

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF090717\
 Data File : BF098343.D
 Acq On : 8 Sep 2017 5:33
 Operator : SJ/JU
 Sample : I5132-02
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 090617-TIDO-F-D-S

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	4.857	468	471	483	rVB	150435	168189	5.21%	0.865%
2	5.293	541	545	557	rBV	1035283	1070780	33.15%	5.508%
3	6.328	717	721	732	rBV2	982220	839337	25.99%	4.317%
4	6.457	739	743	746	rBV	2066631	1773732	54.91%	9.123%
5	6.687	779	782	786	rBV	783252	631067	19.54%	3.246%
6	6.845	805	809	812	rBV	2590490	2327167	72.05%	11.970%
7	7.257	874	879	882	rBV	1824729	1914888	59.28%	9.849%
8	7.969	997	1000	1004	rBV	895474	797064	24.68%	4.100%
9	9.051	1179	1184	1187	rBV	3291365	3230005	100.00%	16.614%
10	9.445	1248	1251	1254	rBV	50503	39113	1.21%	0.201%
11	9.722	1294	1298	1302	rBV	882736	816609	25.28%	4.200%
12	10.516	1428	1433	1436	rBV	1156506	1110640	34.39%	5.713%
13	11.204	1547	1550	1554	rBV	825876	773538	23.95%	3.979%
14	12.616	1784	1790	1793	rBV	48528	36995	1.15%	0.190%
15	12.798	1816	1821	1824	rBV	2652942	2677022	82.88%	13.770%
16	13.698	1970	1974	1982	rBV2	19905	33014	1.02%	0.170%
17	13.839	1994	1998	2002	rBV	764721	697356	21.59%	3.587%
18	15.251	2233	2238	2246	rBV	408915	505042	15.64%	2.598%

