

Data Path : U:\HPCHEM1\BNA_F\DATA\BF011518\
 Data File : BF102077.D
 Acq On : 15 Jan 2018 21:47
 Operator : SJ/JU
 Sample : PB105648BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB105648BL

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-BF010918.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	4.934	479	484	491	rBV	217001	278103	5.30%	0.767%
2	5.334	547	552	555	rBV	3600944	3726751	71.07%	10.279%
3	6.357	720	726	729	rBV	3444242	3583563	68.34%	9.884%
4	6.487	744	748	751	rBV	3804468	3720481	70.95%	10.262%
5	6.722	784	788	791	rBV	1359590	1205481	22.99%	3.325%
6	6.875	810	814	818	rVB	4036806	4056750	77.36%	11.189%
7	7.287	878	884	887	rBV	3225957	3412411	65.07%	9.412%
8	7.998	1000	1005	1009	rBV	1615108	1543910	29.44%	4.258%
9	9.081	1182	1189	1192	rBV	5383897	5244082	100.00%	14.464%
10	9.751	1298	1303	1307	rBV	1567241	1458300	27.81%	4.022%
11	10.539	1431	1437	1440	rBV	1936714	1699115	32.40%	4.686%
12	11.233	1551	1555	1559	rBV	1407339	1148134	21.89%	3.167%
