

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF110217\
 Data File : BF100111.D
 Acq On : 3 Nov 2017 4:06
 Operator : SJ/JU
 Sample : I6274-04
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 110117-TIDO-F-U-S

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-BF102317.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	4.981	489	492	509	rBV	57898	92135	3.86%	0.698%
2	5.392	560	562	585	rBV	352232	582247	24.38%	4.410%
3	6.422	734	737	754	rBV	219996	371369	15.55%	2.813%
4	6.539	754	757	769	rVB	1222858	1188895	49.77%	9.004%
5	6.769	793	796	806	rBV	457063	441239	18.47%	3.342%
6	6.922	818	822	842	rBV	1790730	1675163	70.13%	12.687%
7	7.339	889	893	915	rBV	1199992	1229777	51.49%	9.314%
8	8.051	1010	1014	1033	rBV	656878	606853	25.41%	4.596%
9	9.127	1192	1197	1200	rBV	2495428	2388548	100.00%	18.090%
10	9.804	1304	1312	1325	rBB2	773148	687791	28.80%	5.209%
11	10.474	1423	1426	1431	rVB	36085	31373	1.31%	0.238%
12	10.598	1444	1447	1460	rBV	856297	803279	33.63%	6.084%
13	11.298	1562	1566	1575	rBV	631980	591388	24.76%	4.479%
14	12.880	1830	1835	1838	rBV	1965625	1755237	73.49%	13.294%
15	13.945	2013	2016	2024	rBV	391827	393668	16.48%	2.982%
16	15.409	2260	2265	2276	rVB	268298	364400	15.26%	2.760%

