

Data Path : Z:\HPCHEM1\BNA F\DATA\BF082416\
 Data File : BF089868.D
 Acq On : 24 Aug 2016 12:30
 Operator : UM/SJ
 Sample : H4551-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AST-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-BF081916.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	1.655	4	6	14	rBV	1307967	2388542	9.36%	2.048%
2	1.770	14	16	58	rVB	9607830	25509037	100.00%	21.875%
3	4.079	215	218	222	rBV	1119006	1591213	6.24%	1.365%
4	4.799	277	281	289	rBV	2082758	3166222	12.41%	2.715%
5	5.244	317	320	322	rBV	7266447	8239595	32.30%	7.066%
6	6.307	409	413	415	rBV	8163978	9240968	36.23%	7.924%
7	6.433	421	424	426	rBV	9113724	9283231	36.39%	7.961%
8	6.662	442	444	447	rBV	2211253	2145602	8.41%	1.840%
9	6.822	455	458	460	rBV	8373424	7085592	27.78%	6.076%
10	7.233	491	494	496	rBV	5917683	5640640	22.11%	4.837%
11	7.953	554	557	559	rBV	3432145	2941453	11.53%	2.522%
12	9.039	649	652	654	rBV	8965697	10121959	39.68%	8.680%
13	9.428	684	686	689	rBV	2156029	1970030	7.72%	1.689%
