

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062615\
 Data File : BF080201.D
 Acq On : 26 Jun 2015 20:54
 Operator : TP/IZ
 Sample : G2797-01
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DCW-WW-151-62515

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-BF061715.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	5.510	209	212	216	rBV	30052	31460	2.34%	0.457%
2	5.899	244	246	249	rBV	256723	270714	20.14%	3.931%
3	6.893	331	333	336	rBV	160109	163832	12.19%	2.379%
4	7.064	345	348	350	rBV	555370	569765	42.38%	8.273%
5	7.293	366	368	371	rBV	214760	187898	13.98%	2.728%
6	7.453	380	382	385	rVB	736920	591161	43.97%	8.583%
7	7.853	414	417	419	rBV	577395	476822	35.47%	6.923%
8	8.585	479	481	484	rBV	285321	243423	18.11%	3.534%
9	9.659	573	575	578	rBV	1099375	976957	72.67%	14.185%
10	10.048	608	609	611	rVB2	18167	14772	1.10%	0.214%
11	10.356	634	636	641	rVB	341114	299466	22.27%	4.348%
12	11.145	703	705	708	rBV2	461101	636441	47.34%	9.241%
13	11.865	766	768	770	rBV	355870	321348	23.90%	4.666%
14	13.545	912	915	917	rBV	1196371	1344438	100.00%	19.521%
15	14.585	1004	1006	1008	rBV	14709	23953	1.78%	0.348%
16	14.802	1023	1025	1028	rBV	335883	364629	27.12%	5.294%
17	16.665	1185	1188	1193	rBV	235553	370108	27.53%	5.374%