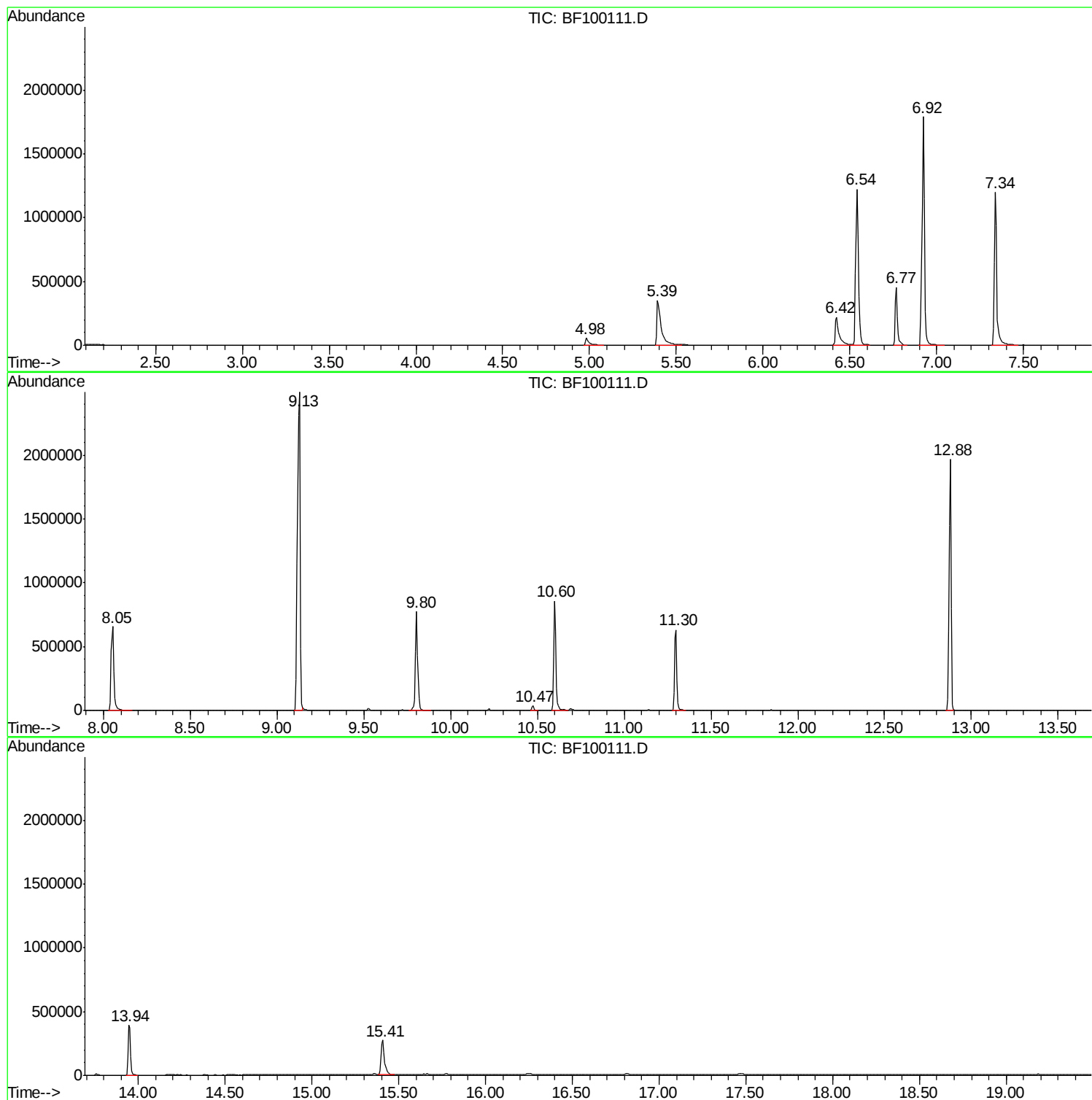
Sum of corrected areas: 13203362

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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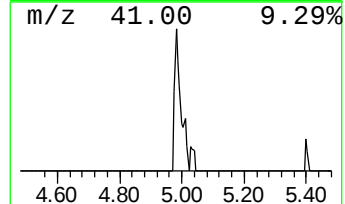
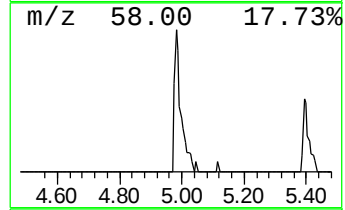
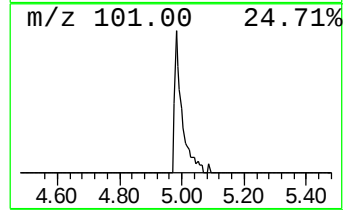
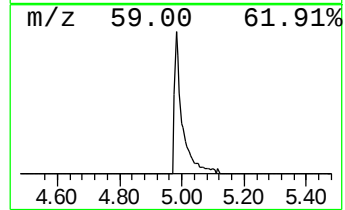
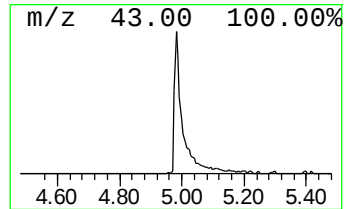
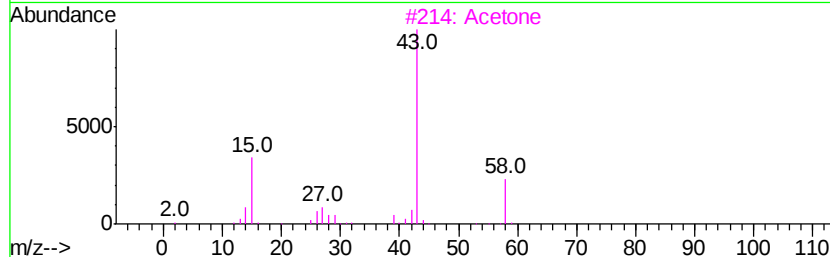
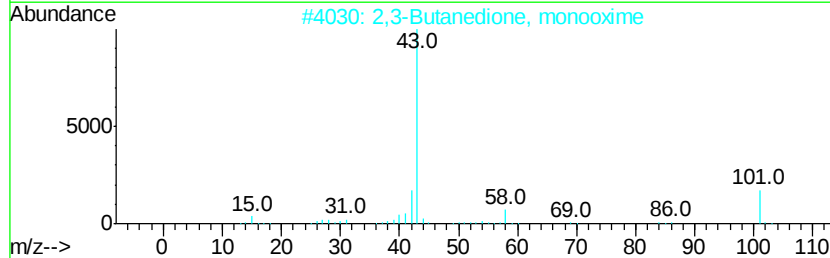
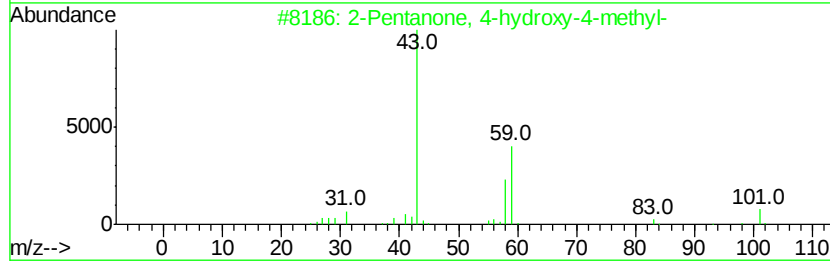
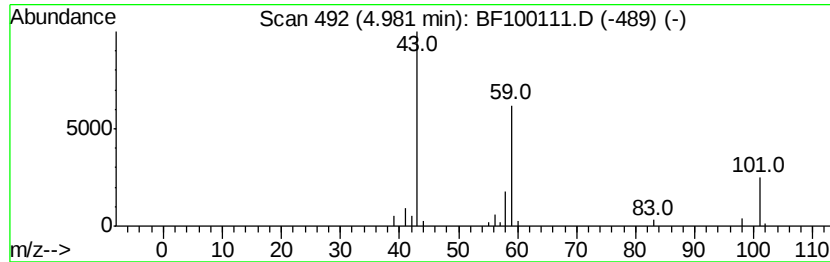
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.98	4.18 ng	92135	1,4-Dichlorobenzene-d4	6.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
3		Acetone	58	C3H6O	000067-64-1	9
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		5-Hexen-2-one	98	C6H10O	000109-49-9	9