14	9.702	708	710	712	rBV	3739577	3241567	12.71%	2.780%
15	10.491	776	779	781	rBV	3431992	4265308	16.72%	3.658%
16	10.628	789	791	796	rVB	276988	307711	1.21%	0.264%
17	11.176	837	839	843	rBV	3701840	3354683	13.15%	2.877%
18	12.777	976	979	981	rBV	6627214	9402119	36.86%	8.063%
19	13.714	1058	1061	1068	rBV	115305	346403	1.36%	0.297%
20	13.817	1068	1070	1078	rVB2	2539686	4018710	15.75%	3.446%
21	15.268	1194	1197	1200	rBV	2092315	2353436	9.23%	2.018%

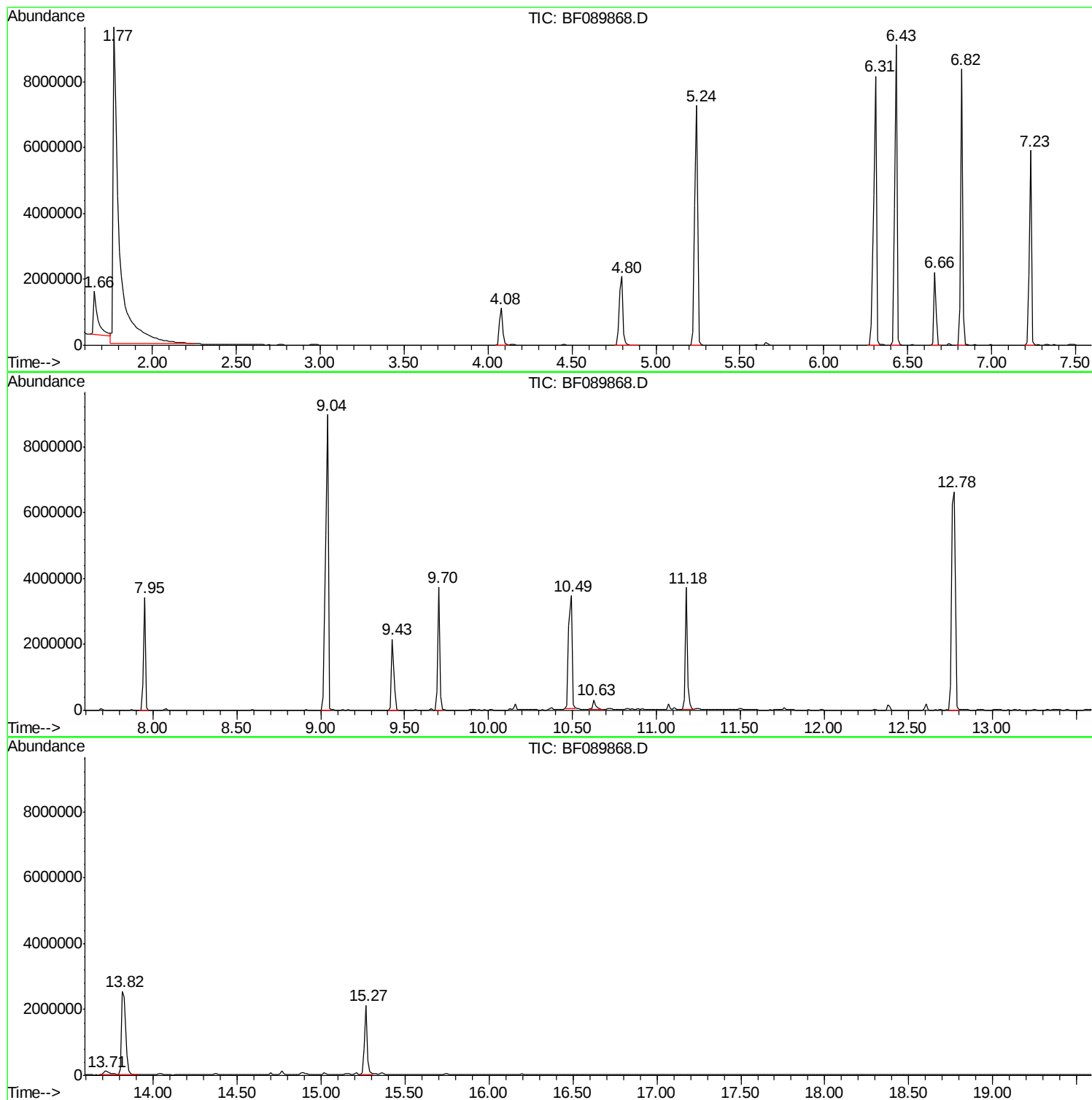
Sum of corrected areas: 116614021

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082416\
Data File : BF089868.D
Acq On : 24 Aug 2016 12:30
Operator : UM/SJ
Sample : H4551-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AST-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082416\
Data File : BF089868.D
Acq On : 24 Aug 2016 12:30
Operator : UM/SJ
Sample : H4551-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AST-1

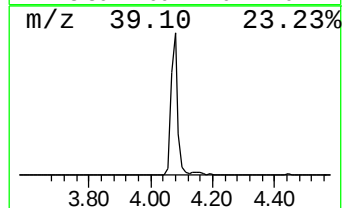
