

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082817\
 Data File : BF098096.D
 Acq On : 29 Aug 2017 5:39
 Operator : SJ/JU
 Sample : I4968-02
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 156-275

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-BF081517.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	4.945	482	486	499	rBV	235006	363253	12.49%	1.494%
2	5.357	551	556	559	rBV	2610236	2726164	93.74%	11.212%
3	6.381	725	730	733	rBV	2640697	2614289	89.89%	10.752%
4	6.510	748	752	756	rBV	2718308	2786890	95.83%	11.462%
5	6.739	787	791	795	rBV	809137	627571	21.58%	2.581%
6	6.898	814	818	821	rBV	2461151	2123063	73.00%	8.731%
7	7.304	882	887	890	rBV	1888325	1813584	62.36%	7.459%
8	8.022	1005	1009	1022	rBV	933606	837081	28.78%	3.443%
9	9.045	1179	1183	1187	rBV	60641	71396	2.46%	0.294%
10	9.098	1187	1192	1195	rBV	3045842	2908172	100.00%	11.960%
11	9.774	1302	1307	1320	rVB	892966	841837	28.95%	3.462%
12	10.563	1436	1441	1456	rBV	1764566	1756982	60.42%	7.226%
13	11.257	1555	1559	1566	rBV	995185	848990	29.19%	3.492%
14	12.845	1824	1829	1832	rBV	2706438	2638401	90.72%	10.851%
15	13.886	2002	2006	2019	rBV	729543	699361	24.05%	2.876%
16	15.309	2243	2248	2260	rVB	503628	658032	22.63%	2.706%

