

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063015\
 Data File : BF080239.D
 Acq On : 30 Jun 2015 16:36
 Operator : TP/IZ
 Sample : G2815-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SWD-1

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-BF061715.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	5.499	208	211	213	rBV	173588	252191	14.57%	2.913%
2	5.887	243	245	248	rBV	349139	414251	23.93%	4.785%
3	6.882	330	332	335	rBV	278420	305331	17.64%	3.527%
4	7.053	345	347	349	rBV	620185	624095	36.05%	7.209%
5	7.282	365	367	369	rBV	245568	210088	12.14%	2.427%
6	7.442	379	381	383	rVB	1046445	838465	48.44%	9.685%
7	7.842	414	416	418	rBV	830474	672595	38.85%	7.769%
8	8.573	478	480	482	rBV	344777	280712	16.22%	3.242%
9	9.648	572	574	577	rVB	1457615	1357110	78.40%	15.676%
10	10.345	633	635	638	rVB	410167	340336	19.66%	3.931%
11	11.133	702	704	707	rBV2	444137	531548	30.71%	6.140%
12	11.854	764	767	769	rBV	314101	356359	20.59%	4.116%
13	13.522	910	913	915	rBV	1621277	1731063	100.00%	19.995%
14	14.768	1020	1022	1025	rBV	393977	383106	22.13%	4.425%
15	16.608	1180	1183	1193	rBV	246349	360058	20.80%	4.159%

