

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051416\
 Data File : BF087033.D
 Acq On : 14 May 2016 21:37
 Operator : UM/SJ
 Sample : PB90575BL
 Misc :
 ALS Vial : 55 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB90575BL

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.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	1.735	11	13	58	rVB	2507106	5039027	100.00%	11.986%
2	4.033	211	214	225	rBV	55715	123216	2.45%	0.293%
3	4.753	274	277	293	rBV	659978	1177006	23.36%	2.800%
4	5.210	314	317	319	rBV	3823169	3909924	77.59%	9.300%
5	5.610	350	352	361	rVB	53460	68420	1.36%	0.163%
6	6.273	408	410	413	rBV	2784383	4041831	80.21%	9.614%
7	6.387	417	420	422	rBV	3853510	4179829	82.95%	9.942%
8	6.604	437	439	442	rBV	713975	933604	18.53%	2.221%
9	6.764	450	453	456	rBV	3262512	2967121	58.88%	7.058%
10	7.187	487	490	492	rBV	2555864	2845669	56.47%	6.769%
11	7.896	550	552	555	rBV	1266047	1221682	24.24%	2.906%
12	8.982	644	647	649	rBV	3905975	4538290	90.06%	10.795%
13	9.645	702	705	708	rBV	1459150	1362615	27.04%	3.241%
14	10.433	771	774	777	rBV	2392024	2425251	48.13%	5.769%
15	11.119	832	834	837	rBV	1253056	1310738	26.01%	3.118%
16	12.536	956	958	961	rBV	64680	89386	1.77%	0.213%
17	12.708	970	973	976	rBV	3853257	3855107	76.50%	9.170%
18	13.245	1018	1020	1022	rBV	67620	56042	1.11%	0.133%
19	13.748	1061	1064	1073	rBV	1068784	1024995	20.34%	2.438%
20	15.108	1180	1183	1192	rBV	698828	872125	17.31%	2.074%