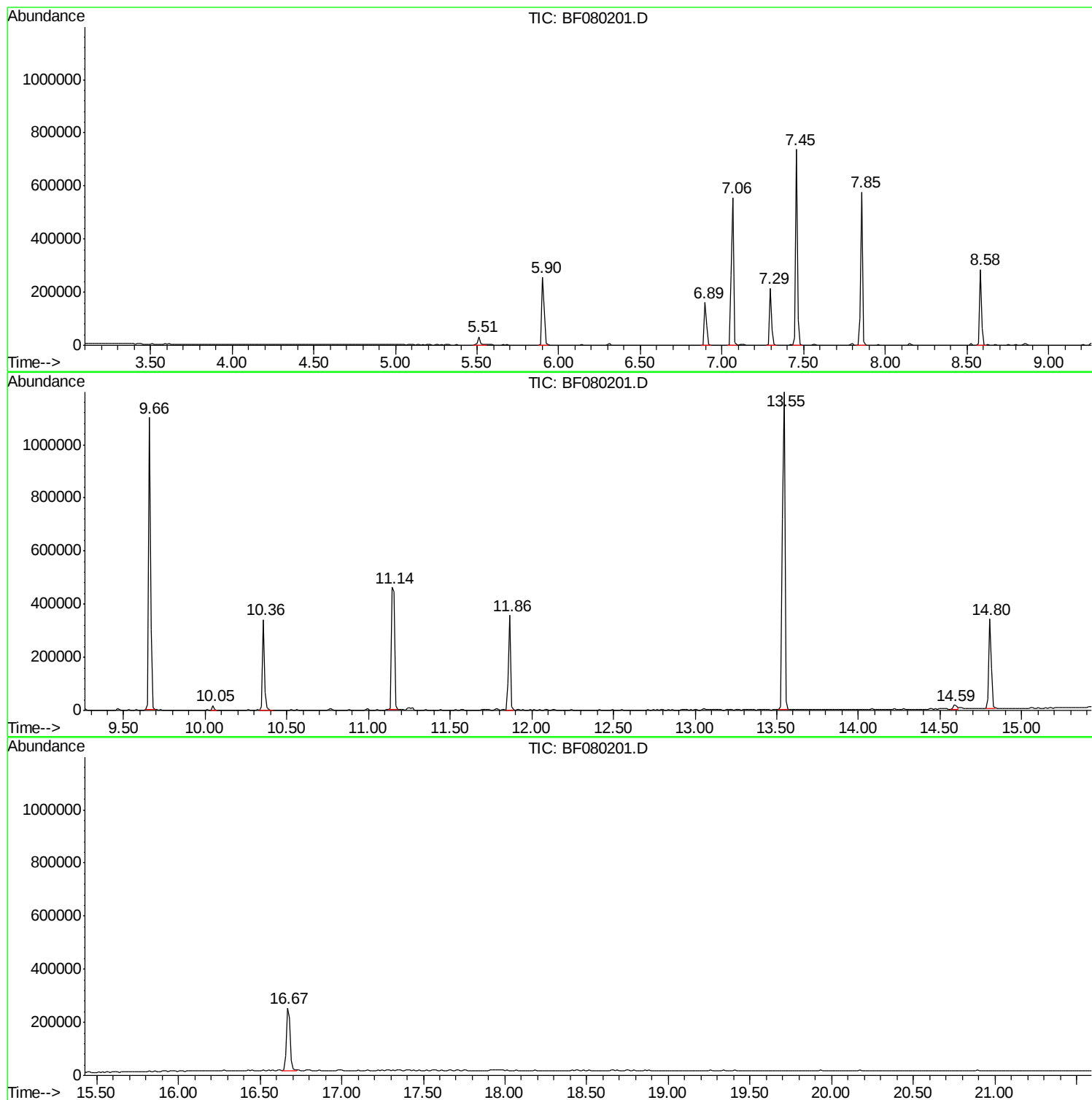
Sum of corrected areas: 6887187

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.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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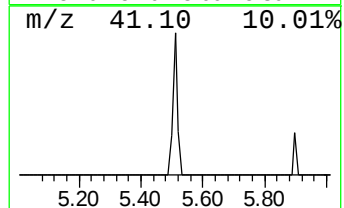
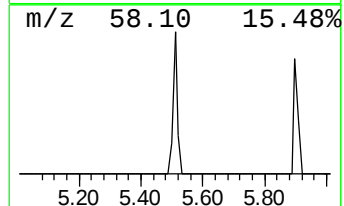
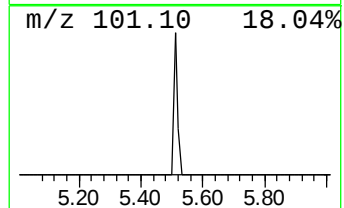
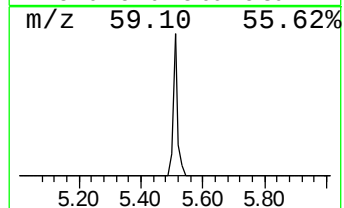
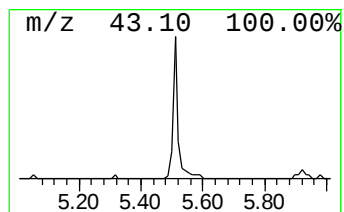
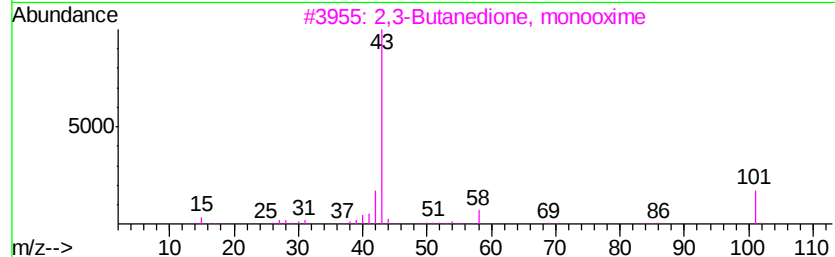
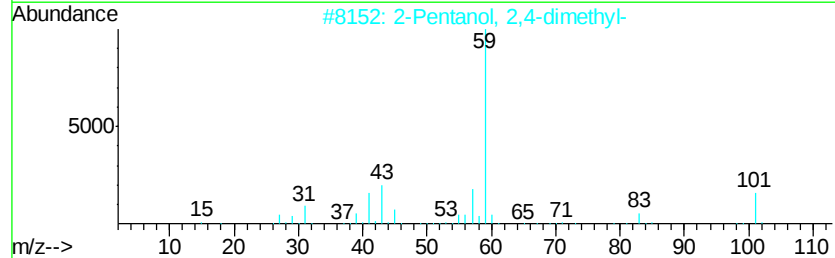
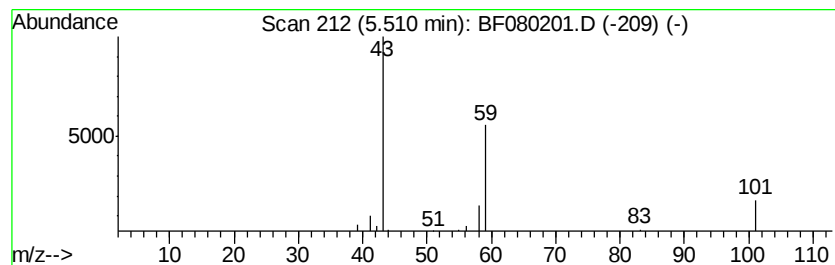
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.51	3.35 ng	31460	1,4-Dichlorobenzene-d4	7.29

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	78
2			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	33
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4			2-Propanone, 1-(1-methylethoxy)-	116	C6H12O2	042781-12-4	9
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9



