

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061116\
 Data File : BF087911.D
 Acq On : 11 Jun 2016 19:34
 Operator : UM/SJ
 Sample : H3387-04
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP16-LF2MW4-9007

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-BF060716.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	1.735	11	13	22	rVV	247727	418489	11.63%	1.617%
2	1.861	22	24	37	rVB	2135770	3388550	94.18%	13.093%
3	4.959	292	295	299	rBV	77304	85993	2.39%	0.332%
4	5.381	330	332	335	rBV	930726	1165086	32.38%	4.502%
5	6.422	421	423	426	rBV	526632	680611	18.92%	2.630%
6	6.570	433	436	438	rBV	2067436	2427467	67.47%	9.379%
7	6.799	454	456	458	rBV	746633	621115	17.26%	2.400%
8	6.959	467	470	472	rVB	2431662	2290403	63.66%	8.850%
9	7.370	503	506	508	rBV	1882878	2040361	56.71%	7.884%
10	8.090	566	569	571	rBV	959853	878305	24.41%	3.394%
11	9.176	660	664	666	rBV	2627279	3597860	100.00%	13.902%
12	9.839	720	722	725	rBB	816788	799264	22.21%	3.088%
13	10.628	788	791	794	rBV2	1500227	2020462	56.16%	7.807%
14	11.325	849	852	854	rBV	1126334	979925	27.24%	3.786%
15	12.914	988	991	994	rBV	2605292	2912463	80.95%	11.253%
16	13.954	1080	1082	1085	rBV	875491	862634	23.98%	3.333%
17	15.371	1203	1206	1209	rBV	582079	711979	19.79%	2.751%

