

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF110817\
 Data File : BF100302.D
 Acq On : 8 Nov 2017 20:04
 Operator : SJ/JU
 Sample : I6311-01 5X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-04-110617

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-BF110817.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	5.116	512	515	524	rBV	168285	176981	10.17%	1.177%
2	5.504	577	581	584	rBV	914852	754192	43.32%	5.017%
3	6.504	747	751	757	rBV	851864	793500	45.58%	5.279%
4	6.663	774	778	781	rBV	947515	811945	46.64%	5.402%
5	6.898	814	818	822	rBV	1564612	1305502	74.98%	8.685%
6	7.051	840	844	848	rVB	733932	660047	37.91%	4.391%
7	7.457	908	913	916	rBV	527610	461183	26.49%	3.068%
8	8.181	1032	1036	1039	rBV	2095412	1714088	98.45%	11.403%
9	9.251	1215	1218	1222	rVB	1183118	938447	53.90%	6.243%
10	9.645	1282	1285	1288	rVB	27572	20318	1.17%	0.135%
11	9.933	1330	1334	1338	rBV	1855315	1741020	100.00%	11.582%
12	10.645	1451	1455	1458	rBV	86254	70183	4.03%	0.467%
13	10.716	1464	1467	1471	rBV	501573	399275	22.93%	2.656%
14	11.422	1583	1587	1590	rBV	1829819	1547387	88.88%	10.294%
15	11.445	1590	1591	1594	rVB	57205	27617	1.59%	0.184%
16	11.810	1650	1653	1656	rBV	76613	56158	3.23%	0.374%
17	12.498	1767	1770	1771	rBV	45521	40736	2.34%	0.271%
18	12.516	1771	1773	1776	rVB	201229	147775	8.49%	0.983%
19	12.598	1785	1787	1790	rBV	36406	31982	1.84%	0.213%
20	12.633	1790	1793	1796	rVB	83962	75526	4.34%	0.502%
21	12.863	1829	1832	1834	rVB	65839	57975	3.33%	0.386%
22	13.004	1852	1856	1859	rBV	707594	629156	36.14%	4.186%
23	13.898	2005	2008	2015	rVB8	22966	38514	2.21%	0.256%
24	14.057	2030	2035	2038	rBV	1283210	1191794	68.45%	7.929%
25	15.098	2210	2212	2220	rVB	45491	88946	5.11%	0.592%
26	15.539	2282	2287	2300	rVB	919503	1251257	71.87%	8.324%