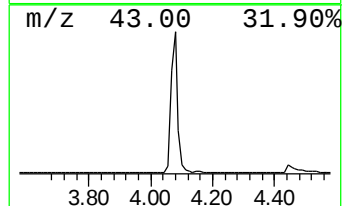
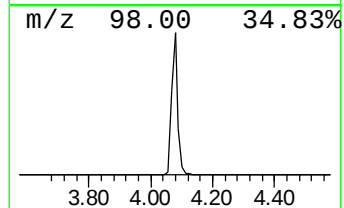
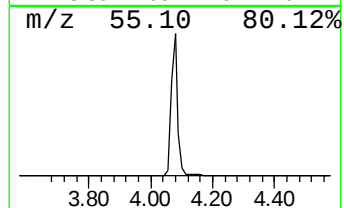
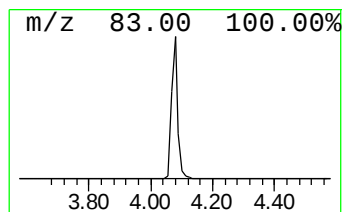
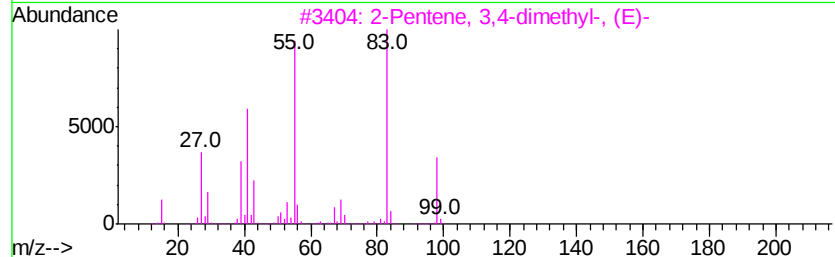
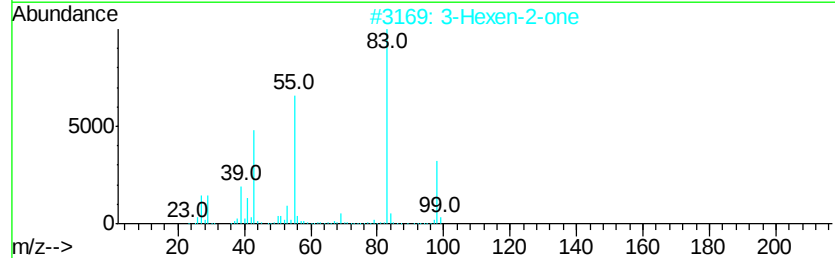
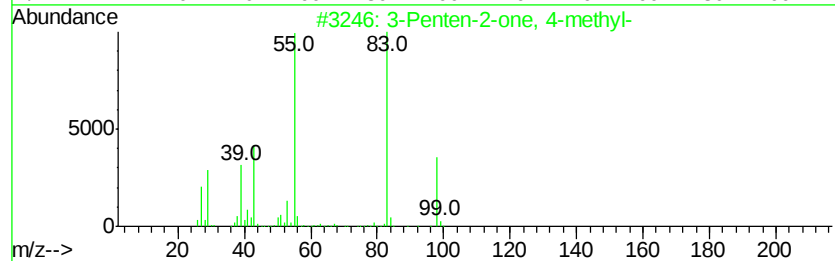
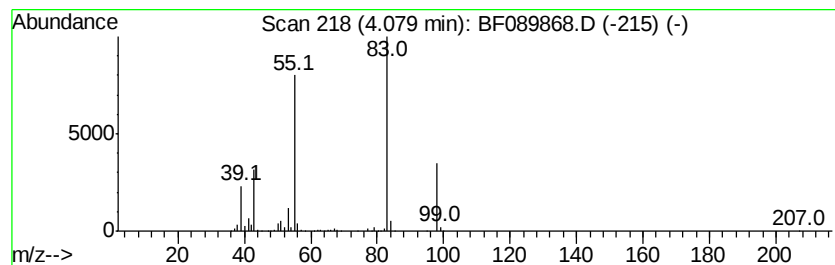
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 3-Penten-2-one, 4-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.08	14.83 ng	1591210	1,4-Dichlorobenzene-d4	6.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	91
2		3-Hexen-2-one	98	C6H10O	000763-93-9	91
3		2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	90
4		2-Pentene, 3,4-dimethyl-, (Z)-	98	C7H14	004914-91-4	86
5		2,4-Azetidinedione, 3,3-diethyl-	141	C7H11N2O2	042282-85-9	78



Data Path : Z:\HPCHEM1\BNA F\DATA\BF082416\
Data File : BF089868.D
Acq On : 24 Aug 2016 12:30
Operator : UM/SJ
Sample : H4551-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AST-1

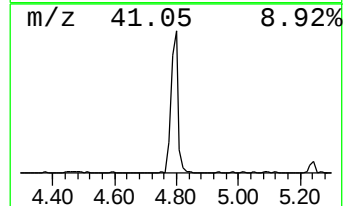