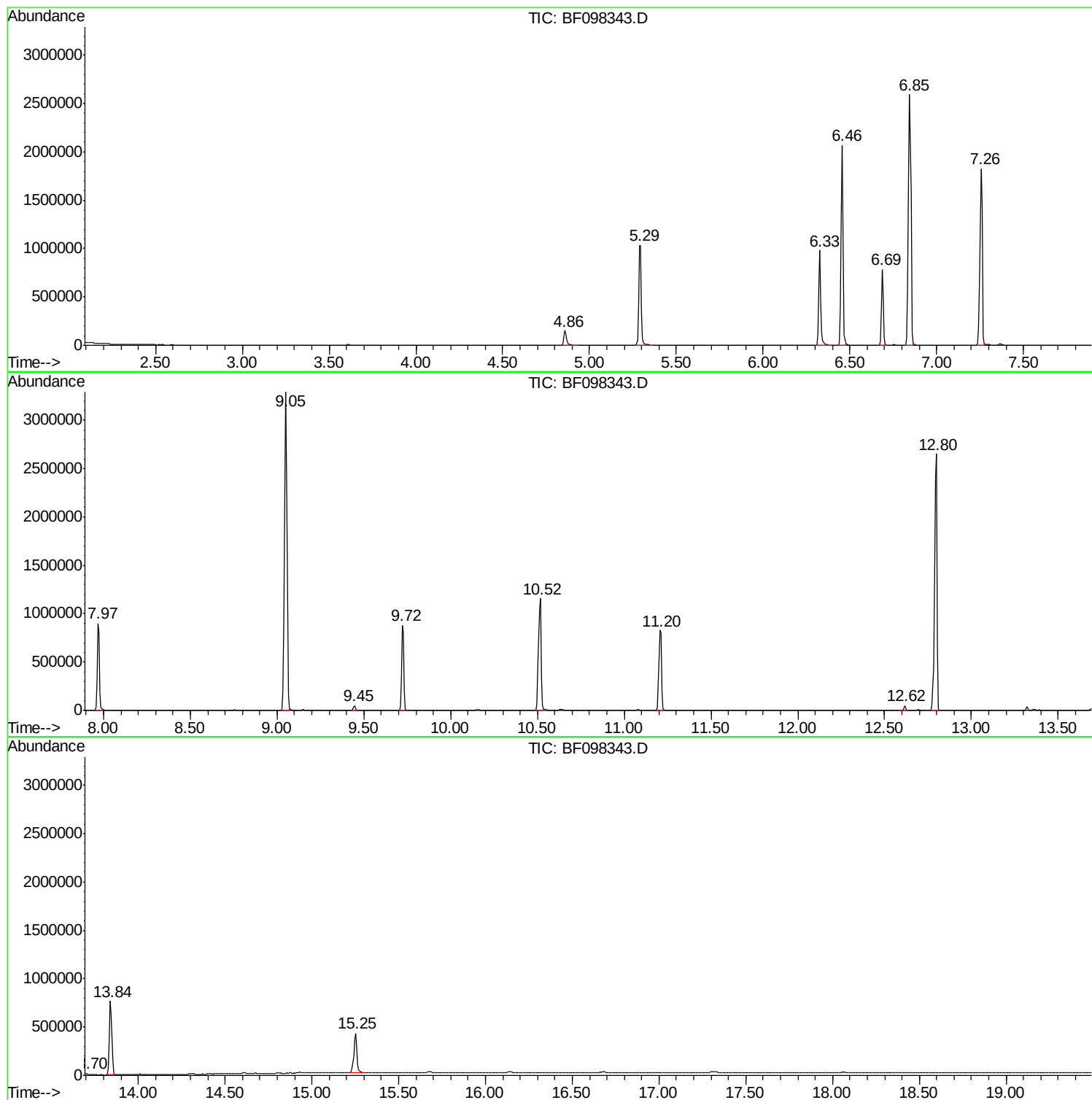
Sum of corrected areas: 19441558

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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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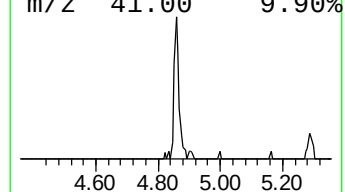
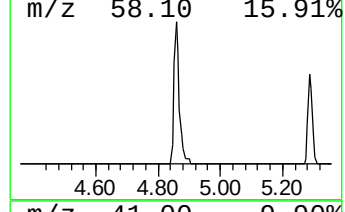
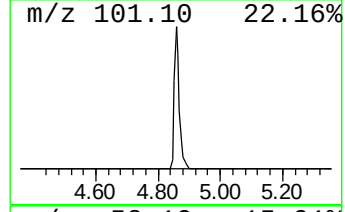
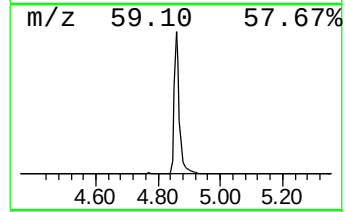
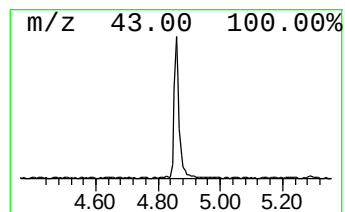
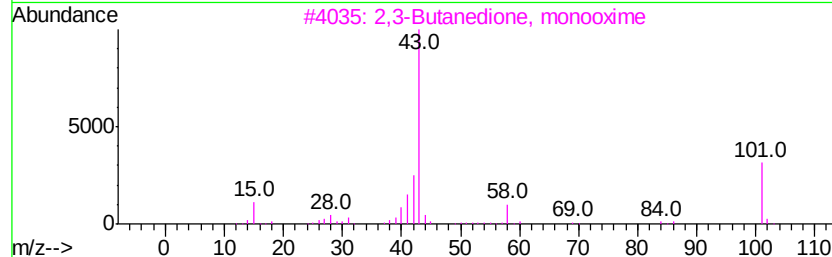
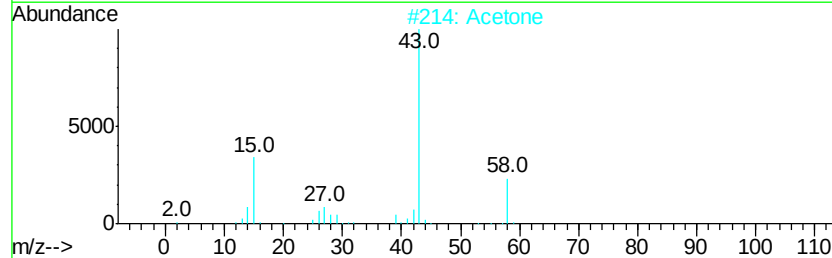
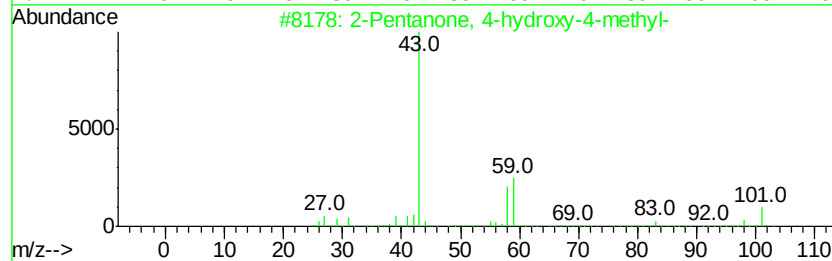
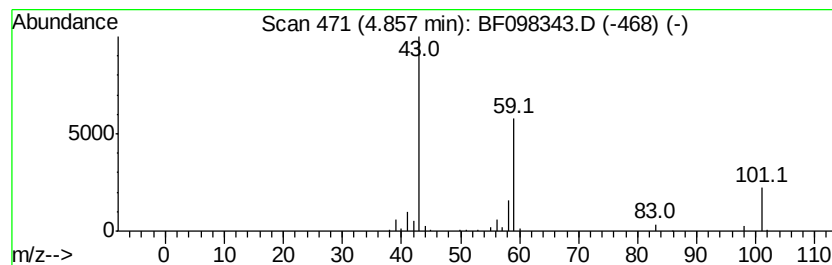
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
4.86	5.33 ng	168189	1,4-Dichlorobenzene-d4	6.69	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	Acetone	58	C3H6O	000067-64-1	10
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4	2-Propanone, 1-(1-methylethoxy)-	116	C6H12O2	042781-12-4	9
5	3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	9