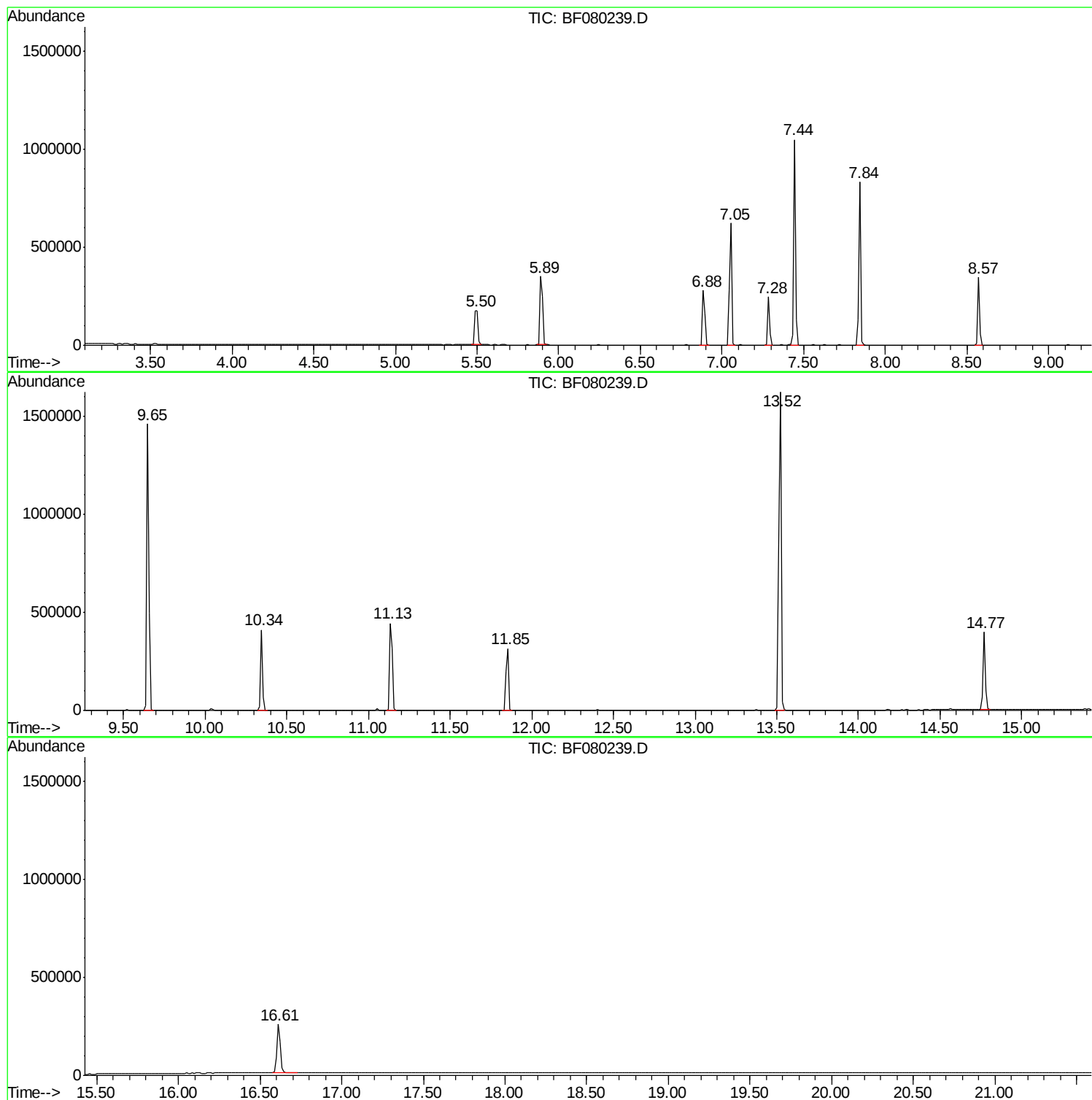
Sum of corrected areas: 8657308

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.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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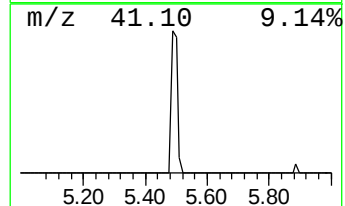
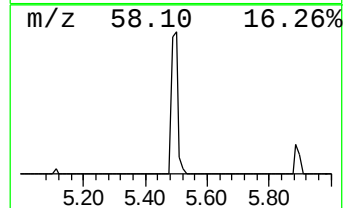
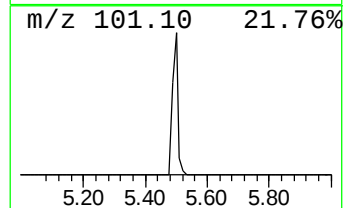
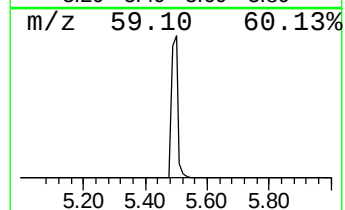
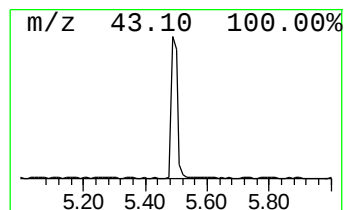
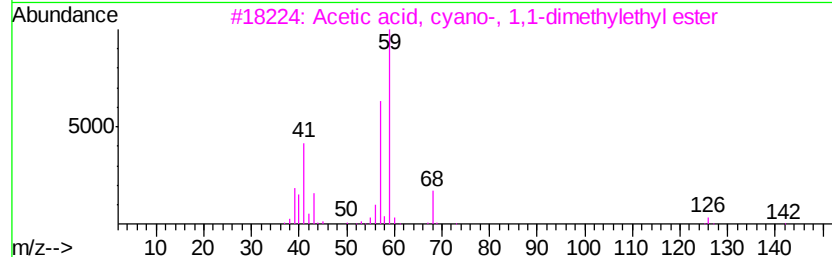
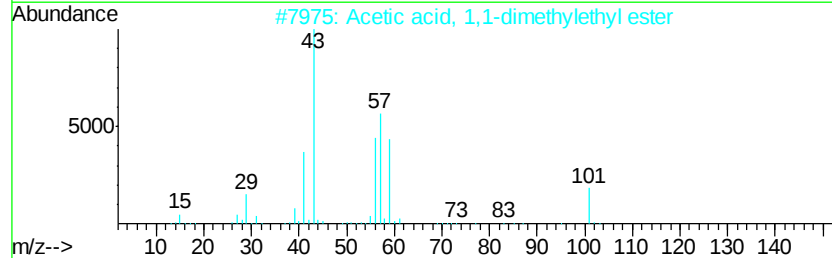
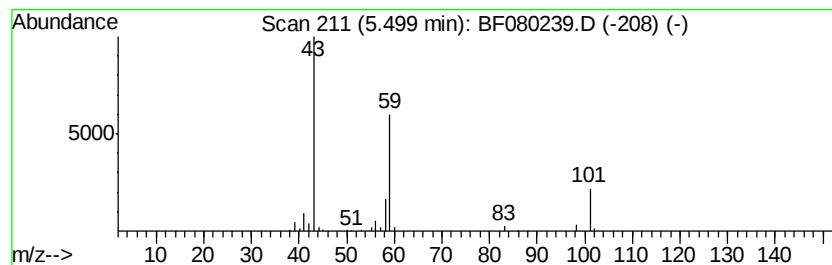
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.50	24.01 ng	252191	1,4-Dichlorobenzene-d4	7.28