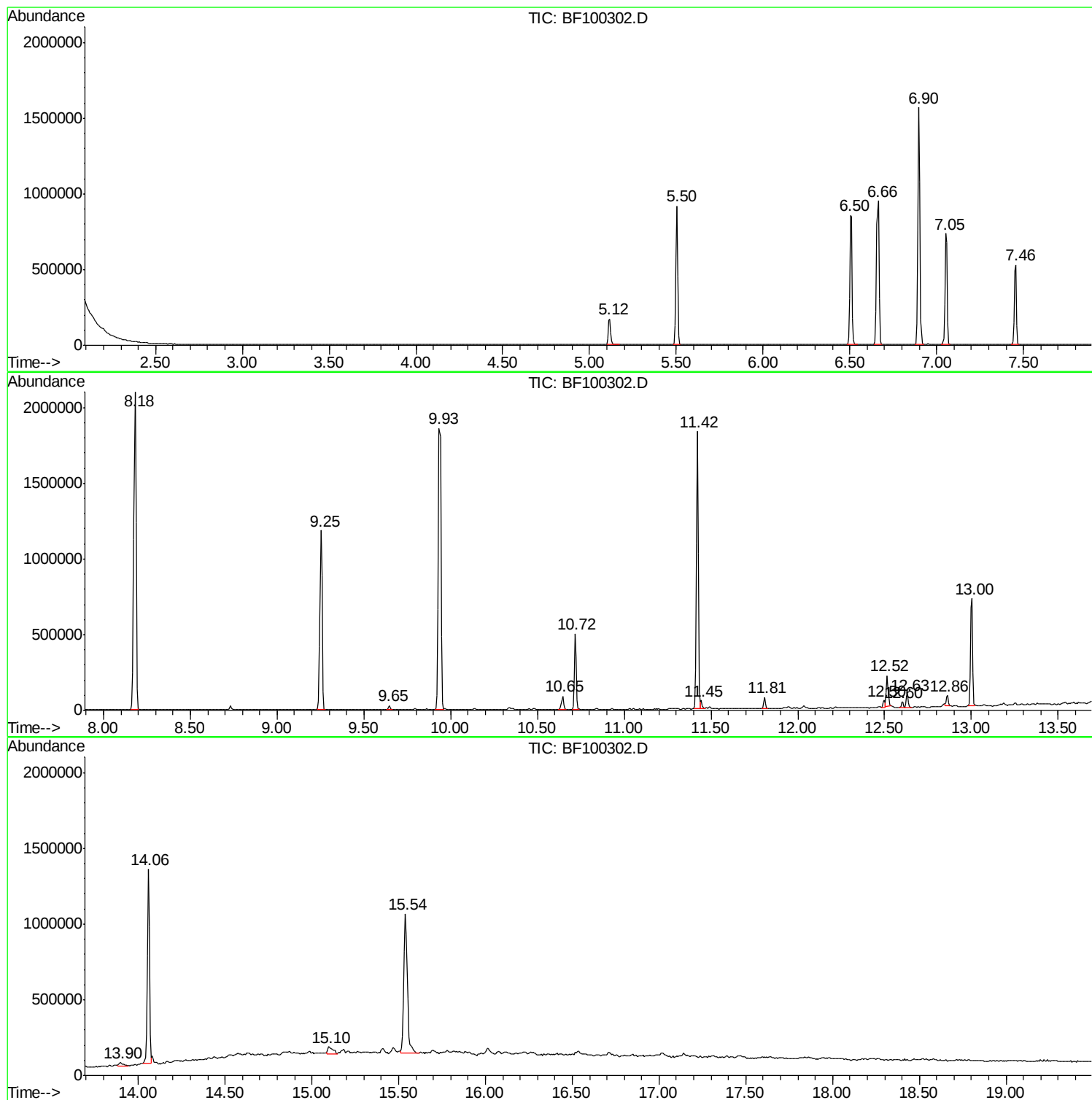
Sum of corrected areas: 15031504

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110817\
Data File : BF100302.D
Acq On : 8 Nov 2017 20:04
Operator : SJ/JU
Sample : I6311-01 5X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TR-04-110617

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110817\
Data File : BF100302.D
Acq On : 8 Nov 2017 20:04
Operator : SJ/JU
Sample : I6311-01 5X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TR-04-110617

