

Data Path : U:\HPCHEM1\BNA F\DATA\BF082217\
 Data File : BF097961.D
 Acq On : 22 Aug 2017 18:44
 Operator : SJ/JU
 Sample : PB101706BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB101706BL

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	496	rVB	278221	360232	10.78%	1.439%
2	5.357	552	556	560	rBV	2011739	2052713	61.45%	8.200%
3	6.380	724	730	733	rBV	2003974	2080101	62.26%	8.309%
4	6.516	748	753	756	rBV	2180626	2073226	62.06%	8.282%
5	6.745	788	792	796	rBV	958330	764532	22.89%	3.054%
6	6.904	814	819	822	rBV	2455892	2277909	68.19%	9.099%
7	7.310	882	888	891	rBV	1879146	1906301	57.06%	7.615%
8	8.027	1005	1010	1014	rBV	1098035	1029408	30.81%	4.112%
9	9.104	1187	1193	1197	rBV	2982286	3323551	99.49%	13.276%
10	9.780	1303	1308	1312	rBV2	1215993	1123929	33.64%	4.490%
11	10.568	1436	1442	1445	rBV	1418689	1381806	41.36%	5.520%
12	11.263	1555	1560	1563	rBV	1400000	1209312	36.20%	4.831%
13	12.851	1825	1830	1834	rBV	2935063	3340726	100.00%	13.345%
14	13.892	2003	2007	2011	rBV	1147025	1137048	34.04%	4.542%
15	15.315	2243	2249	2253	rBV	644974	771754	23.10%	3.083%
16	15.356	2253	2256	2262	rVB	28061	36902	1.10%	0.147%
17	15.768	2322	2326	2333	rVB2	24712	36909	1.10%	0.147%
18	16.245	2402	2407	2416	rVB2	27202	48893	1.46%	0.195%
19	16.803	2497	2502	2512	rVB2	20293	41608	1.25%	0.166%
20	17.456	2608	2613	2624	rVB6	15007	37435	1.12%	0.150%

