

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030216\
 Data File : BF085113.D
 Acq On : 2 Mar 2016 12:19
 Operator : SJ/IZ
 Sample : PB88622BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB88622BL

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-BF030116.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	2.378	6	8	13	rVB	47887	68219	1.17%	0.154%
2	5.201	252	255	262	rBV	653382	877123	15.01%	1.975%
3	5.601	287	290	292	rBV	4439433	4748169	81.23%	10.690%
4	6.607	375	378	380	rBV	4365983	4613407	78.92%	10.387%
5	6.756	388	391	393	rBV	4750481	4835951	82.73%	10.888%
6	6.984	409	411	414	rVB	1228626	1075054	18.39%	2.420%
7	7.144	422	425	427	rBV	4138219	3656609	62.55%	8.233%
8	7.545	457	460	463	rBV	2708746	2826818	48.36%	6.364%
9	8.276	521	524	526	rBV	1125260	1437477	24.59%	3.236%
10	9.350	615	618	620	rBV	5387676	5568261	95.26%	12.537%
11	10.025	675	677	680	rVB	1571912	1521701	26.03%	3.426%
12	10.813	743	746	749	rBV	3373654	3208864	54.89%	7.225%
13	11.511	805	807	809	rBV	1996783	1644788	28.14%	3.703%
14	13.099	943	946	948	rBV	6106557	5845493	100.00%	13.161%
15	14.151	1035	1038	1040	rBV	1102007	1481823	25.35%	3.336%
16	15.625	1164	1167	1169	rBV	711773	1006426	17.22%	2.266%