13	12.698	1798	1804	1807	rBV	72873	93707	1.79%	0.258%
14	12.822	1820	1825	1828	rBV	3394163	3171428	60.48%	8.747%
15	13.863	1998	2002	2006	rBV	1045073	988306	18.85%	2.726%
16	15.280	2238	2243	2247	rBV	784173	925458	17.65%	2.553%

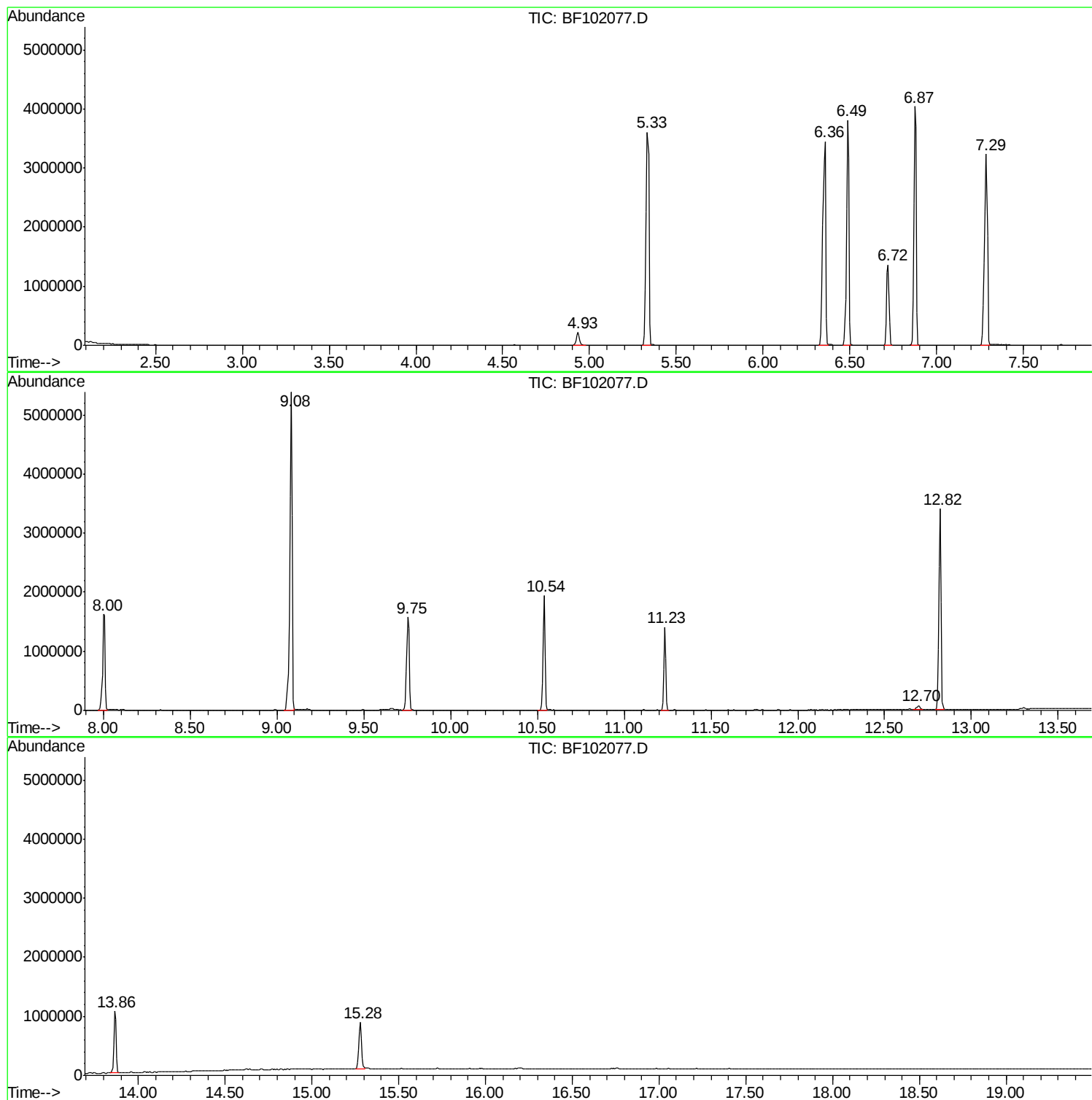
Sum of corrected areas: 36255980

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF010918.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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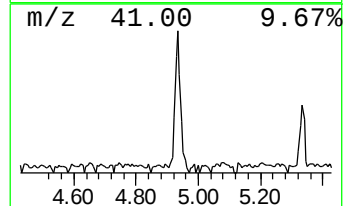
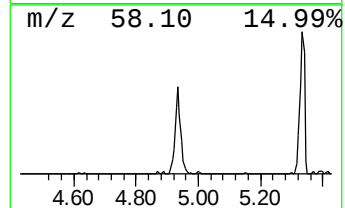
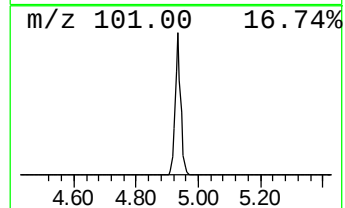
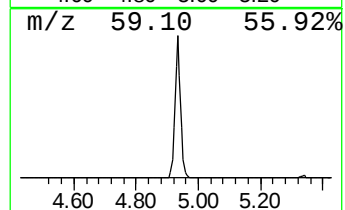
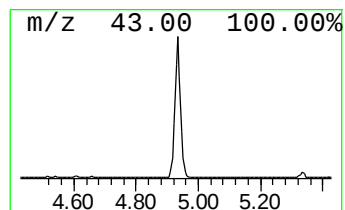
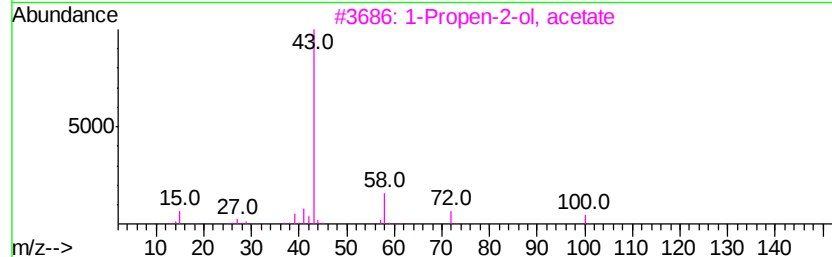
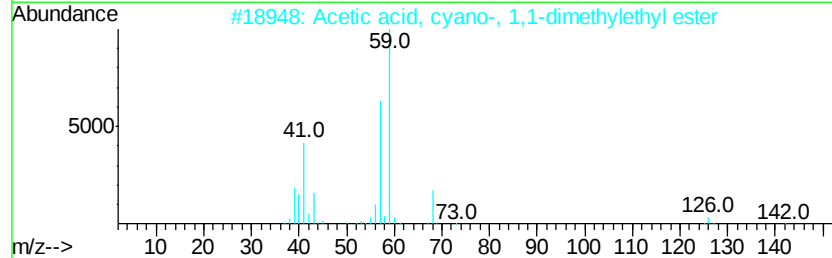
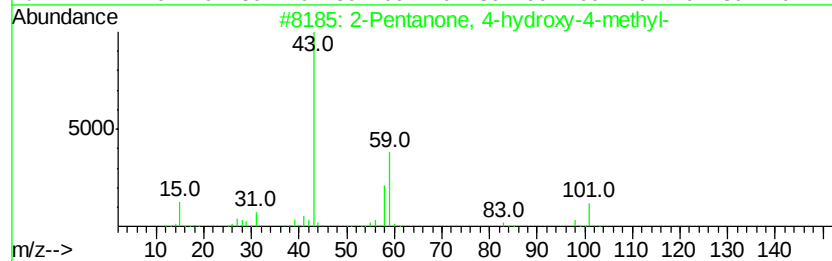
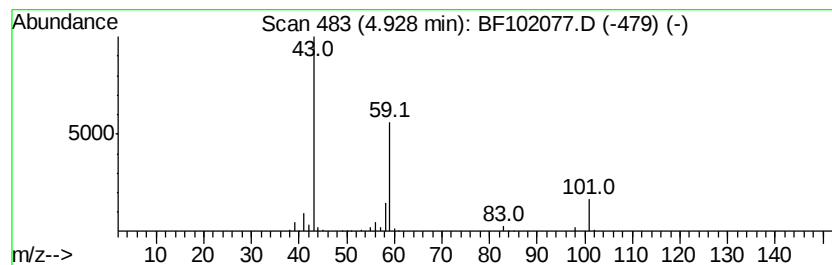
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.93	4.61 ng	278103	1,4-Dichlorobenzene-d4	6.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
