

Data Path : Z:\HPCHEM1\BNA_F\Data\BF030217\
Data File : BF093335.D
Acq On : 2 Mar 2017 21:55
Operator : SJ/MA
Sample : PB97297BL
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PB97297BL

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-BF013017.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	4.956	249	251	267	rBV	489390	813299	17.04%	2.165%
2	5.390	287	289	304	rBV	3619195	4205121	88.13%	11.192%
3	6.441	378	381	384	rBV	3445520	3713113	77.82%	9.883%
4	6.567	389	392	394	rBV	3653836	4166180	87.31%	11.089%
5	6.796	409	412	414	rBV	814036	886351	18.58%	2.359%
6	6.944	423	425	428	rBV	2491172	3139710	65.80%	8.357%
7	7.367	459	462	464	rBV	2647846	2737965	57.38%	7.287%
8	8.087	522	525	534	rBV	1241629	1297409	27.19%	3.453%
9	9.173	617	620	622	rBV	4186827	4771552	100.00%	12.700%
10	9.847	676	679	682	rBV	1437315	1402509	29.39%	3.733%
11	10.648	746	749	751	rBV	2126220	2545403	53.35%	6.775%
12	11.333	807	809	812	rBV	1125666	1351782	28.33%	3.598%
13	12.922	945	948	951	rBV	3884938	4146986	86.91%	11.038%
14	13.974	1037	1040	1047	rVB	1248375	1329673	27.87%	3.539%
15	15.391	1161	1164	1174	rBV	759125	1063966	22.30%	2.832%