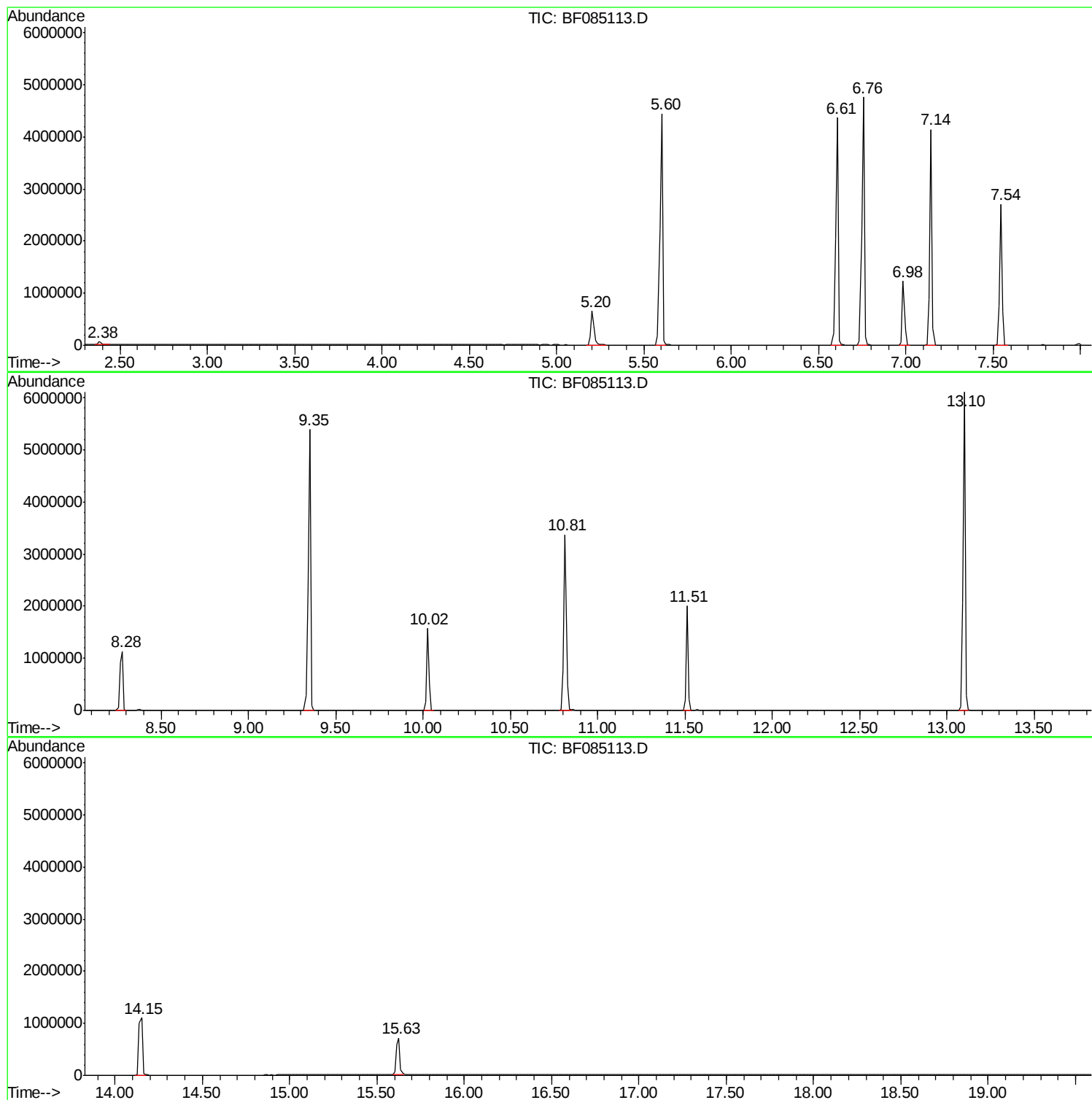
Sum of corrected areas: 44416183

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030116.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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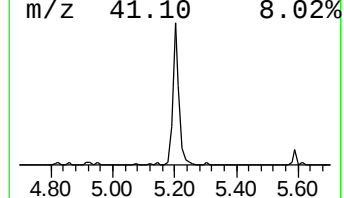
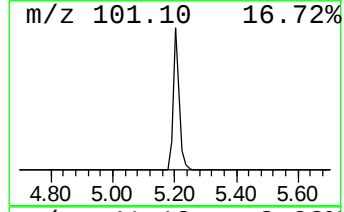
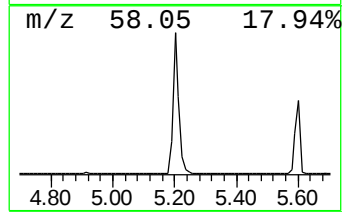
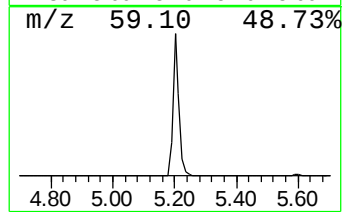
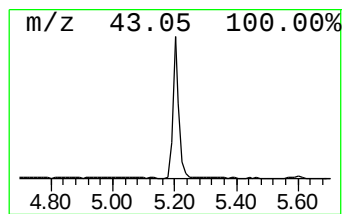
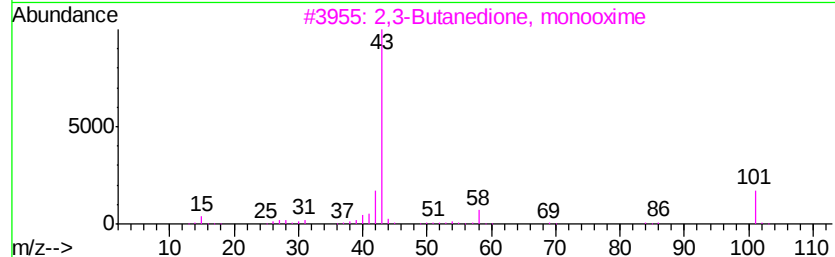
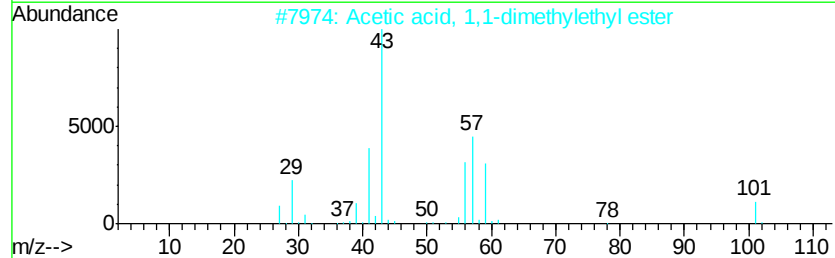
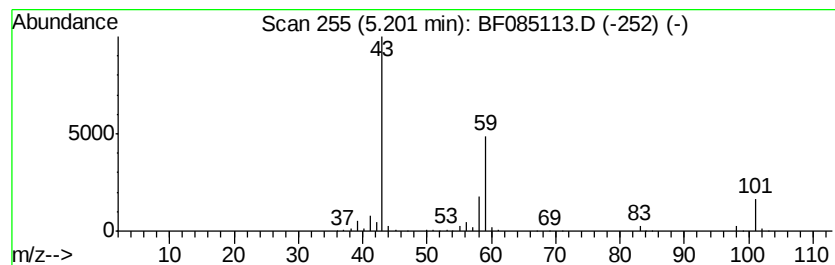
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	16.32 ng	877123	1,4-Dichlorobenzene-d4	6.98

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59	
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28	
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16	
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10	
5	Butane	58	C4H10	000106-97-8	9	