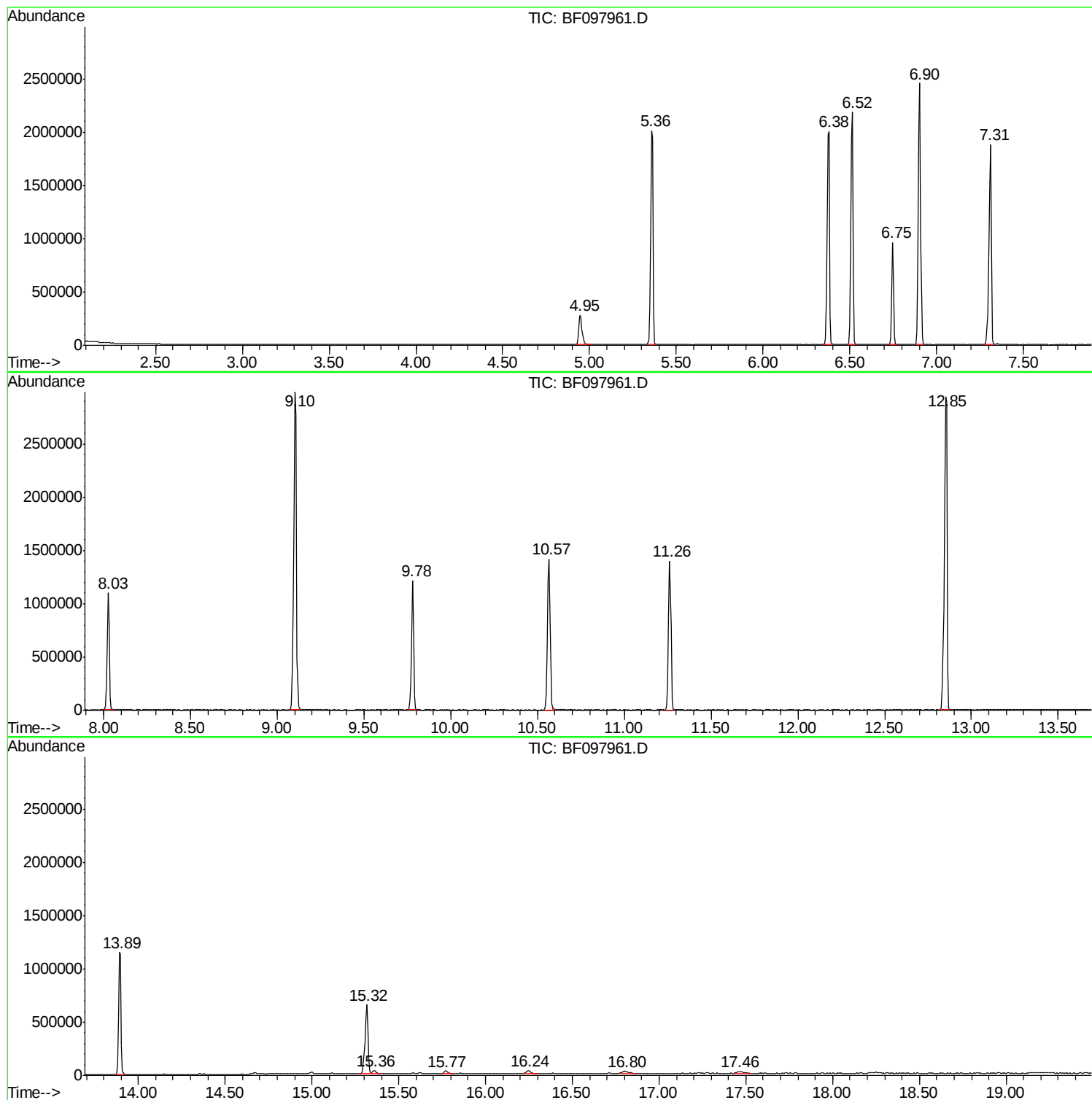
Sum of corrected areas: 25034295

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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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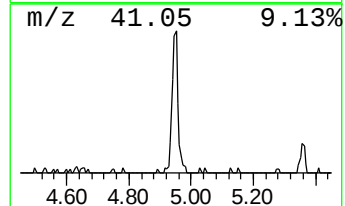
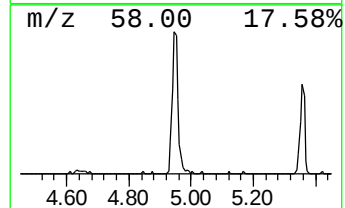
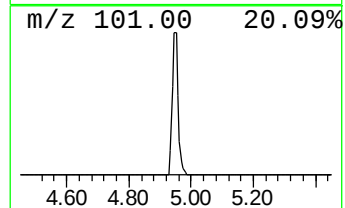
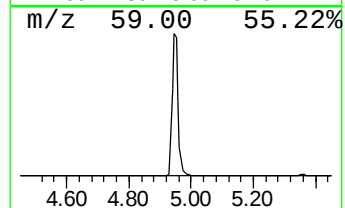
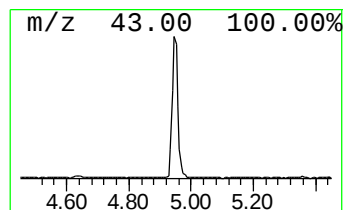
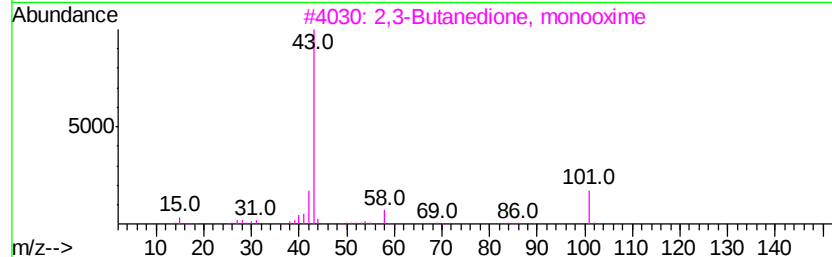
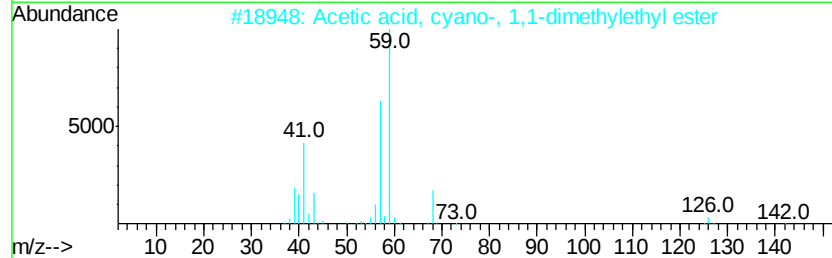
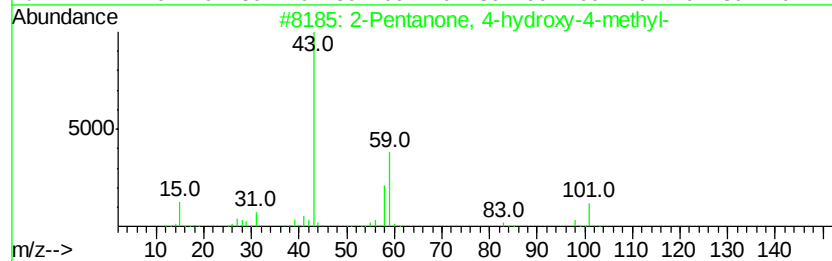
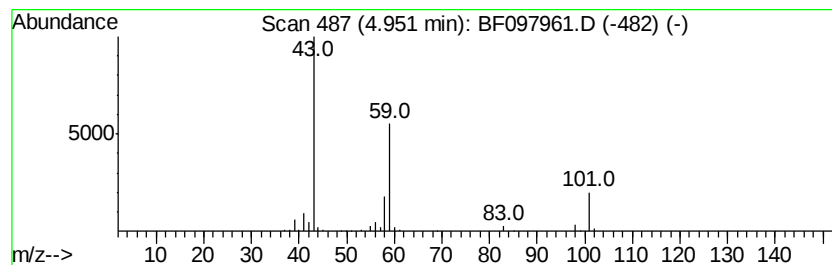
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.95	9.42 ng	360232	1,4-Dichlorobenzene-d4	6.75

