

Data Path : Z:\HPCHEM1\BNA_F\Data\BF031817\
 Data File : BF093740.D
 Acq On : 18 Mar 2017 8:59
 Operator : SJ/MA
 Sample : PB97670BL
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB97670BL

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-BF031817.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	5.030	272	275	283	rVB	566483	983340	15.99%	2.085%
2	5.453	309	312	315	rBV	4210760	4684172	76.15%	9.931%
3	6.482	398	402	404	rBV	4226603	5573555	90.61%	11.817%
4	6.619	411	414	416	rBV	4234603	5393582	87.68%	11.435%
5	6.847	431	434	436	rVB	1281331	1111143	18.06%	2.356%
6	7.007	445	448	450	rBV	3983489	4248288	69.06%	9.007%
7	7.407	480	483	485	rBV	3180647	3459055	56.23%	7.334%
8	8.128	543	546	549	rVB	1490992	1407849	22.89%	2.985%
9	9.213	637	641	643	rBV	4672766	6151384	100.00%	13.042%
10	9.876	697	699	702	rVB	1313505	1549628	25.19%	3.285%
11	10.665	765	768	771	rBV	2827939	3083825	50.13%	6.538%
12	11.362	826	829	831	rBV	1627451	1630084	26.50%	3.456%
13	12.951	965	968	970	rBV	4716310	5201486	84.56%	11.028%
14	13.991	1056	1059	1061	rBV	1255946	1467914	23.86%	3.112%
15	15.397	1179	1182	1185	rBV	858043	1220590	19.84%	2.588%