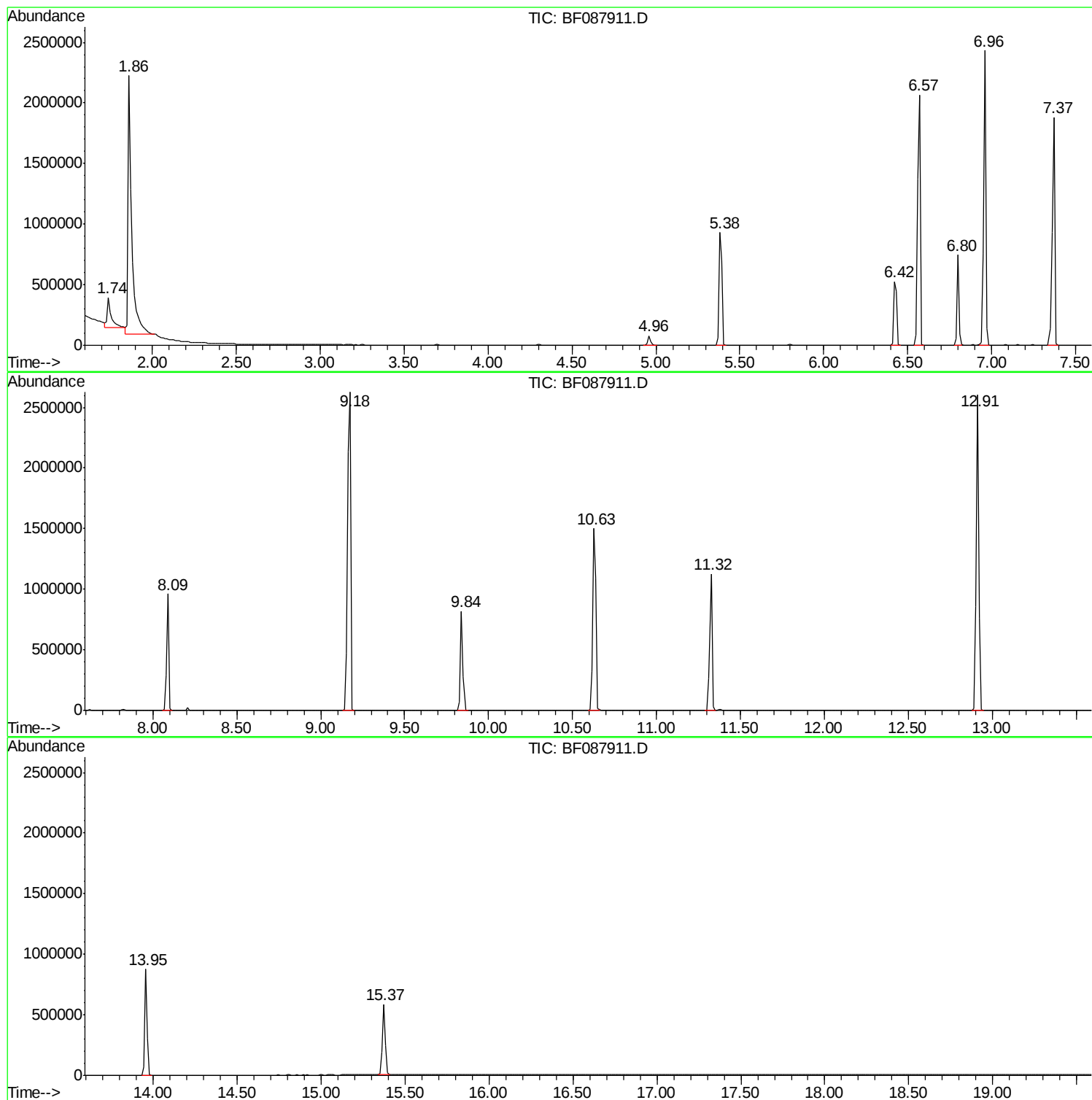
Sum of corrected areas: 25880967

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.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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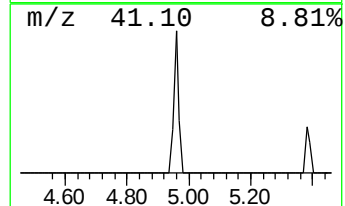
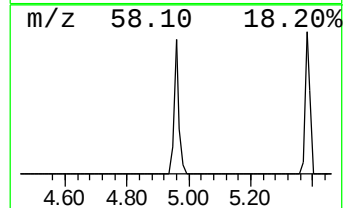
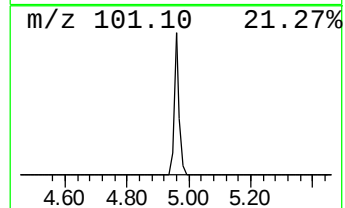
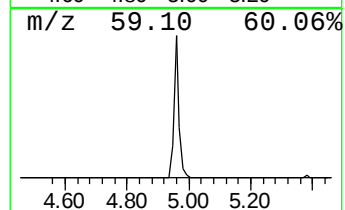
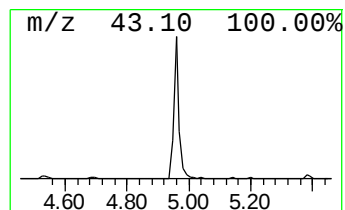
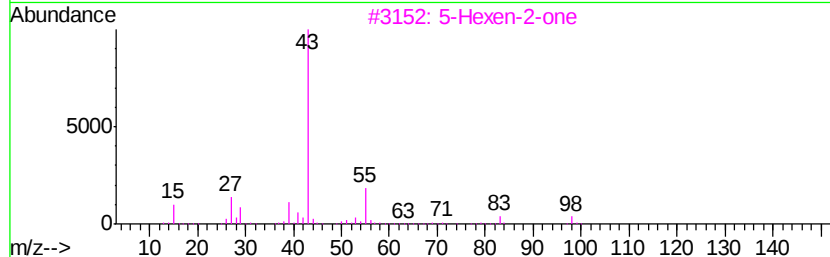
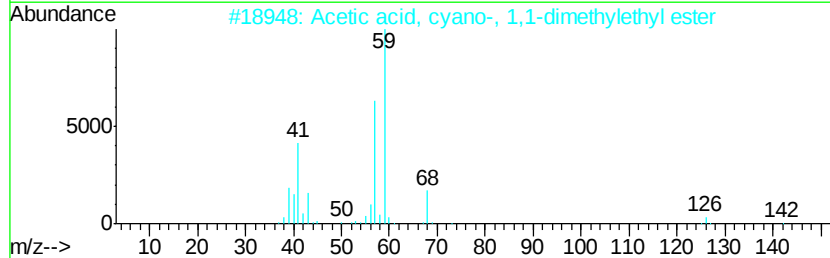
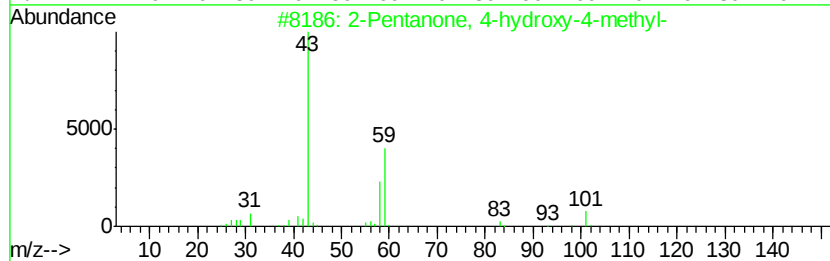
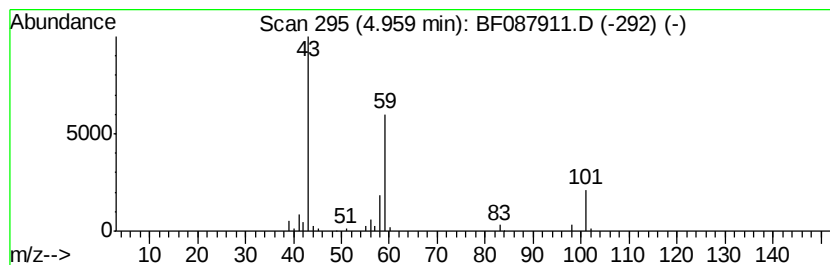
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Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.96	2.77 ng	85993	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	56
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
3		5-Hexen-2-one	98	C6H10O	000109-49-9	10