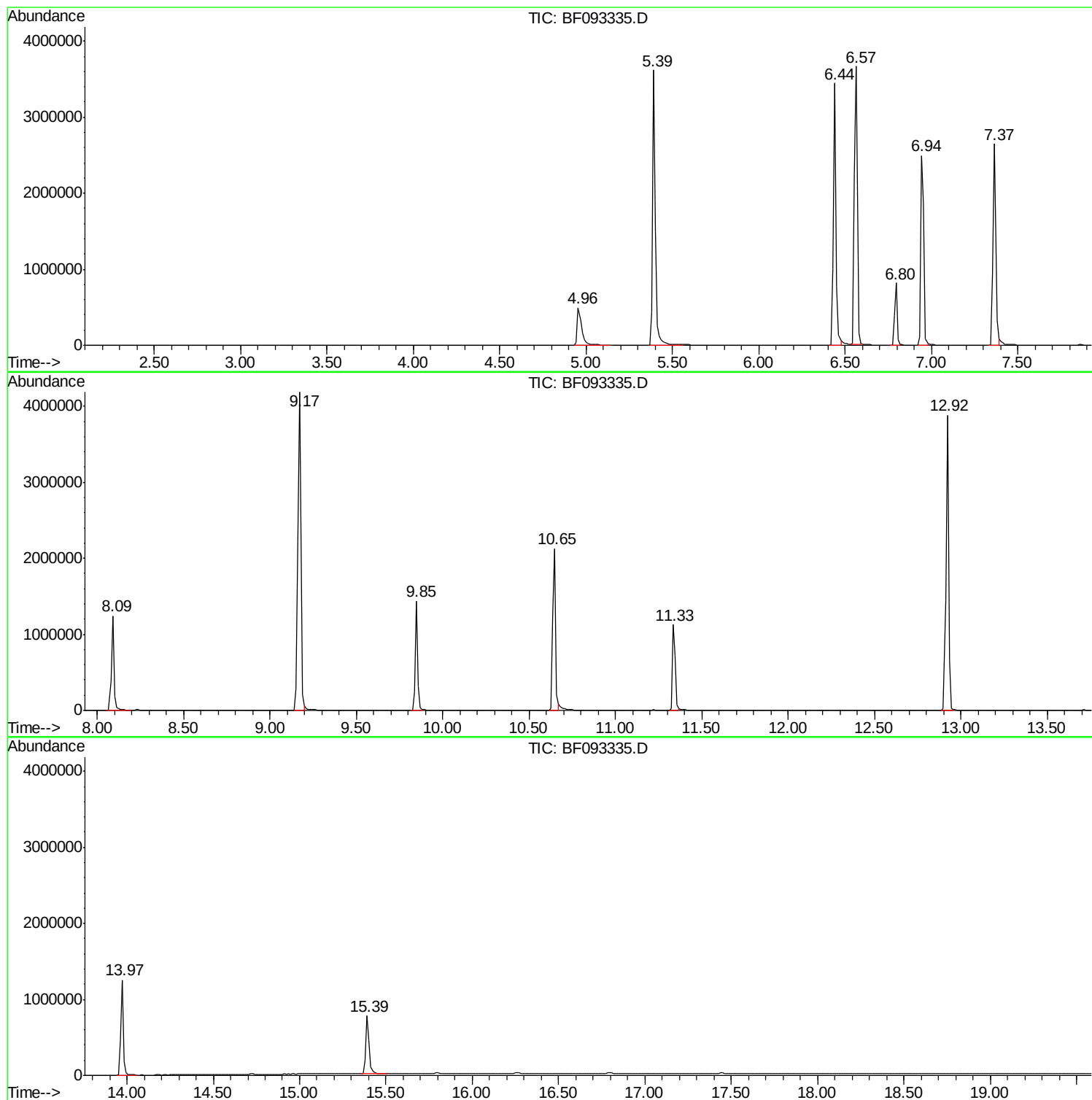
Sum of corrected areas: 37571019

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.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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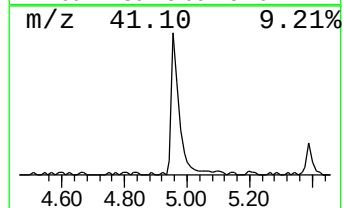
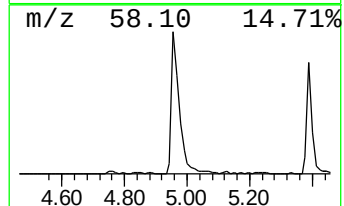
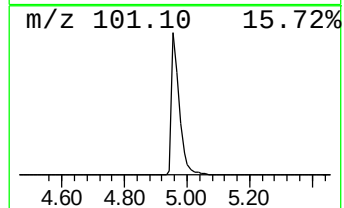
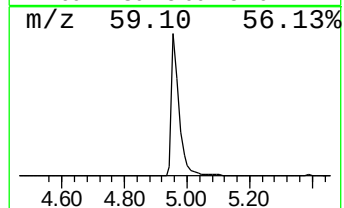
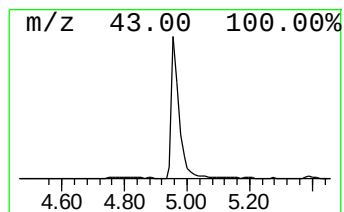
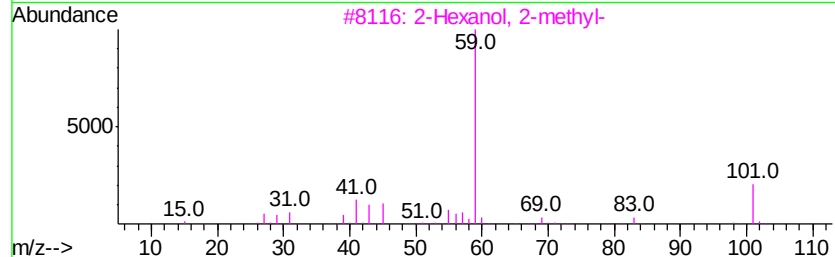
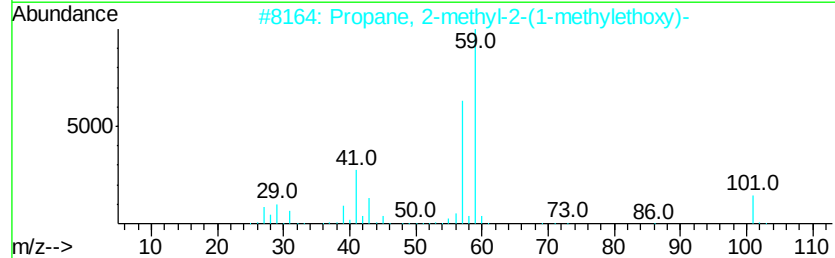
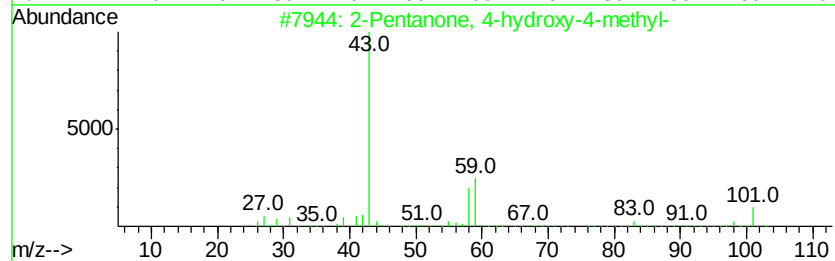
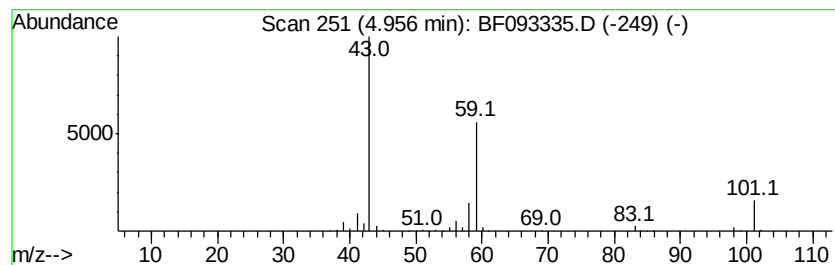
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.96	18.35 ng	813299	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	36
3		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
4		3-Hexanol	102	C6H14O	000623-37-0	33
5		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33



