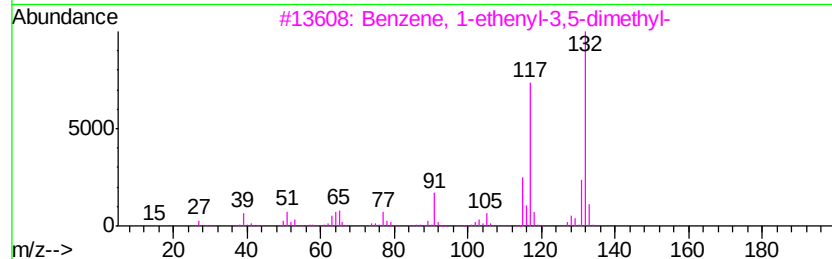
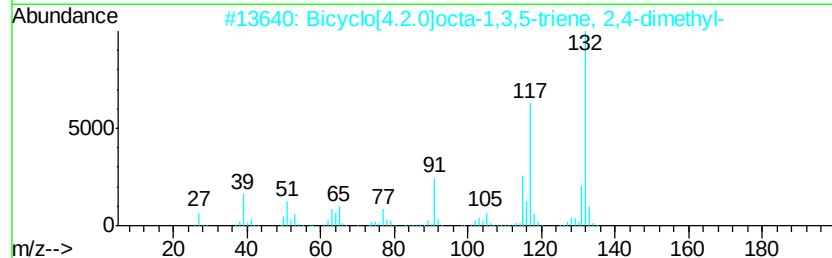
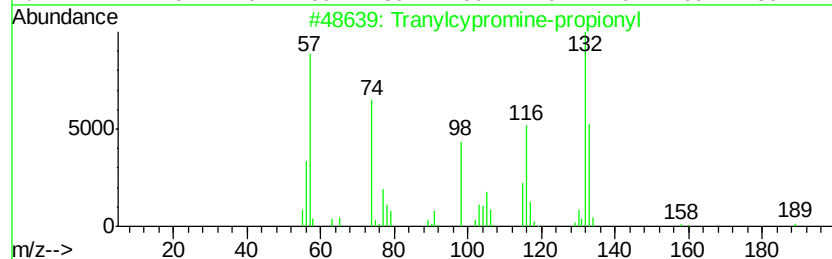
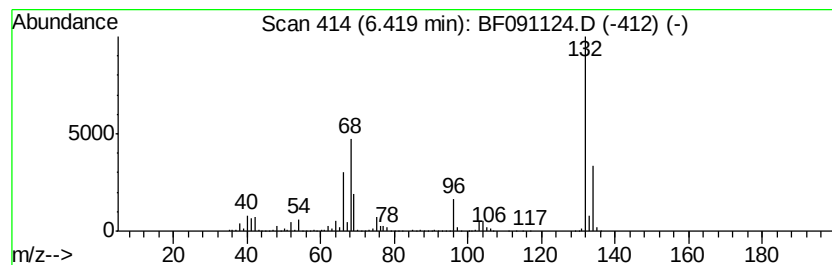
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Peak Number 3 unknown6.42 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.42	90.79 ng	5185390	1,4-Dichlorobenzene-d4	6.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
2		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
3		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
4		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.81	17.9	ng	1021890	1	6.65	1142340	20.0
unknown6.42	6.42	90.8	ng	5185390	1	6.65	1142340	20.0