

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101717\
 Data File : BF099594.D
 Acq On : 17 Oct 2017 21:18
 Operator : SJ/JU
 Sample : I5869-02
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-5S-10-11-17

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-BF092317.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	5.122	499	516	518	rBV	2878367	10027452	100.00%	38.795%
2	6.469	742	745	755	rBV2	358846	412597	4.11%	1.596%
3	6.598	764	767	776	rVB	252251	227873	2.27%	0.882%
4	6.828	802	806	809	rBV	830767	743198	7.41%	2.875%
5	6.981	828	832	836	rBV	2426486	2455688	24.49%	9.501%
6	7.398	897	903	906	rBV	1759991	1919295	19.14%	7.426%
7	8.104	1019	1023	1027	rBV	1069609	1021442	10.19%	3.952%
8	9.186	1201	1207	1210	rBV	3364827	3487265	34.78%	13.492%
9	9.863	1317	1322	1325	rVV	1261065	1081937	10.79%	4.186%
10	11.351	1571	1575	1578	rBV	1048790	914806	9.12%	3.539%
11	12.933	1839	1844	1847	rBV	2307566	2225621	22.20%	8.611%
12	13.992	2020	2024	2027	rBV	661812	631429	6.30%	2.443%
13	15.451	2265	2272	2279	rBV	552865	698681	6.97%	2.703%

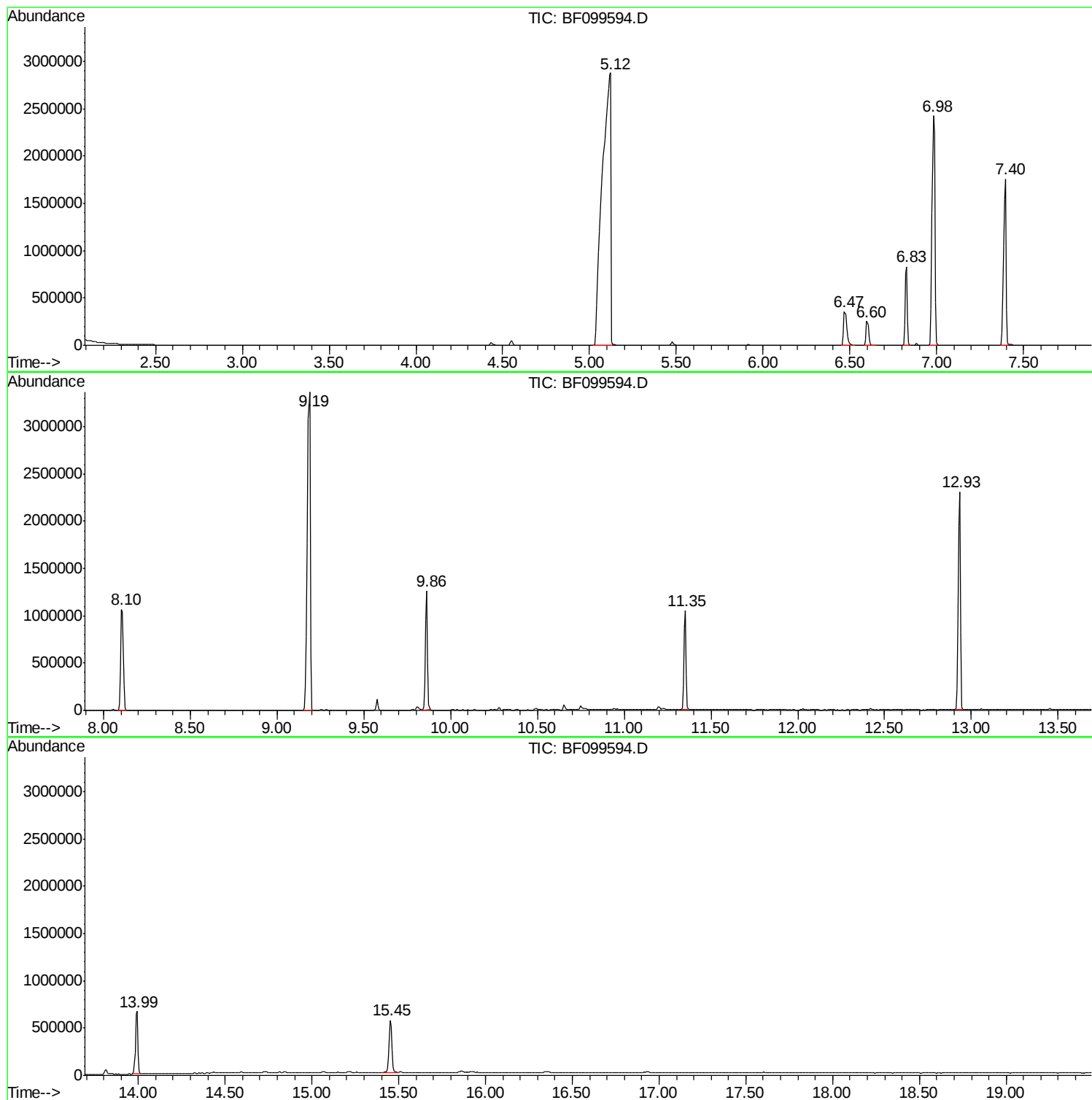
Sum of corrected areas: 25847284

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.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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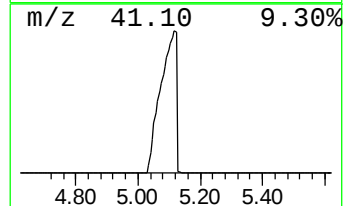
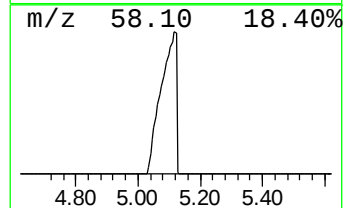
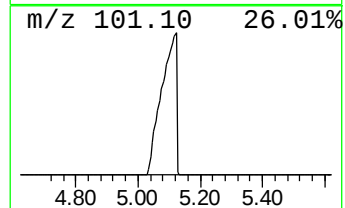
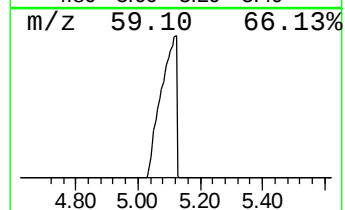
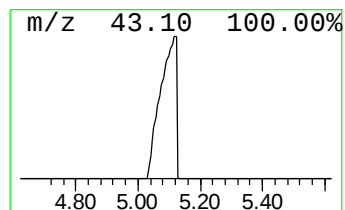
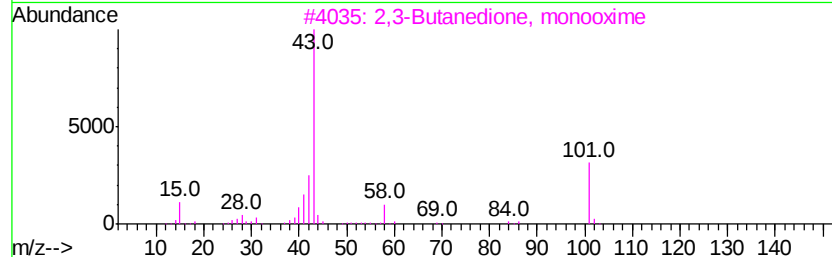
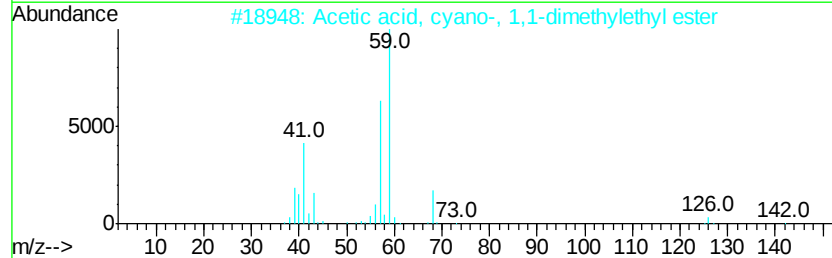
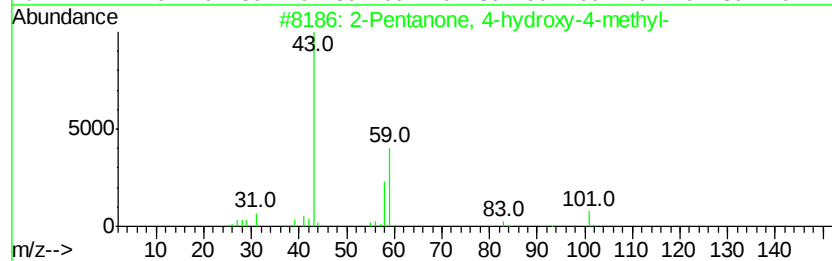
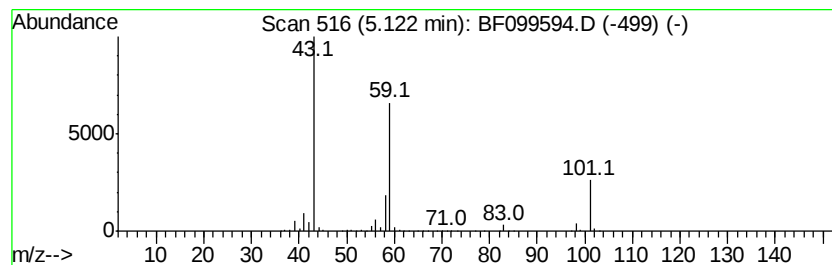