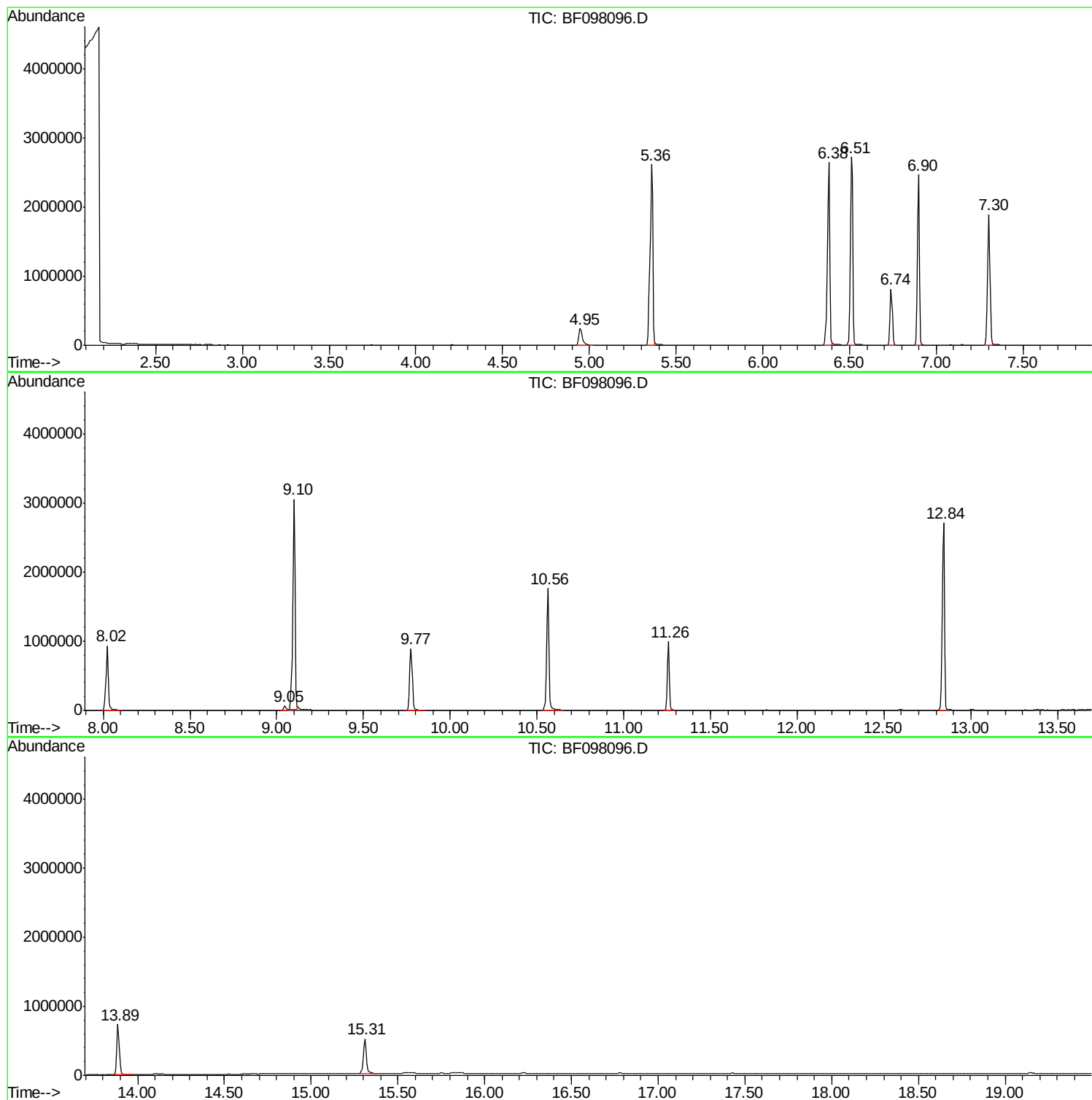
Sum of corrected areas: 24315066

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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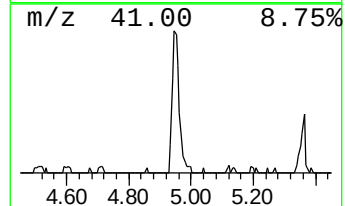
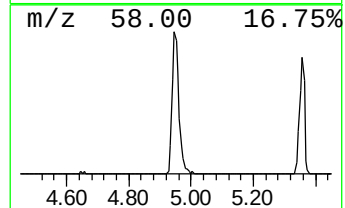
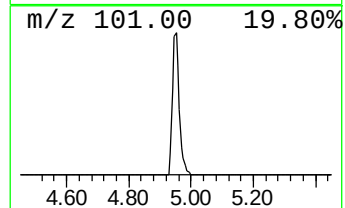
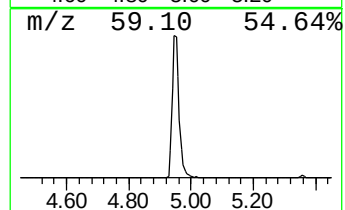
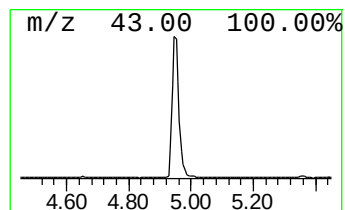
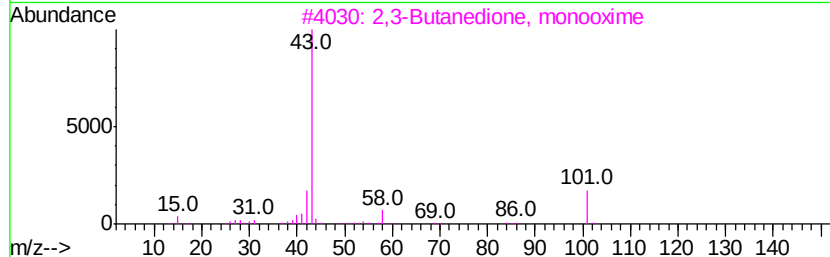
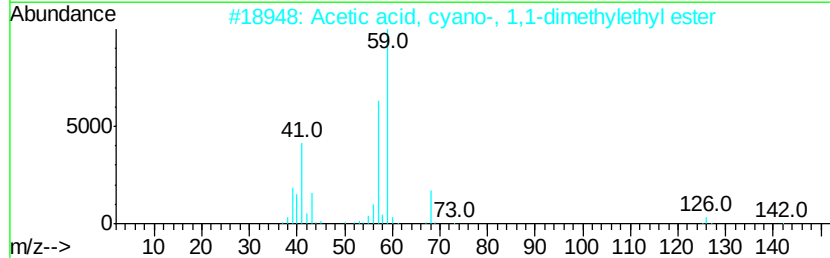
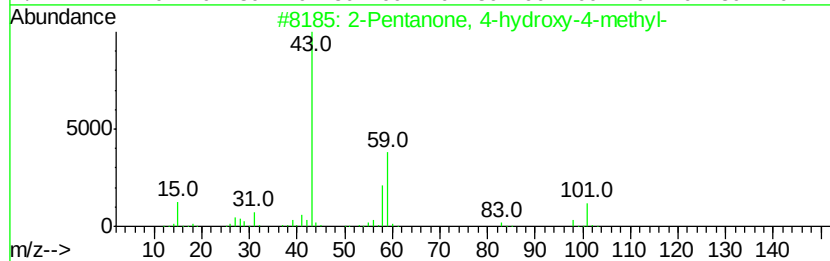
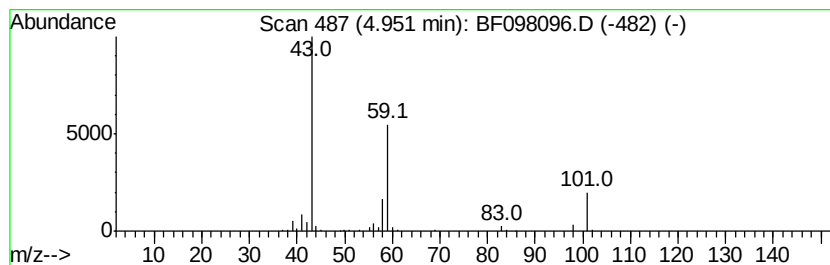
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.95	11.58 ng	363253	1,4-Dichlorobenzene-d4	6.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
4		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	9