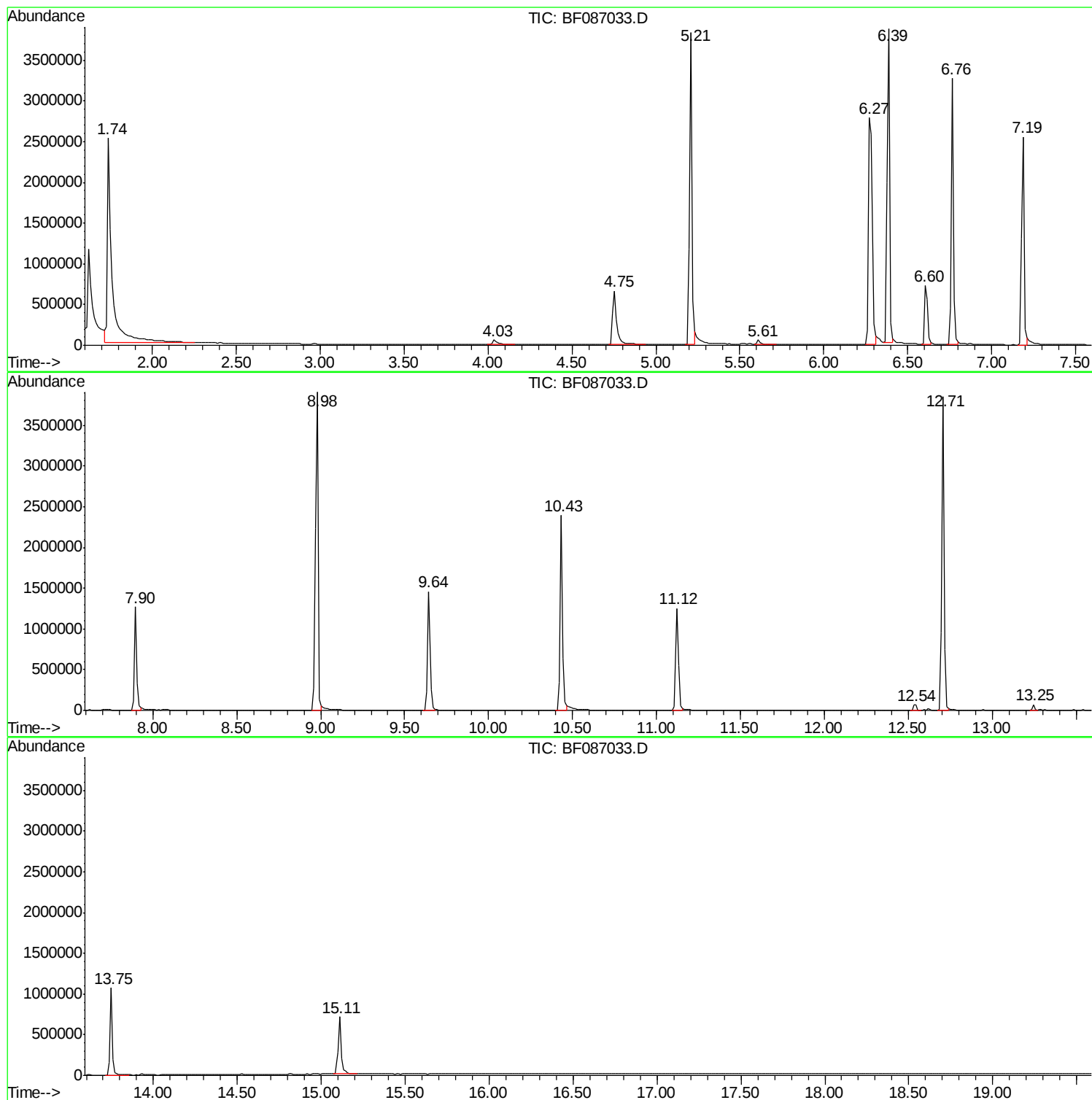
Sum of corrected areas: 42041878

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051416\
Data File : BF087033.D
Acq On : 14 May 2016 21:37
Operator : UM/SJ
Sample : PB90575BL
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PB90575BL

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051416\
Data File : BF087033.D
Acq On : 14 May 2016 21:37
Operator : UM/SJ
Sample : PB90575BL
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PB90575BL