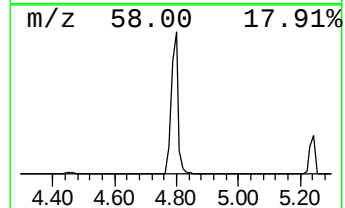
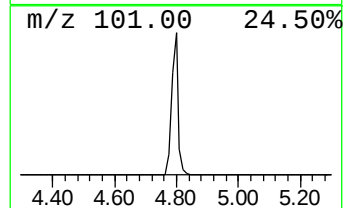
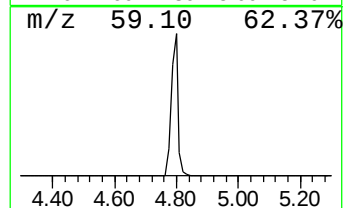
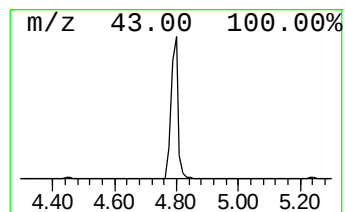
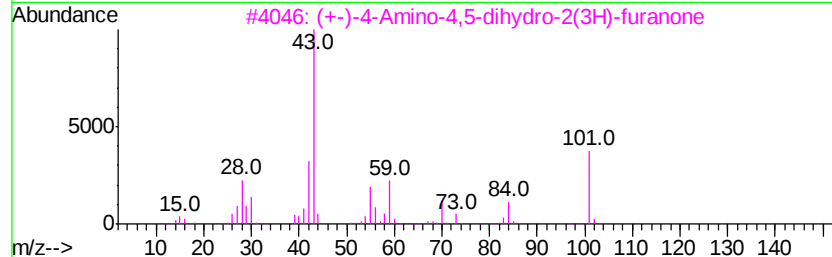
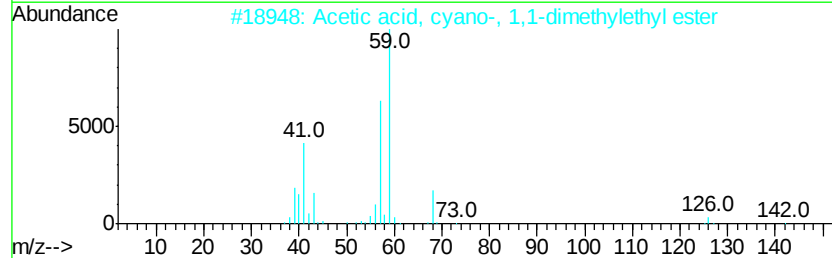
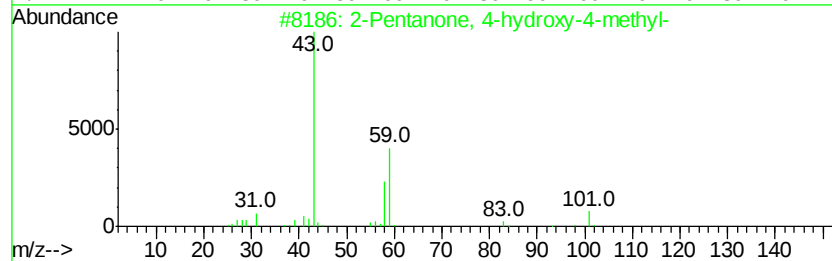
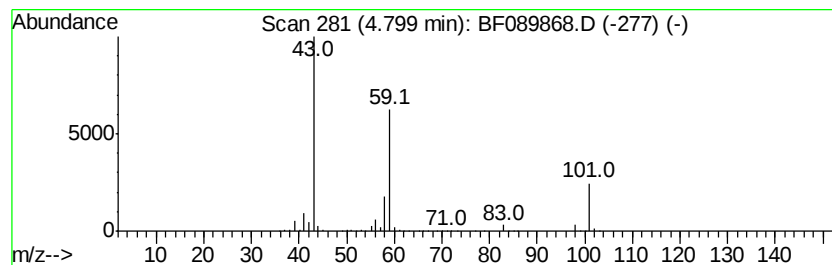
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.80	29.51 ng	3166220	1,4-Dichlorobenzene-d4	6.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
3		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
4		5-Hexen-2-one	98	C6H10O	000109-49-9	9
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082416\
Data File : BF089868.D
Acq On : 24 Aug 2016 12:30
Operator : UM/SJ
Sample : H4551-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AST-1

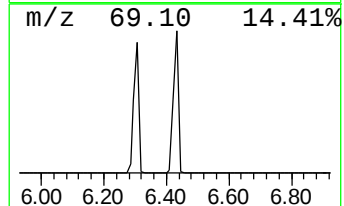
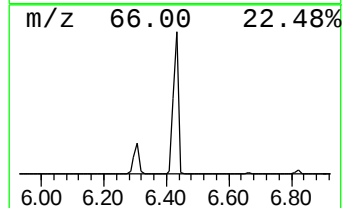
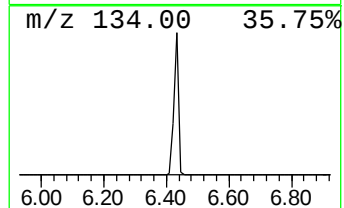
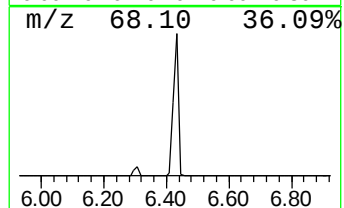
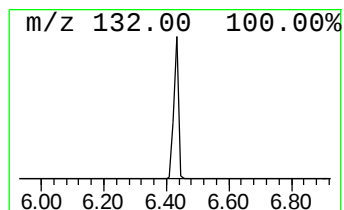