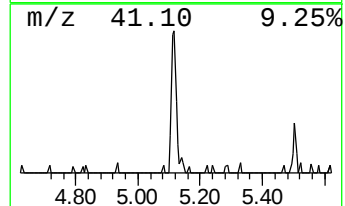
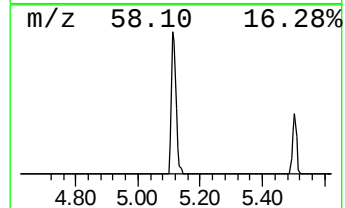
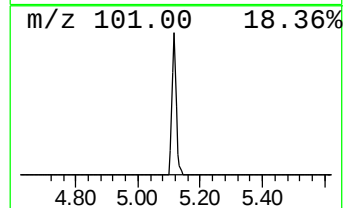
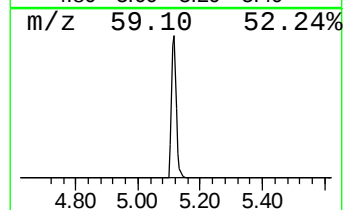
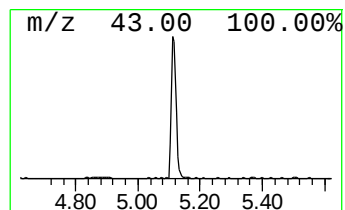
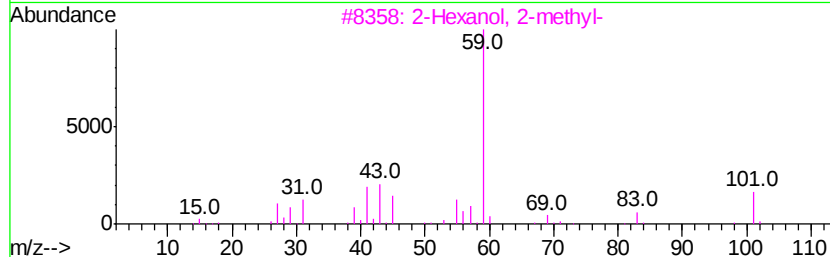
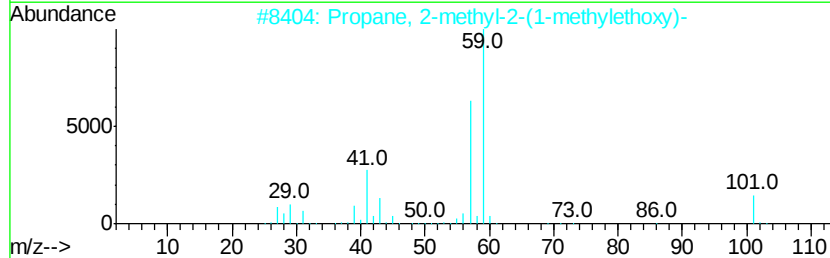
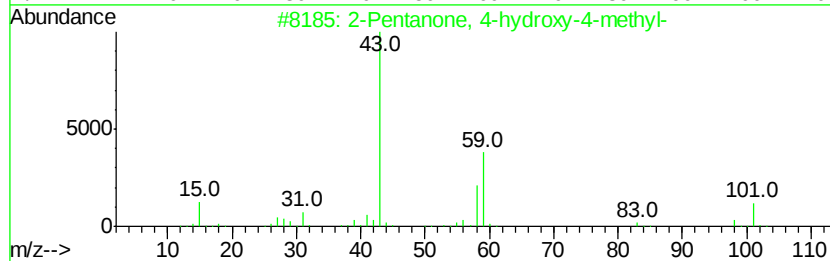
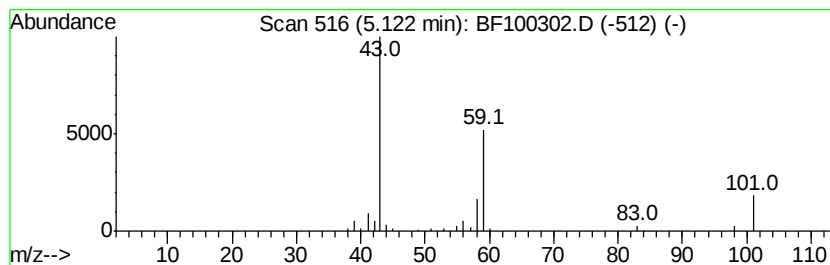
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.12	2.71 ng	176981	1,4-Dichlorobenzene-d4	6.90

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	33
3			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF110817\
Data File : BF100302.D
Acq On : 8 Nov 2017 20:04
Operator : SJ/JU
Sample : I6311-01 5X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TR-04-110617