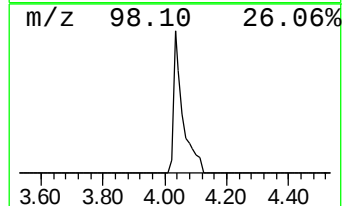
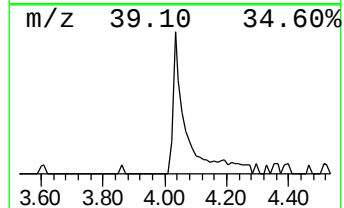
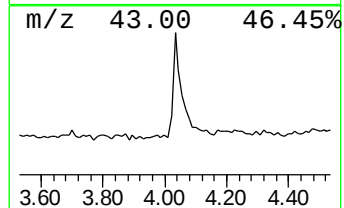
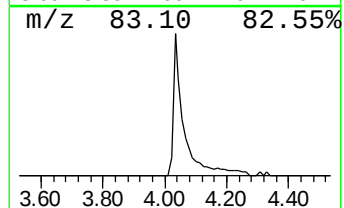
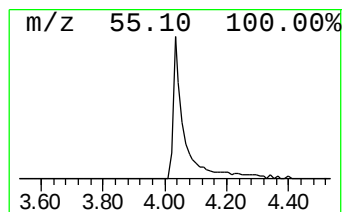
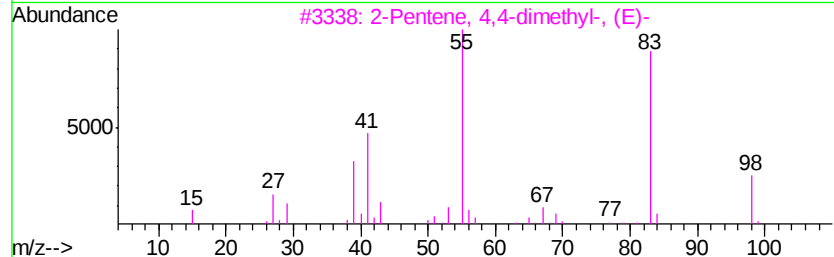
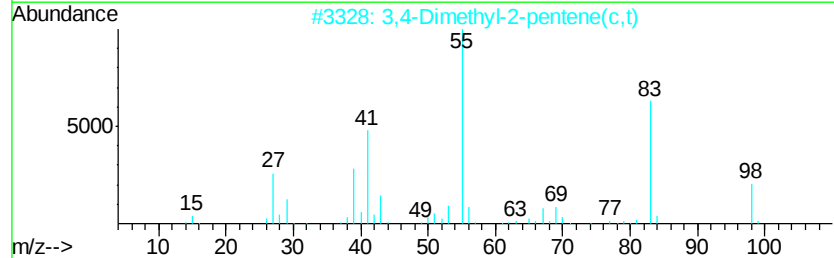
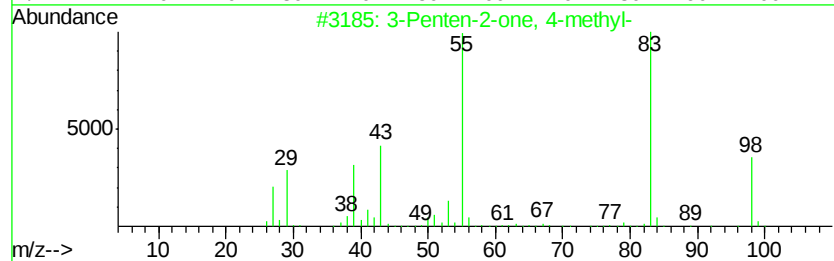
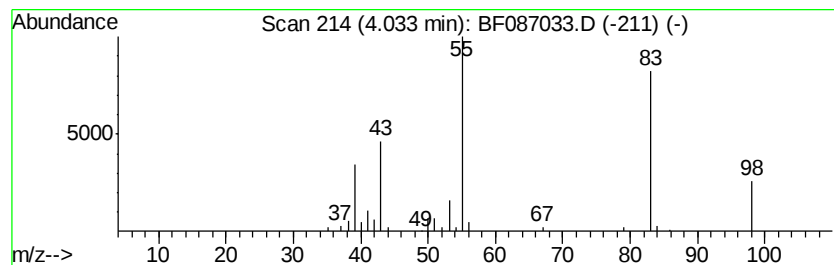
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 3-Penten-2-one, 4-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.03	2.64 ng	123216	1,4-Dichlorobenzene-d4	6.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	90
2		3,4-Dimethyl-2-pentene(c,t)	98	C7H14	1000118-16-5	72
3		2-Pentene, 4,4-dimethyl-, (E)-	98	C7H14	000690-08-4	72
4		2-Pentene, 3,4-dimethyl-, (Z)-	98	C7H14	004914-91-4	72
5		2-Pentene, 4,4-dimethyl-, (Z)-	98	C7H14	000762-63-0	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051416\
Data File : BF087033.D
Acq On : 14 May 2016 21:37
Operator : UM/SJ
Sample : PB90575BL
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PB90575BL