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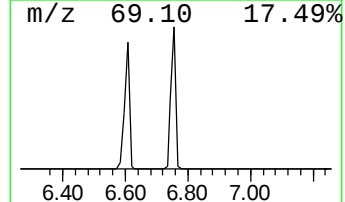
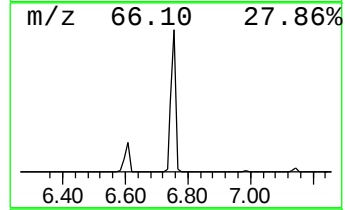
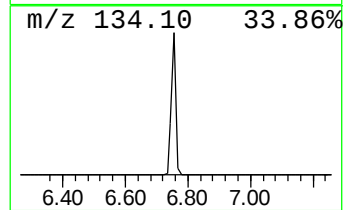
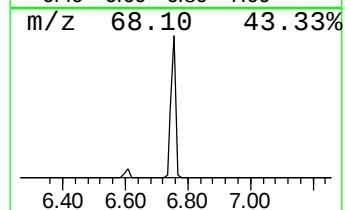
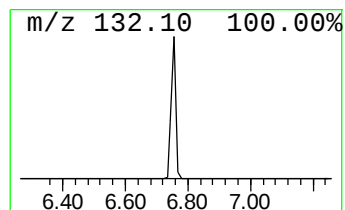
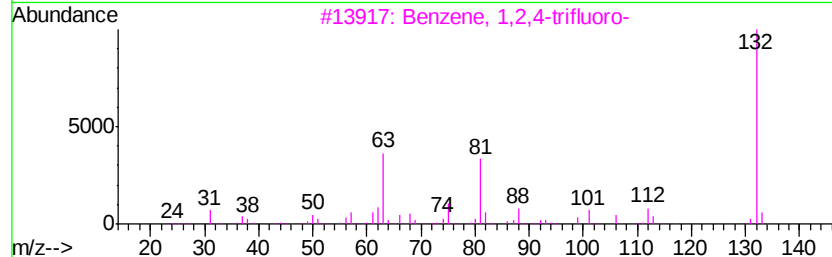
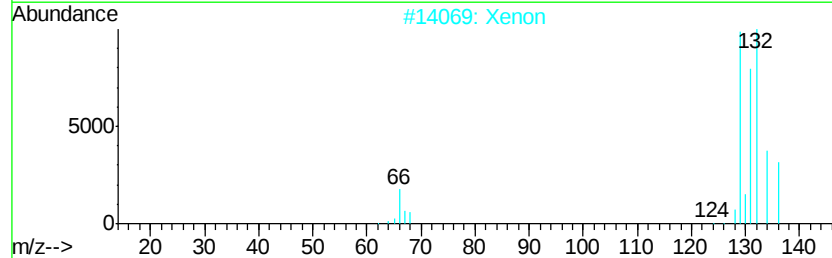
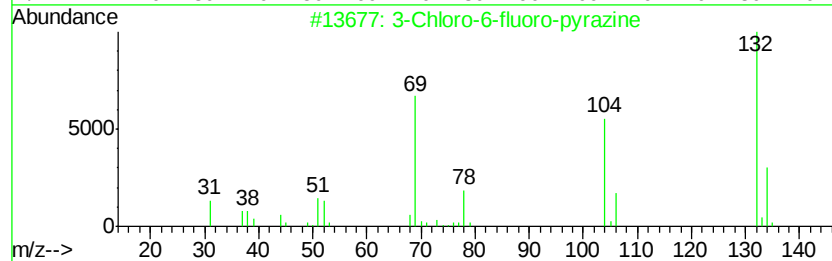
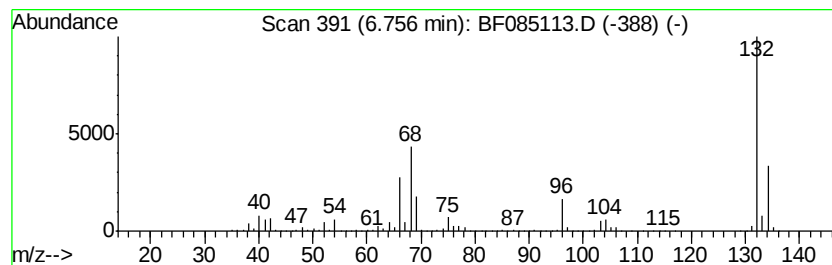
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Peak Number 2 unknown6.76 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.76	89.97 ng	4835950	1,4-Dichlorobenzene-d4	6.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Xenon	132	Xe	007440-63-3	9
3		Benzene, 1,2,4-trifluoro-	132	C6H3F3	000367-23-7	9
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.20	16.3	ng	877123	1	6.98	1075050	20.0
unknown6.76	6.76	90.0	ng	4835950	1	6.98	1075050	20.0