

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030416\
 Data File : BF085206.D
 Acq On : 4 Mar 2016 16:54
 Operator : SJ/IZ
 Sample : PB88703BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB88703BL

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-BF030316.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	2.356	4	6	13	rVB	51188	75102	1.81%	0.218%
2	5.179	251	253	258	rVB	493007	796048	19.16%	2.314%
3	5.579	285	288	291	rBV	2539783	3247900	78.16%	9.442%
4	6.596	373	377	379	rBV	3091243	3565051	85.79%	10.364%
5	6.733	387	389	392	rBV	2818708	3406419	81.97%	9.903%
6	6.973	407	410	412	rBV	986881	972309	23.40%	2.827%
7	7.122	421	423	426	rVB	2128294	2768709	66.63%	8.049%
8	7.533	456	459	461	rBV	2289356	2438422	58.68%	7.089%
9	8.253	519	522	525	rBV	1471602	1326317	31.92%	3.856%
10	9.328	613	616	619	rBV	3253039	4070673	97.96%	11.834%
11	10.013	673	676	678	rBV	1532401	1490282	35.86%	4.333%
12	10.802	742	745	747	rBV	2329464	2494896	60.04%	7.253%
13	11.499	803	806	808	rBV	1567643	1546464	37.21%	4.496%
14	13.088	942	945	947	rBV	3149212	4155539	100.00%	12.081%
15	14.128	1034	1036	1039	rBV	1131847	1185597	28.53%	3.447%
16	15.603	1162	1165	1168	rBV	681364	857704	20.64%	2.494%