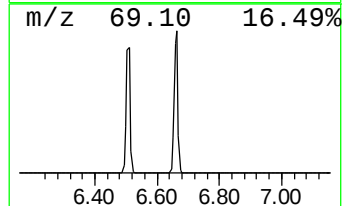
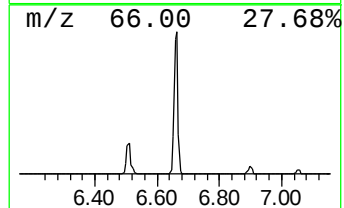
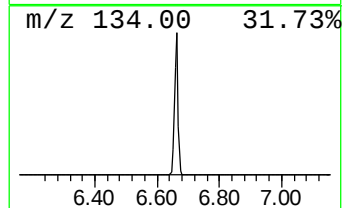
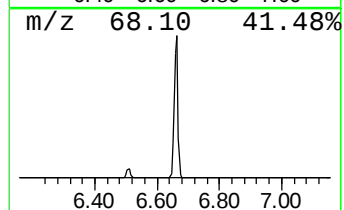
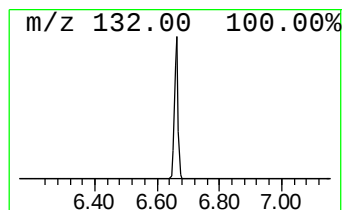
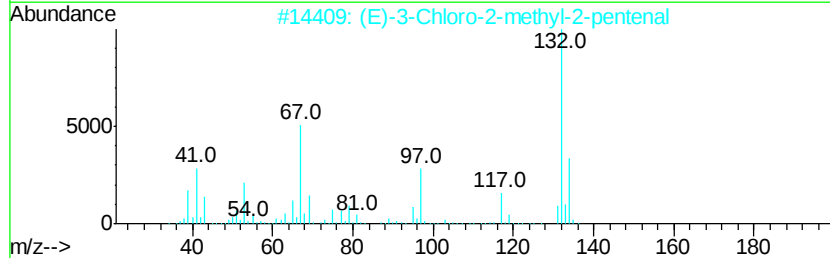
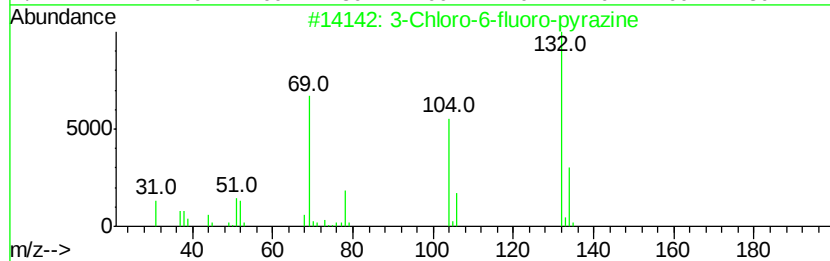
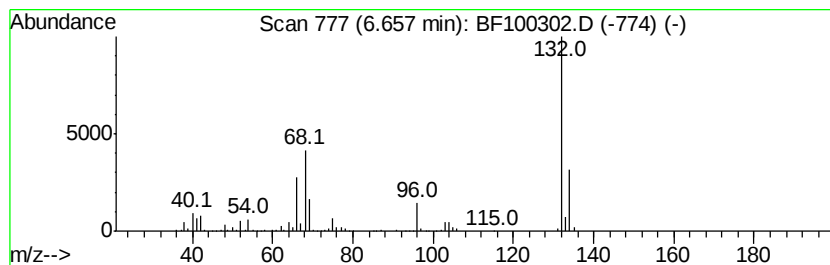
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.66 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.66	12.44 ng	811945	1,4-Dichlorobenzene-d4	6.90

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		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF110817\
Data File : BF100302.D
Acq On : 8 Nov 2017 20:04
Operator : SJ/JU
Sample : I6311-01 5X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TR-04-110617

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110817.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
2-Pentanone, 4-hy...	5.12	2.7	ng	176981	1	6.90	1305500	20.0
unknown6.66	6.66	12.4	ng	811945	1	6.90	1305500	20.0