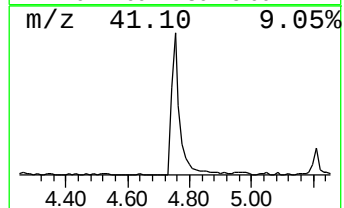
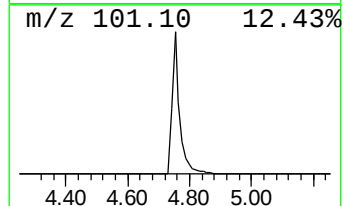
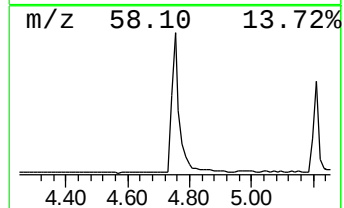
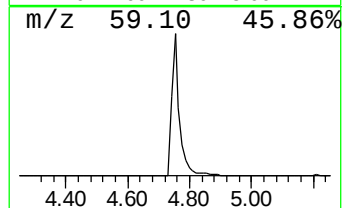
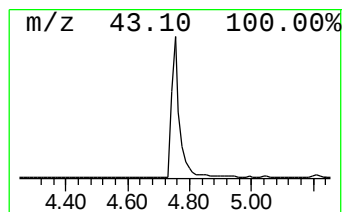
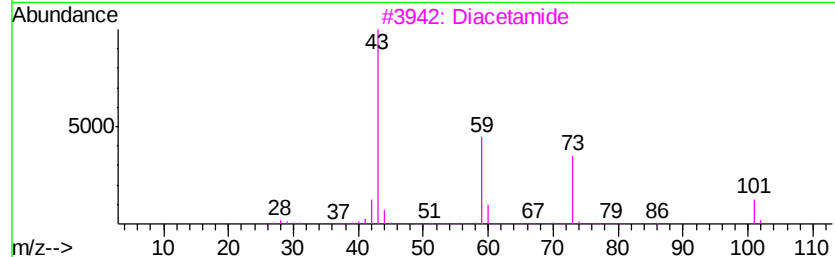
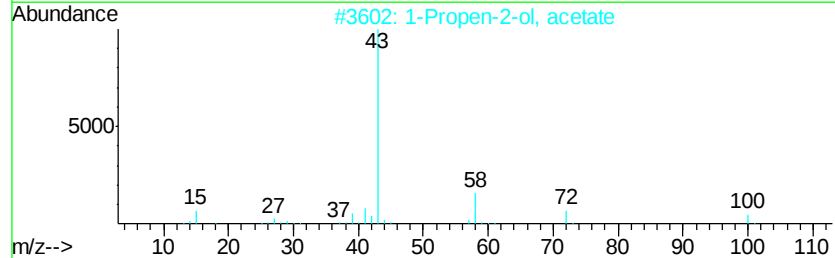
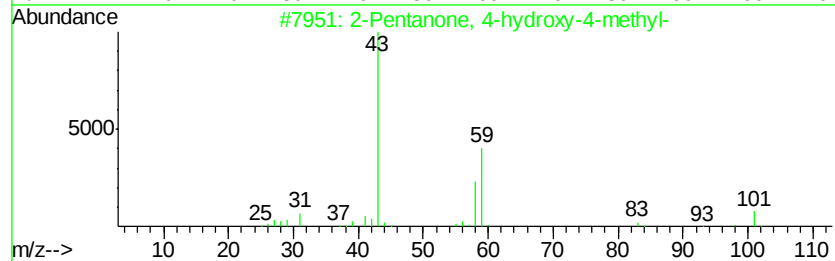
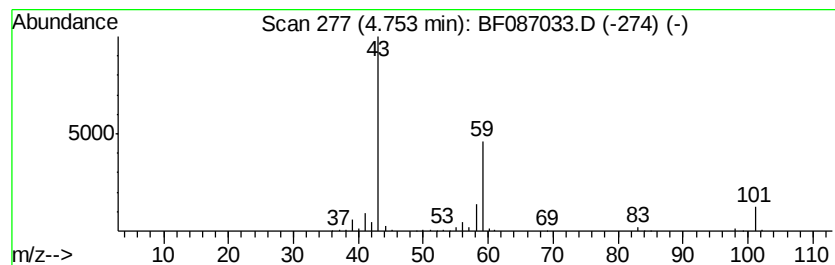
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.75	25.21 ng	1177010	1,4-Dichlorobenzene-d4	6.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
3		Diacetamide	101	C4H7NO2	000625-77-4	9
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		2-Propanone, 1-(1-methylethoxy)-	116	C6H12O2	042781-12-4	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051416\
Data File : BF087033.D
Acq On : 14 May 2016 21:37
Operator : UM/SJ
Sample : PB90575BL
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PB90575BL