4		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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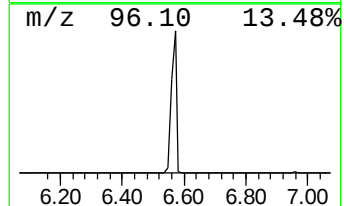
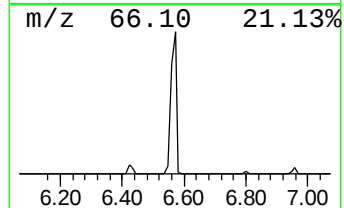
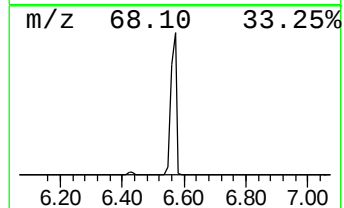
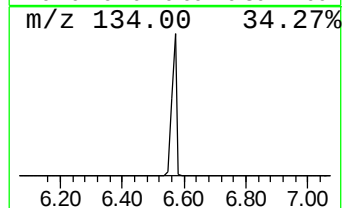
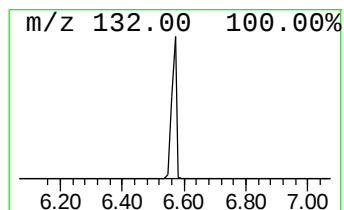
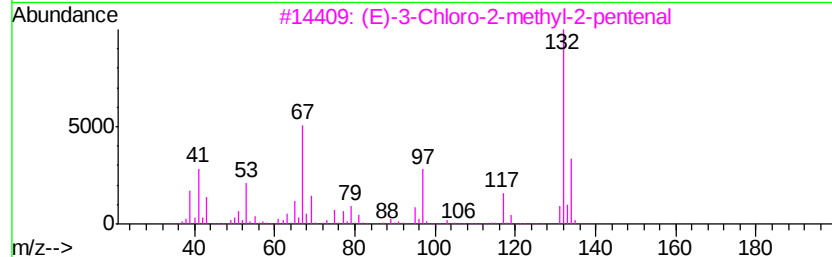
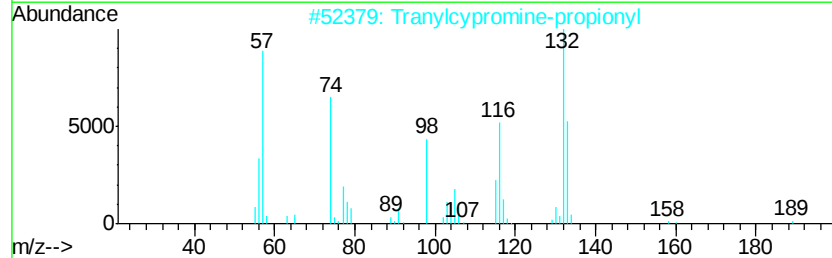
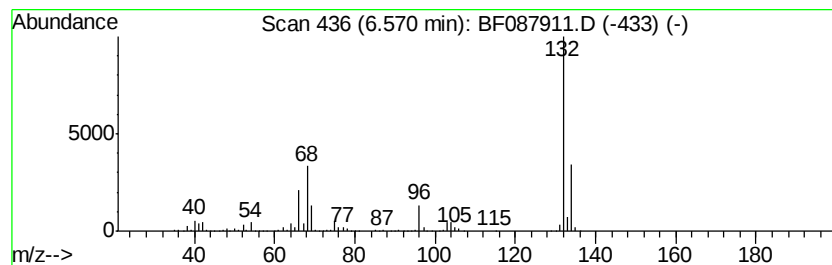
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Peak Number 4 unknown6.57 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.57	78.16 ng	2427470	1,4-Dichlorobenzene-d4	6.80

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
4			2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
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...	4.96	2.8	ng	85993	1	6.80	621115	20.0
unknown6.57	6.57	78.2	ng	2427470	1	6.80	621115	20.0