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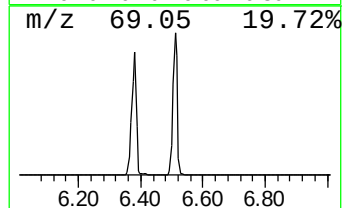
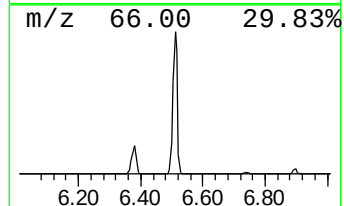
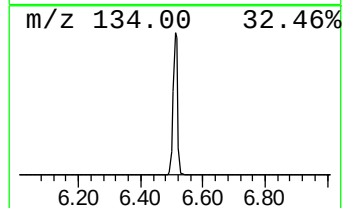
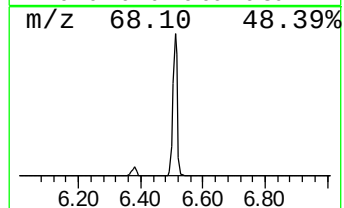
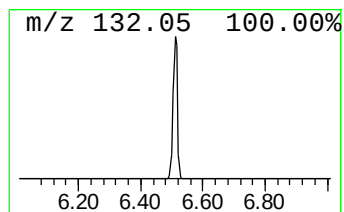
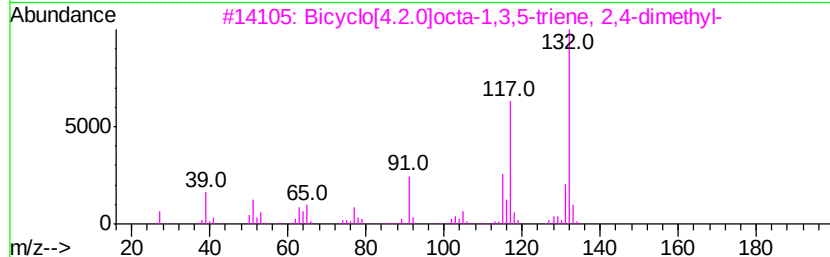
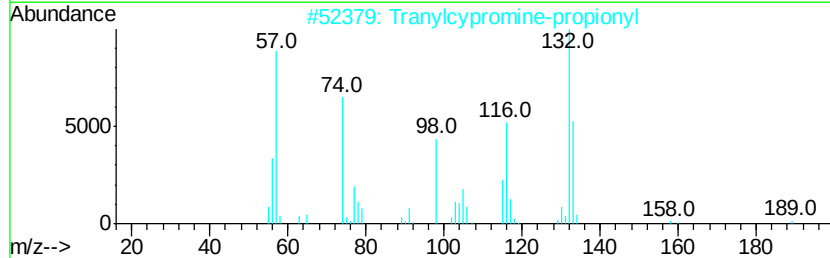
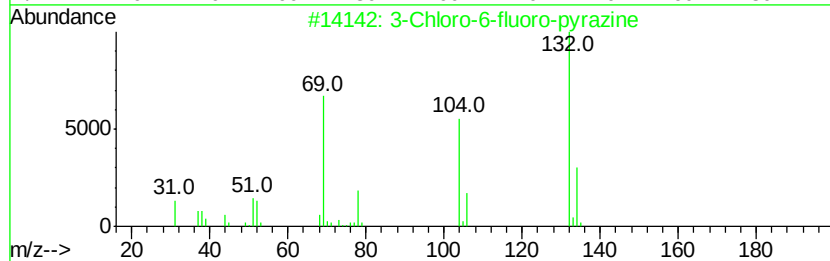
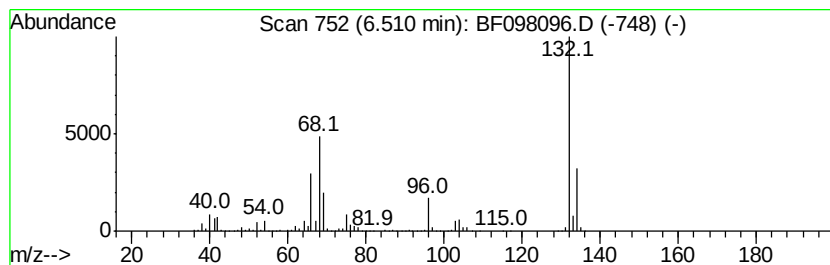
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Peak Number 2 unknown6.51 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.51	88.82 ng	2786890	1,4-Dichlorobenzene-d4	6.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	9
5		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.95	11.6	ng	363253	1	6.74	627571	20.0
unknown6.51	6.51	88.8	ng	2786890	1	6.74	627571	20.0