3			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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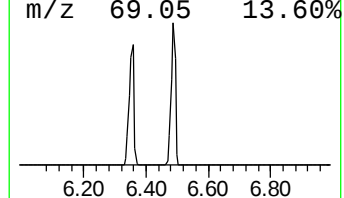
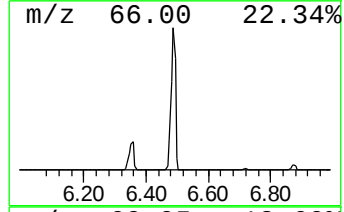
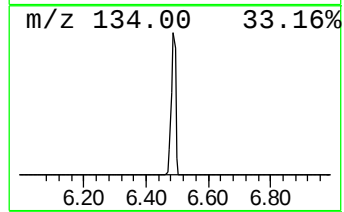
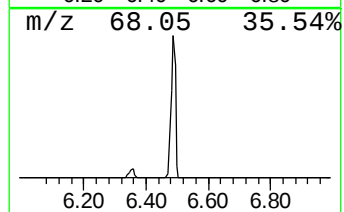
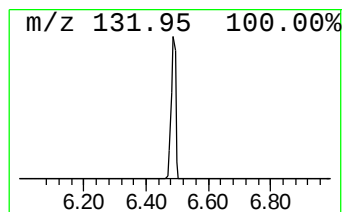
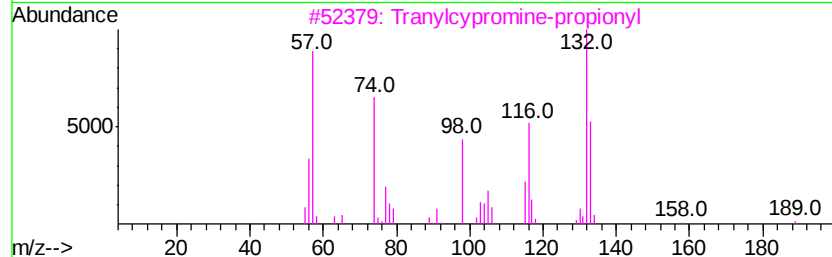
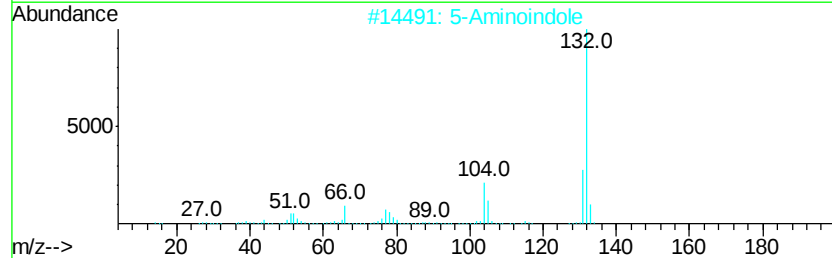
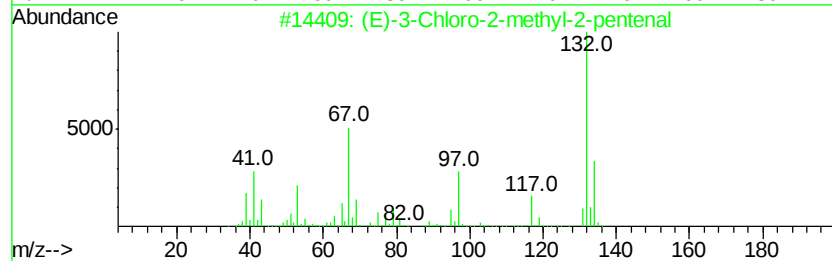
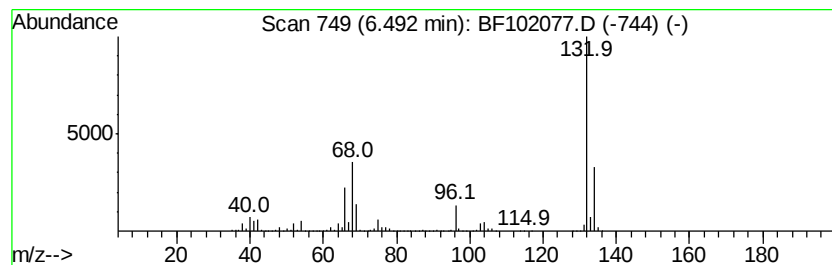
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Peak Number 2 unknown6.49 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.49	61.73 ng	3720480	1,4-Dichlorobenzene-d4	6.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
2	5-Aminoindole	132	C8H8N2	005192-03-0	12
3	Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4	Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5	1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.93	4.6	ng	278103	1	6.72	1205480	20.0
unknown6.49	6.49	61.7	ng	3720480	1	6.72	1205480	20.0