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
4			5-Hexen-2-one	98	C6H10O	000109-49-9	9
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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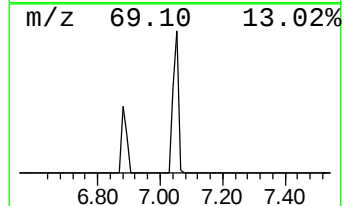
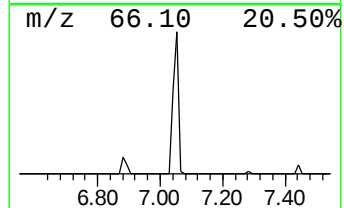
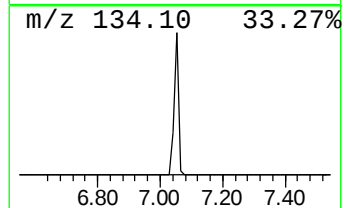
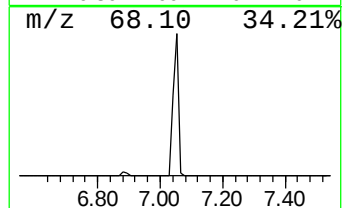
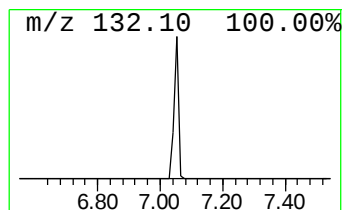
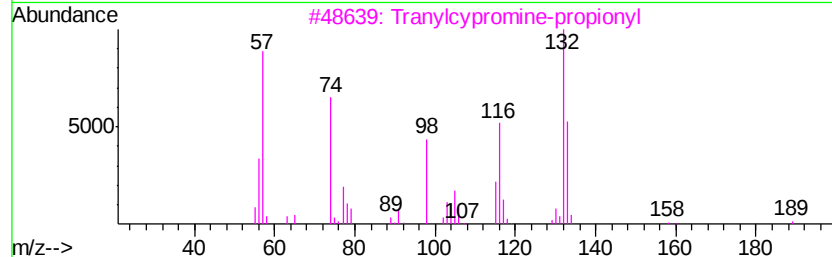
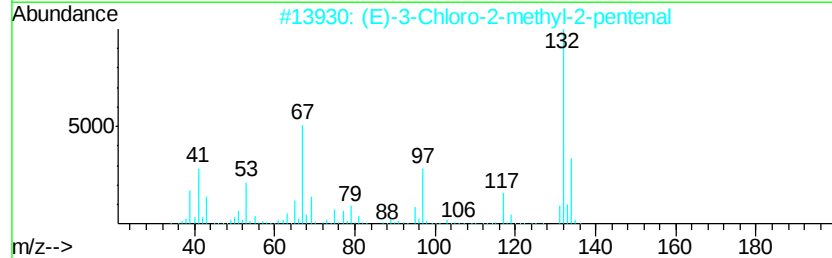
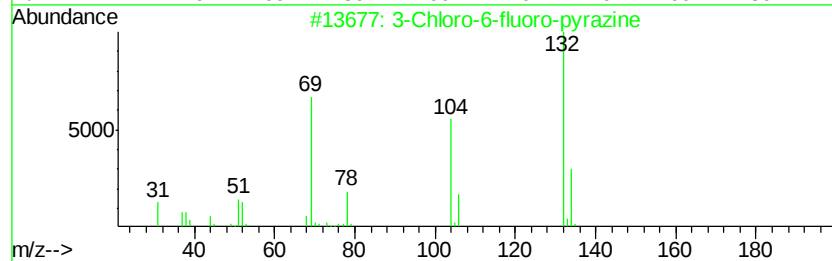
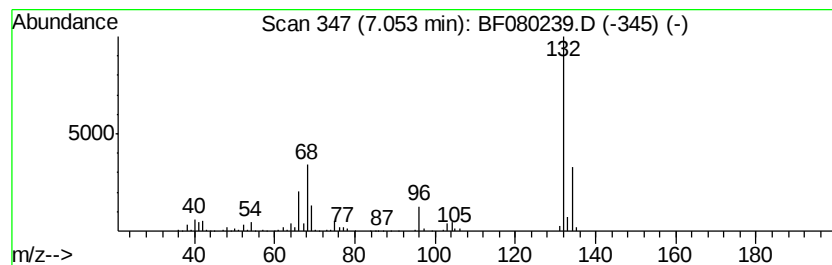
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Peak Number 2 unknown7.05 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.05	59.41 ng	624095	1,4-Dichlorobenzene-d4	7.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		Amphetaminil	250	C17H18N2	017590-01-1	12
5		5-Aminoindole	132	C8H8N2	005192-03-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.50	24.0	ng	252191	1	7.28	210088	20.0
unknown7.05	7.05	59.4	ng	624095	1	7.28	210088	20.0