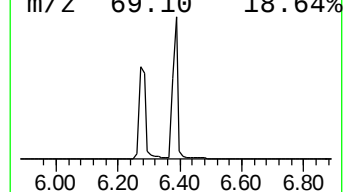
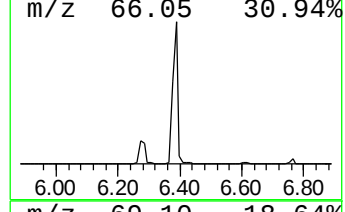
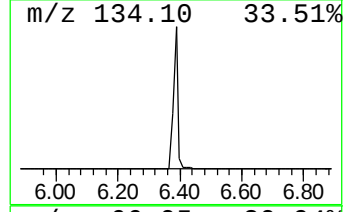
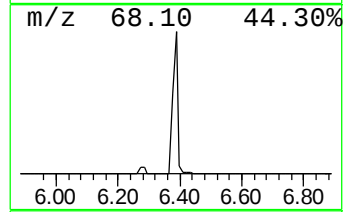
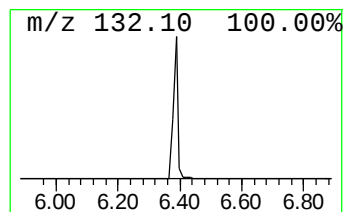
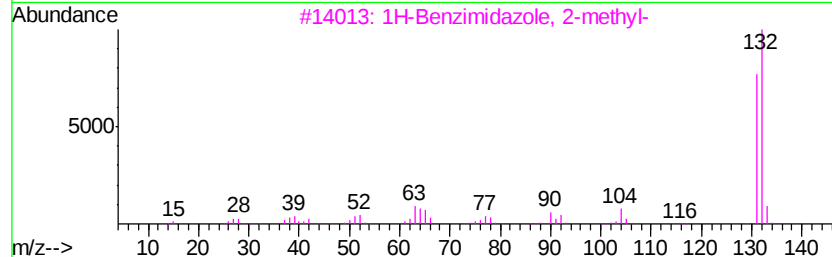
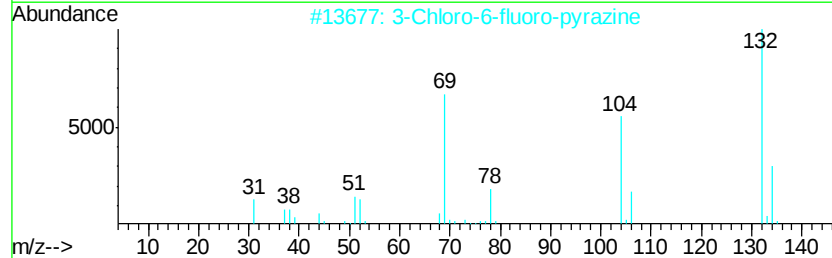
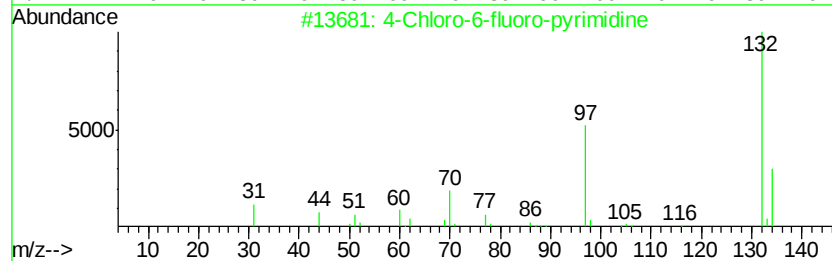
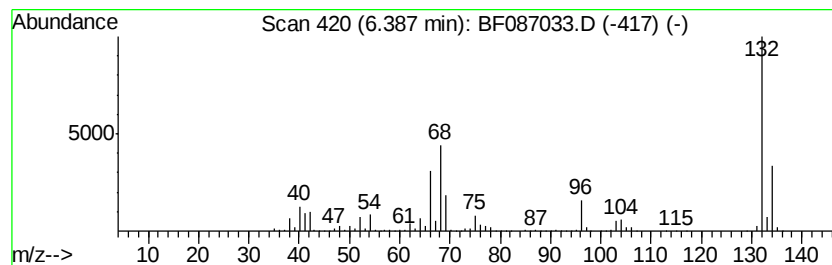
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 unknown6.39 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.39	89.54 ng	4179830	1,4-Dichlorobenzene-d4	6.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
5		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051416\
Data File : BF087033.D
Acq On : 14 May 2016 21:37
Operator : UM/SJ
Sample : PB90575BL
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PB90575BL

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3-Penten-2-one, 4...	4.03	2.6	ng	123216	1	6.62	933604	20.0
2-Pentanone, 4-hy...	4.75	25.2	ng	1177010	1	6.62	933604	20.0
unknown6.39	6.39	89.5	ng	4179830	1	6.62	933604	20.0