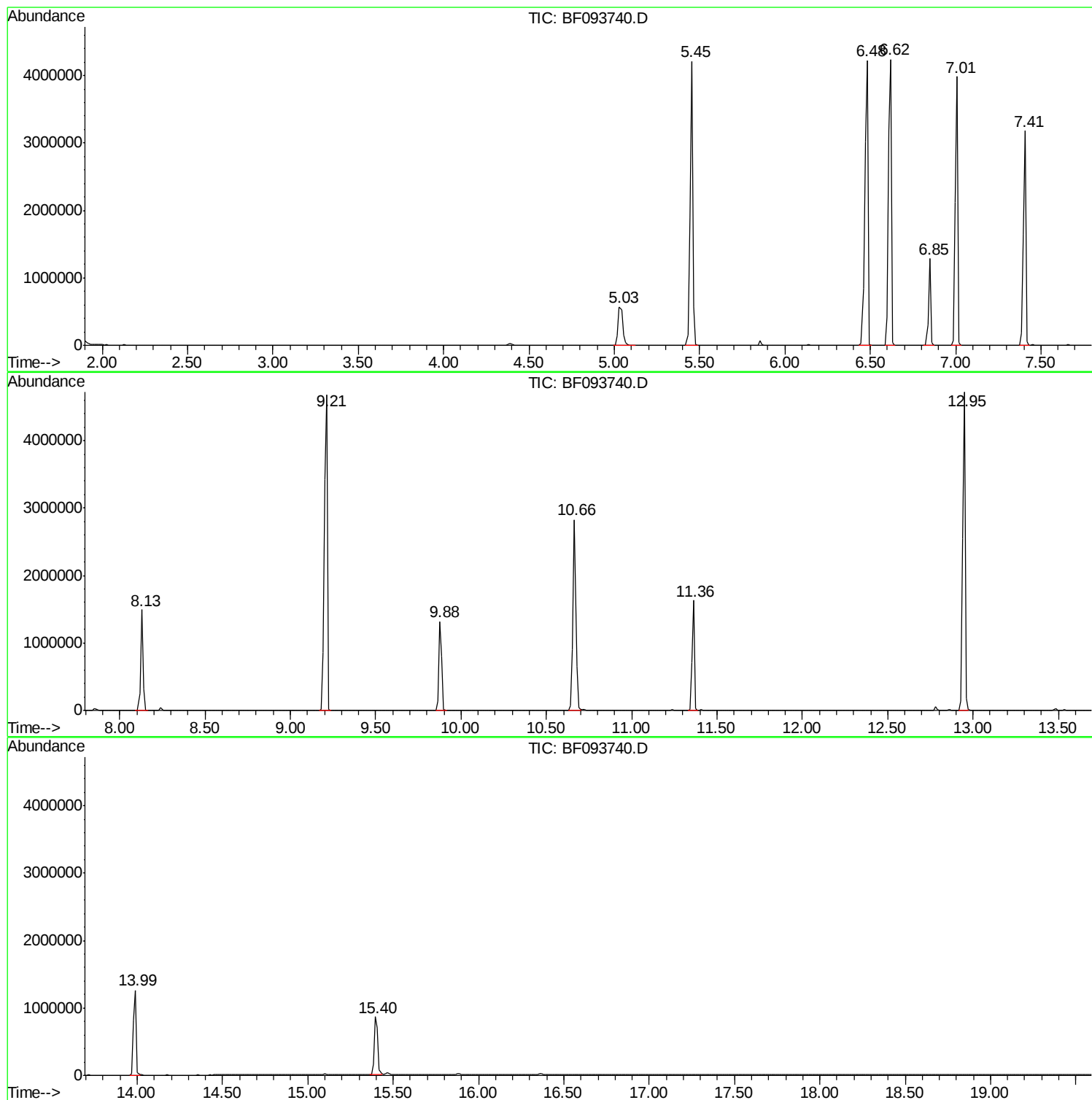
Sum of corrected areas: 47165895

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031817.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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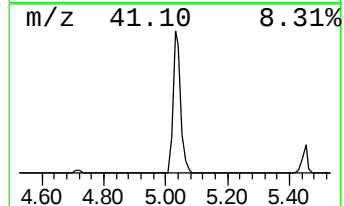
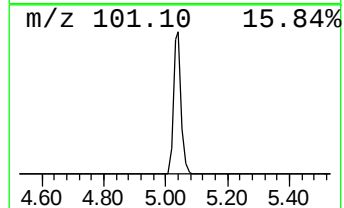
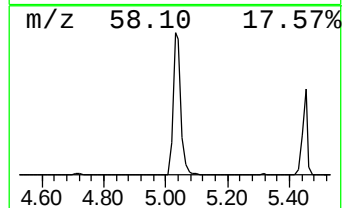
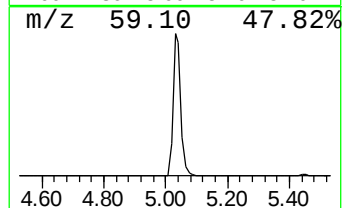
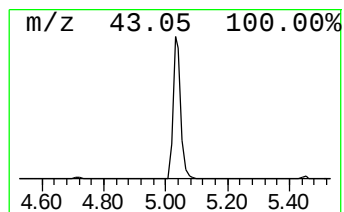
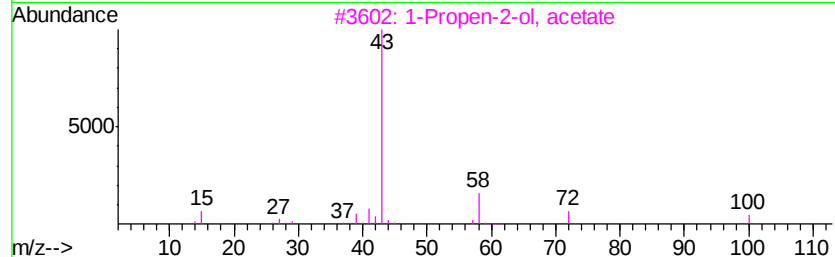
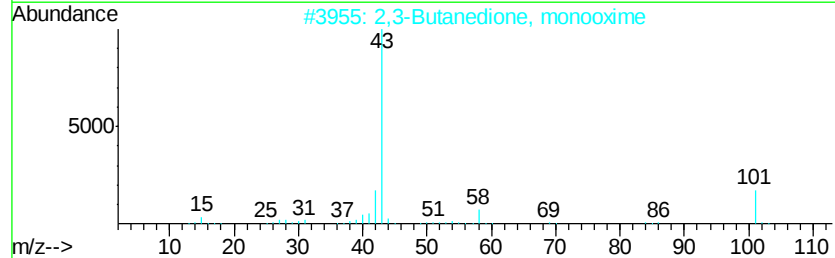
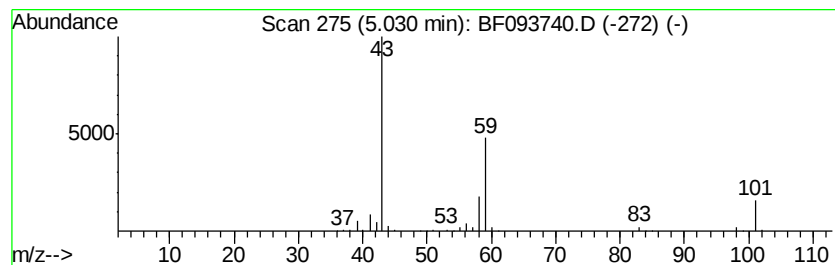
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.03	17.70 ng	983340	1,4-Dichlorobenzene-d4	6.85	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
3	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4	Acetic acid, 2-propenyl ester	100	C5H8O2	000591-87-7	9
5	Formamide, N,N-diethyl-	101	C5H11NO	000617-84-5	7