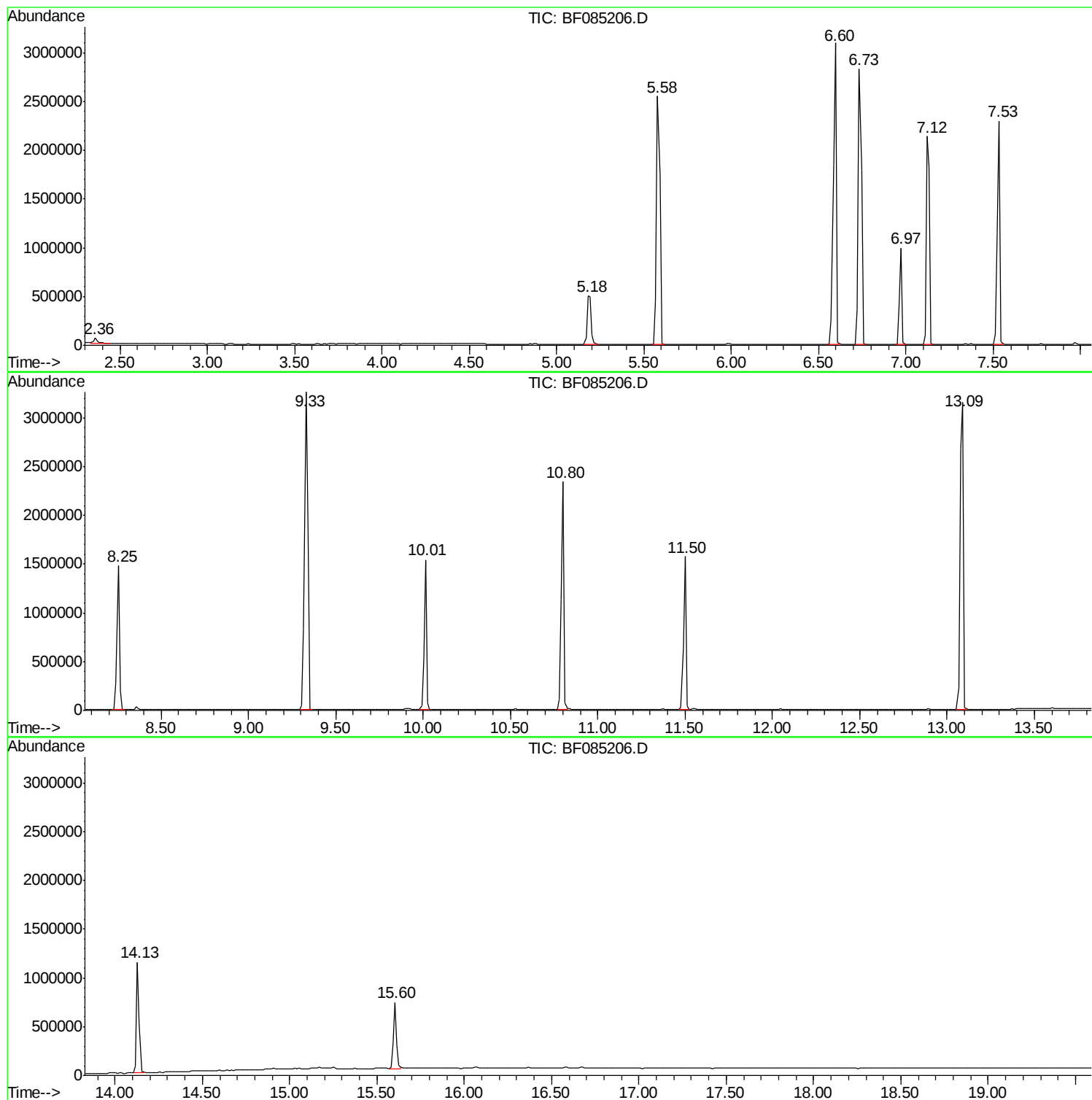
Sum of corrected areas: 34397432

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030316.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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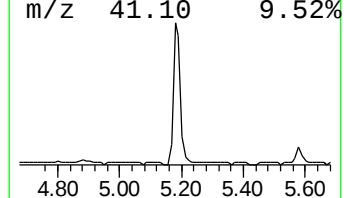
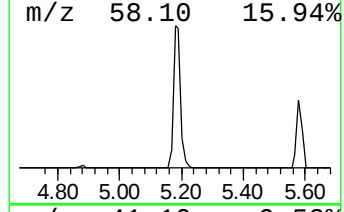
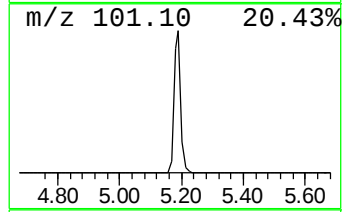
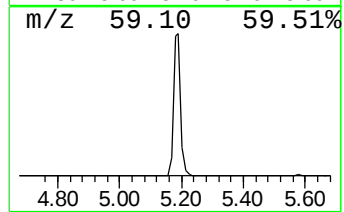
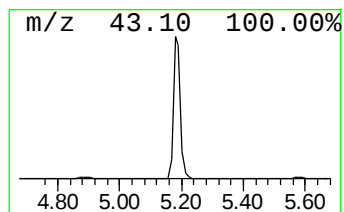
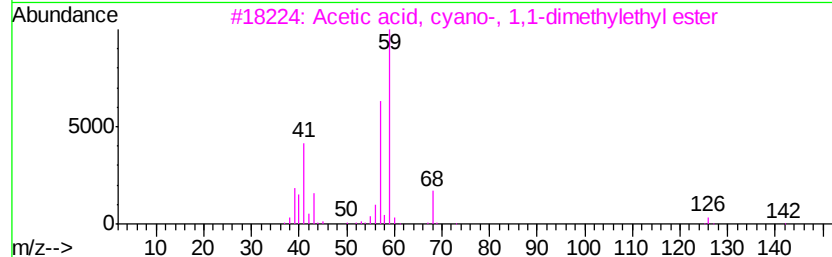
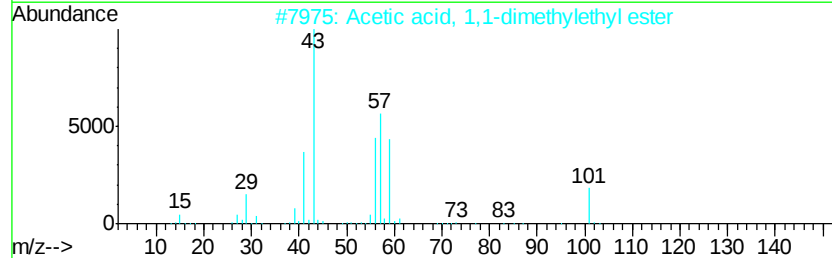
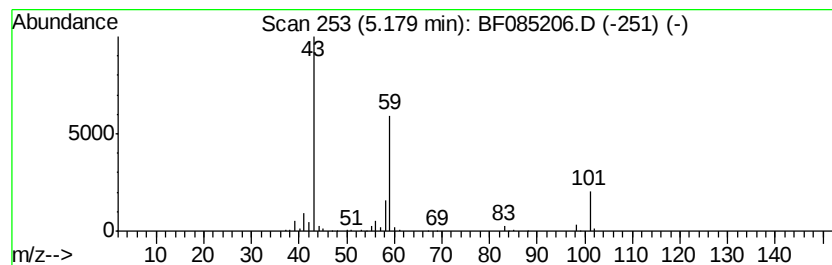
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	16.37 ng	796048	1,4-Dichlorobenzene-d4	6.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
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	23
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