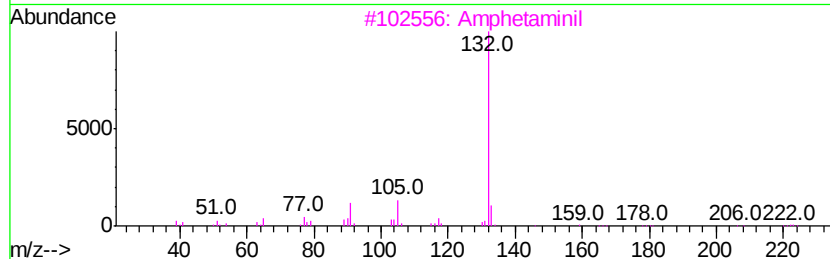
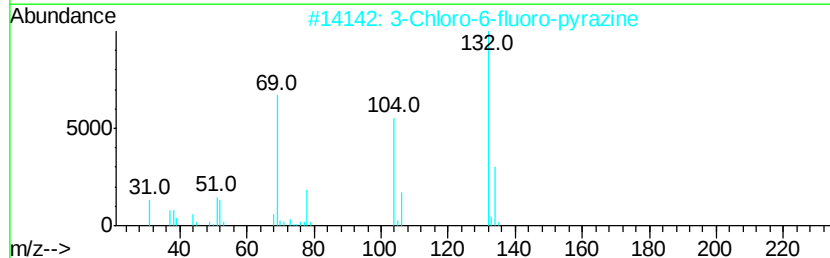
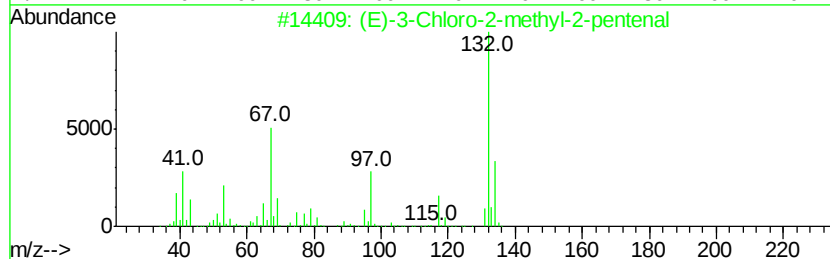
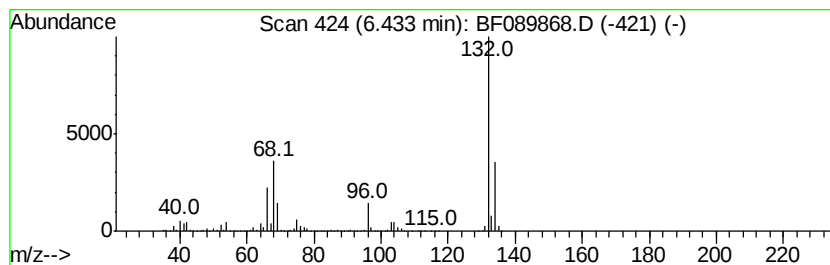
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 unknown6.43 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.43	86.53 ng	9283230	1,4-Dichlorobenzene-d4	6.66

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
2			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
3			Amphetaminil	250	C17H18N2	017590-01-1	12
4			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5			Xenon	132	Xe	007440-63-3	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082416\
Data File : BF089868.D
Acq On : 24 Aug 2016 12:30
Operator : UM/SJ
Sample : H4551-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AST-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3-Penten-2-one, 4...	4.08	14.8	ng	1591210	1	6.66	2145600	20.0
2-Pentanone, 4-hy...	4.80	29.5	ng	3166220	1	6.66	2145600	20.0
unknown6.43	6.43	86.5	ng	9283230	1	6.66	2145600	20.0