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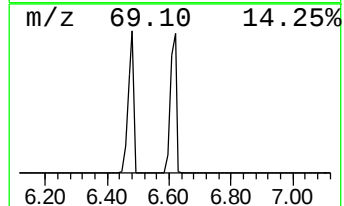
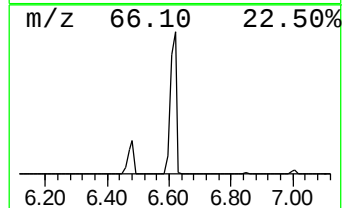
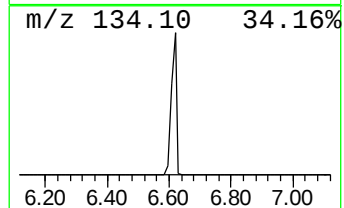
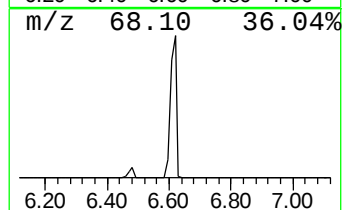
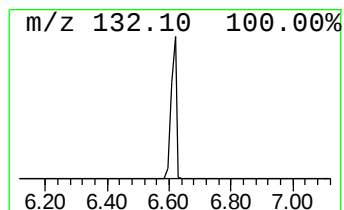
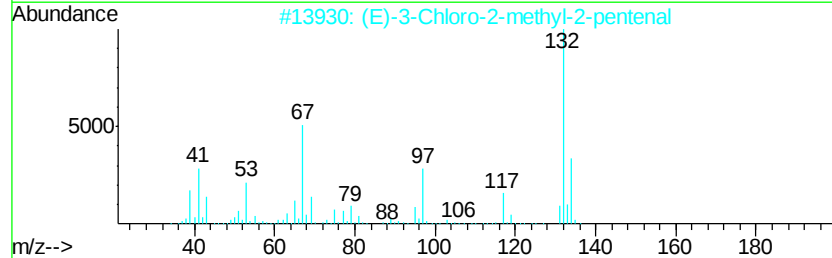
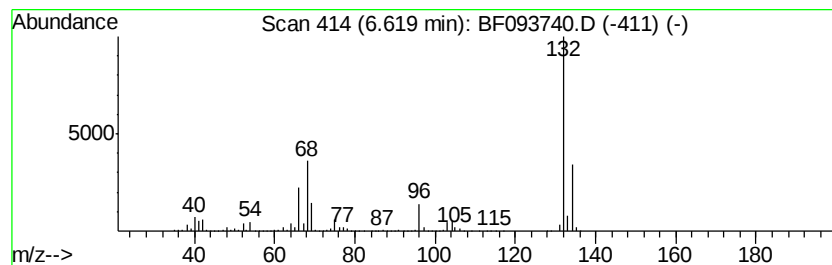
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Peak Number 2 unknown6.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	97.08 ng	5393580	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	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...	5.03	17.7	ng	983340	1	6.85	1111140	20.0
unknown6.62	6.62	97.1	ng	5393580	1	6.85	1111140	20.0