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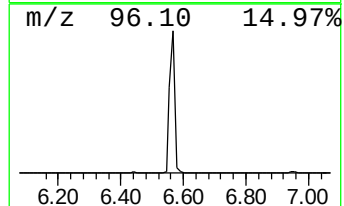
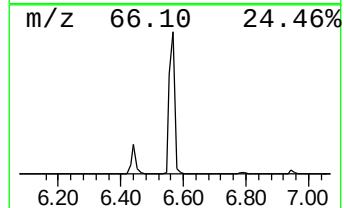
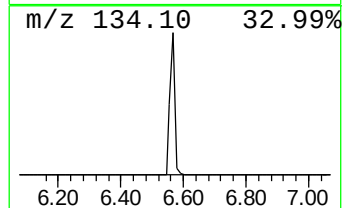
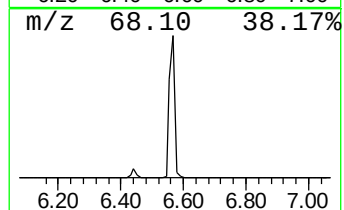
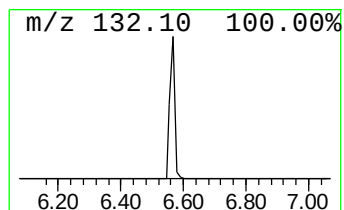
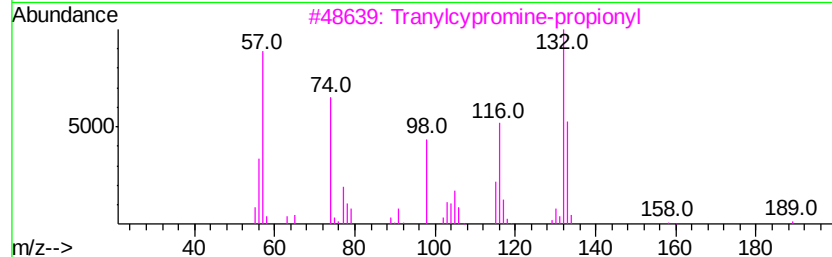
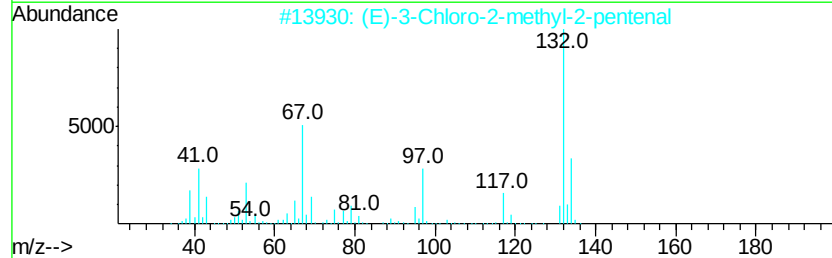
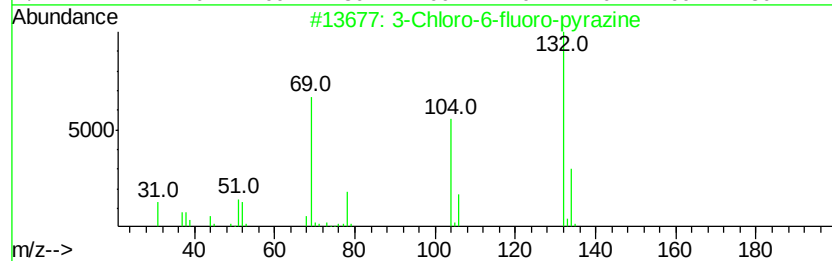
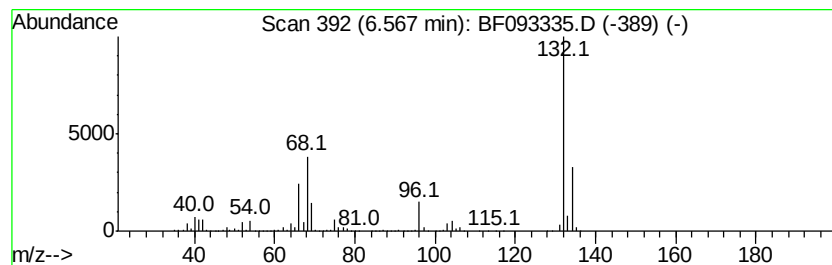
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Peak Number 2 unknown6.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.57	94.01 ng	4166180	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	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...	4.96	18.4	ng	813299	1	6.80	886351	20.0
unknown6.57	6.57	94.0	ng	4166180	1	6.80	886351	20.0