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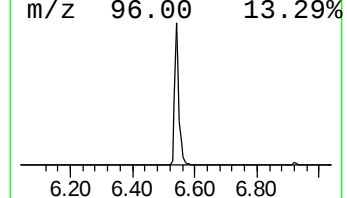
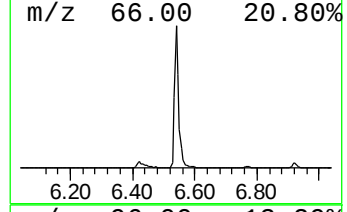
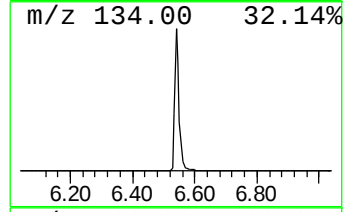
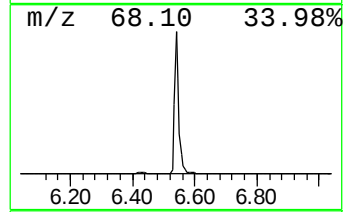
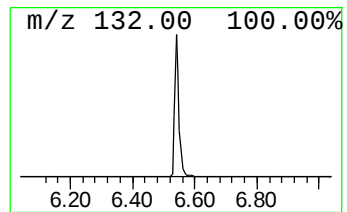
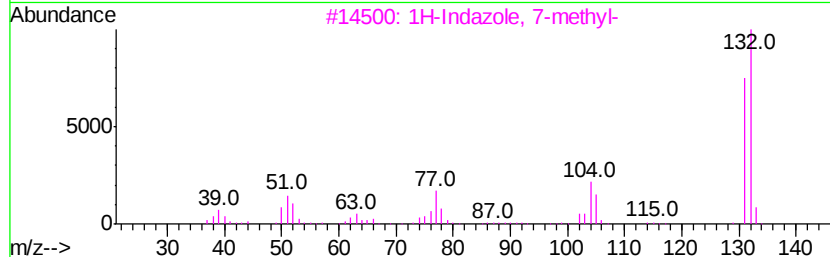
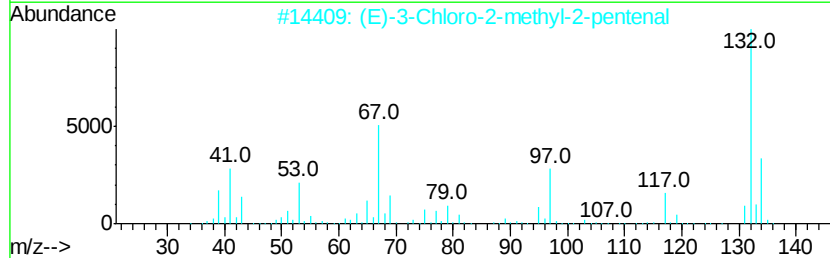
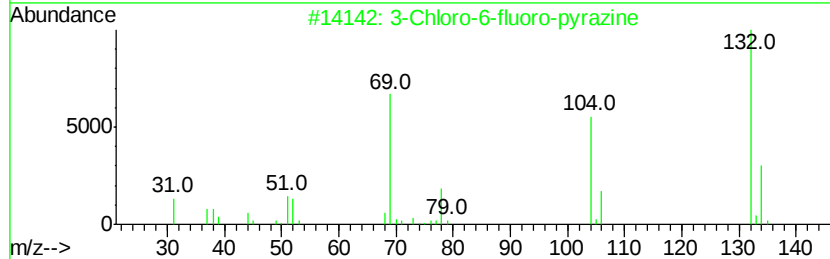
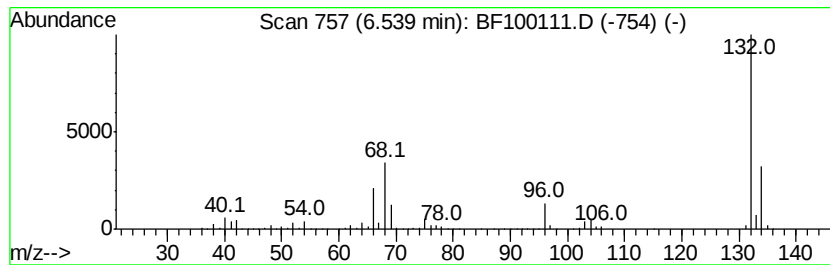
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Peak Number 2 unknown6.54 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.54	53.89 ng	1188900	1,4-Dichlorobenzene-d4	6.77

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		1H-Indazole, 7-methyl-	132	C8H8N2	1000316-00-7	10
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		Xenon	132	Xe	007440-63-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.98	4.2	ng	92135	1	6.77	441239	20.0
unknown6.54	6.54	53.9	ng	1188900	1	6.77	441239	20.0