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.12	269.85 ng	10027500	1,4-Dichlorobenzene-d4	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4		5-Hexen-2-one	98	C6H10O	000109-49-9	9
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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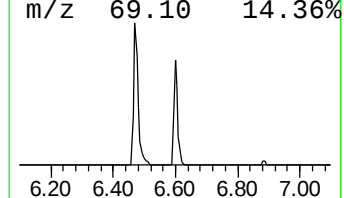
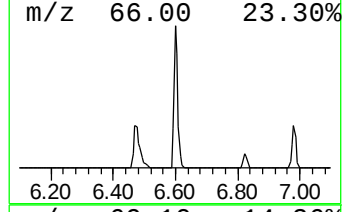
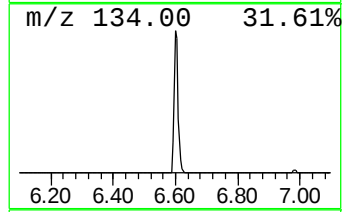
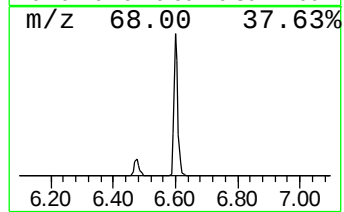
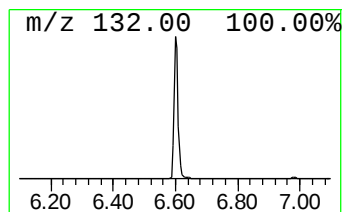
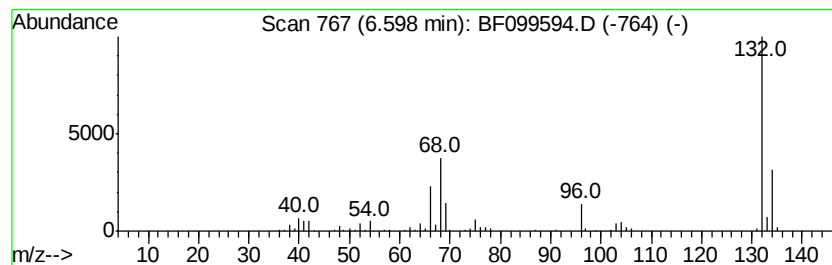
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TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.60 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.60	6.13 ng	227873	1,4-Dichlorobenzene-d4	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.12	269.9	ng	10027500	1	6.83	743198	20.0
unknown6.60	6.60	6.1	ng	227873	1	6.83	743198	20.0