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		Butane	58	C4H10	000106-97-8	9



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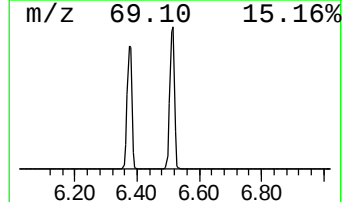
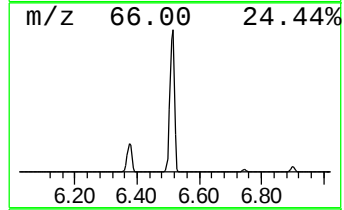
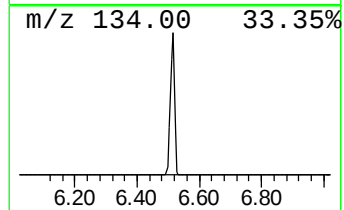
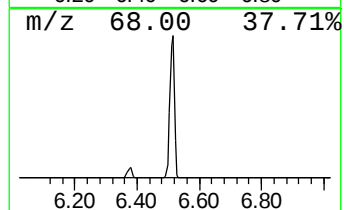
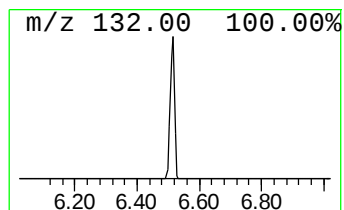
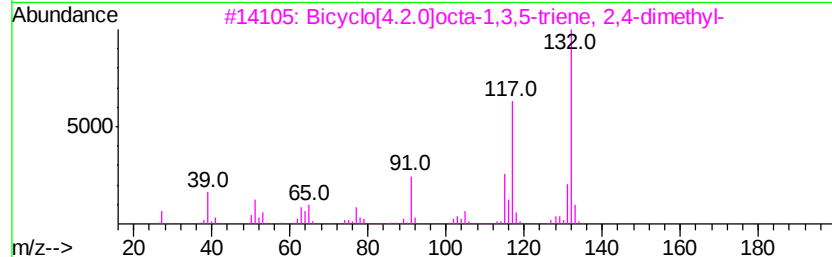
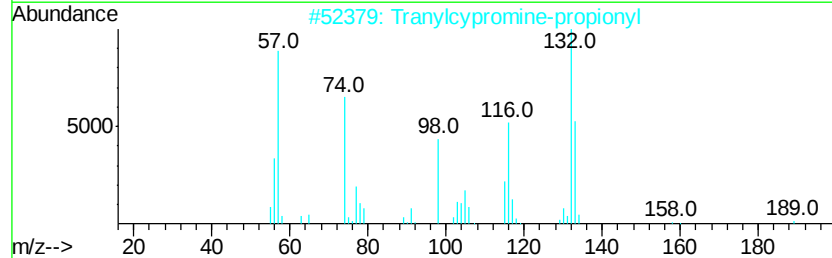
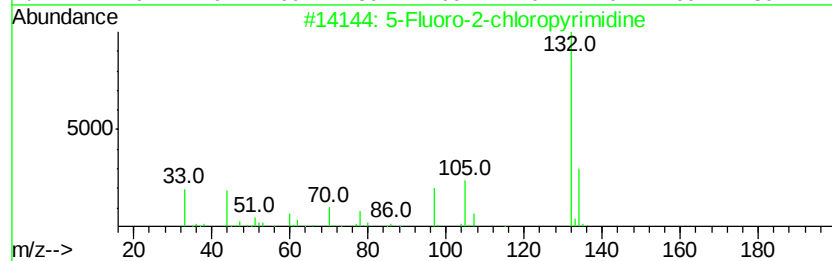
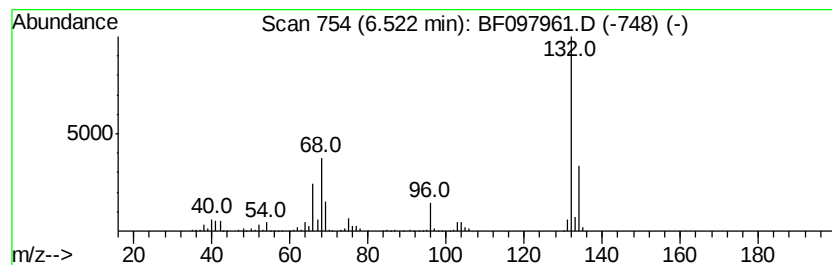
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Peak Number 2 unknown6.52 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.52	54.24 ng	2073230	1,4-Dichlorobenzene-d4	6.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	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		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.95	9.4	ng	360232	1	6.75	764532	20.0
unknown6.52	6.52	54.2	ng	2073230	1	6.75	764532	20.0