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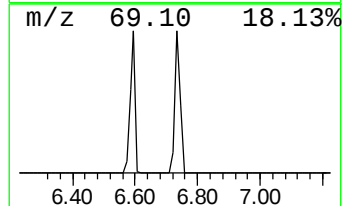
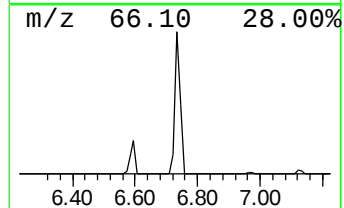
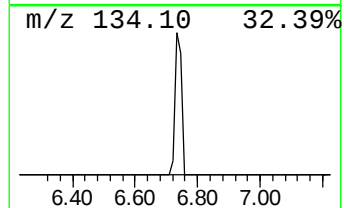
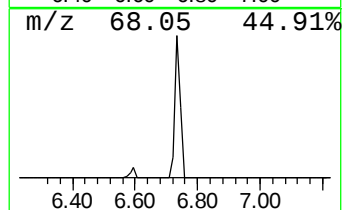
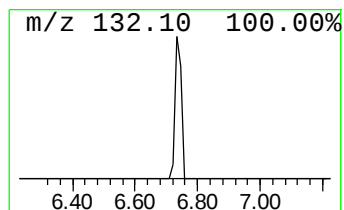
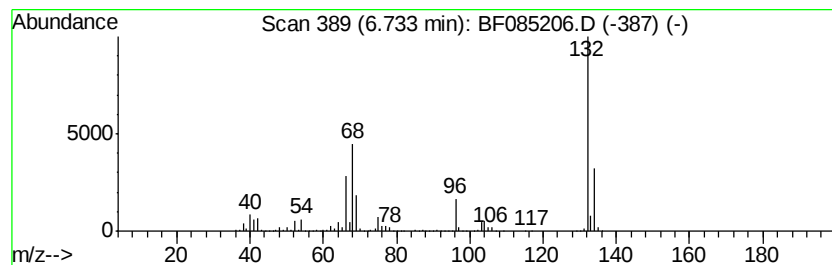
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Peak Number 2 unknown6.73 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.73	70.07 ng	3406420	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.18	16.4	ng	796048	1	6.97	972309	20.0
unknown6.73	6.73	70.1	ng	3406420	1	6.97	972309	20.0