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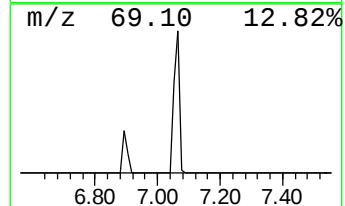
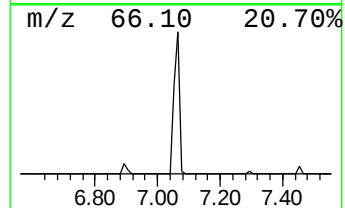
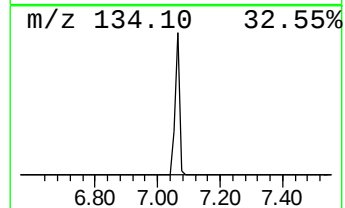
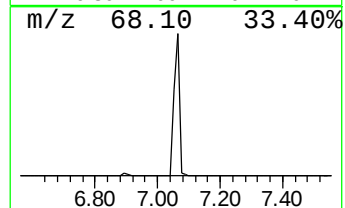
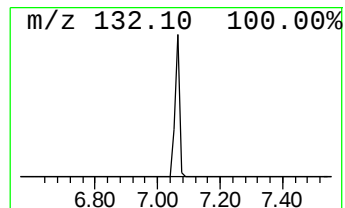
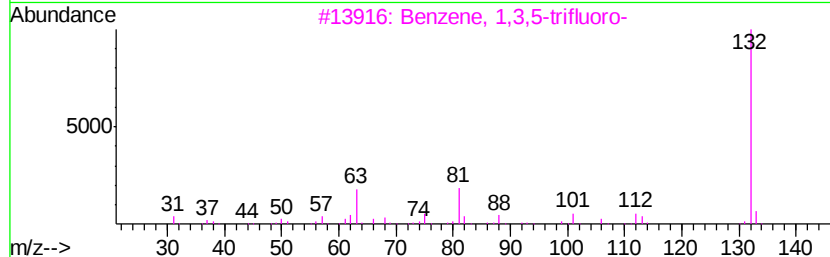
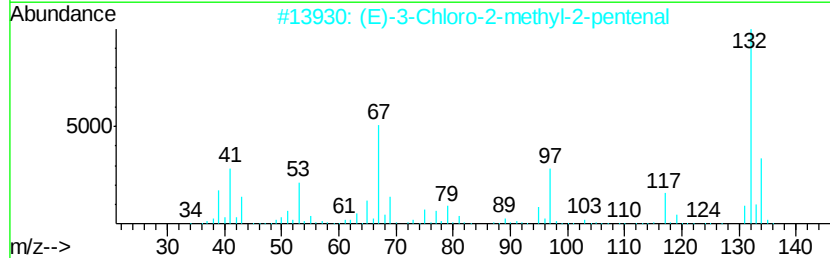
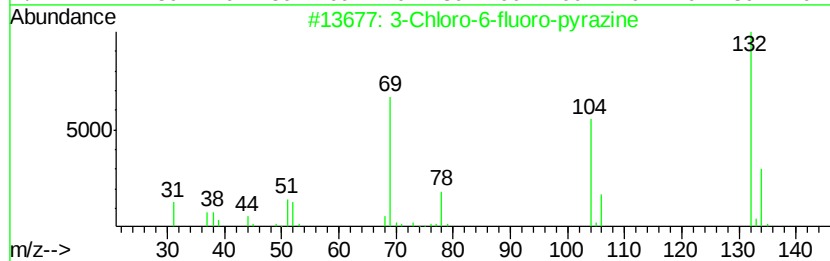
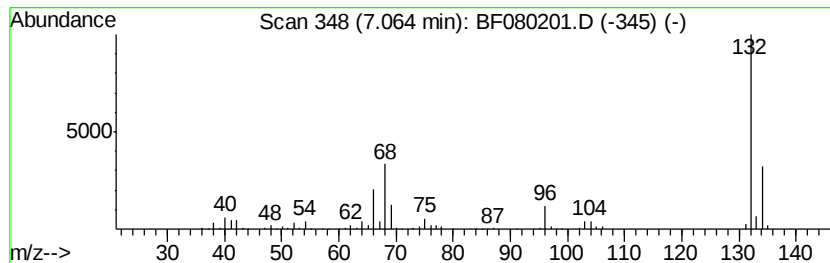
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Peak Number 2 unknown7.06 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.06	60.65 ng	569765	1,4-Dichlorobenzene-d4	7.29

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		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.51	3.4	ng	31460	1	7.29	187898	20.0
unknown7.06	7.06	60.6	ng	569765	1	7.29	187898	20.0