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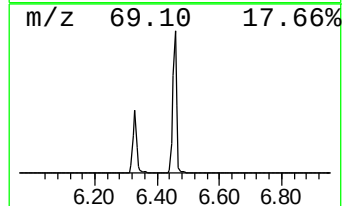
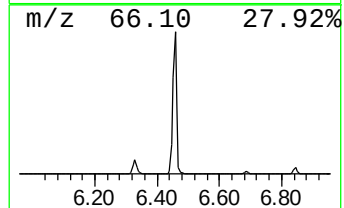
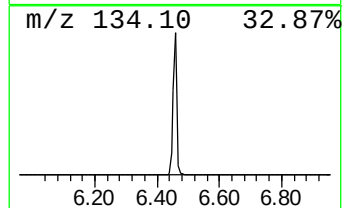
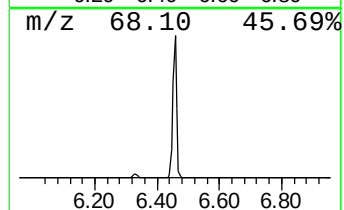
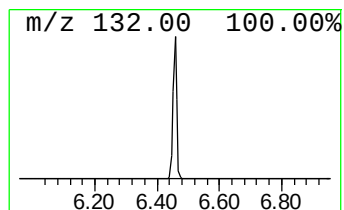
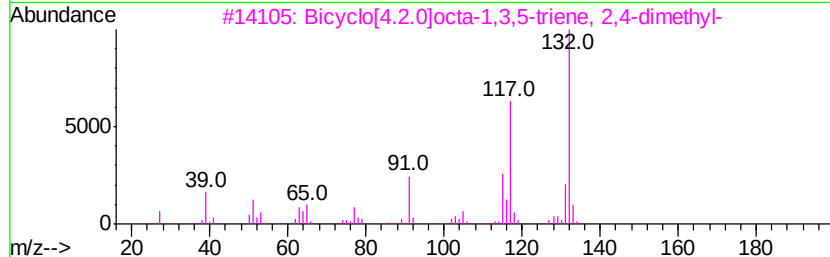
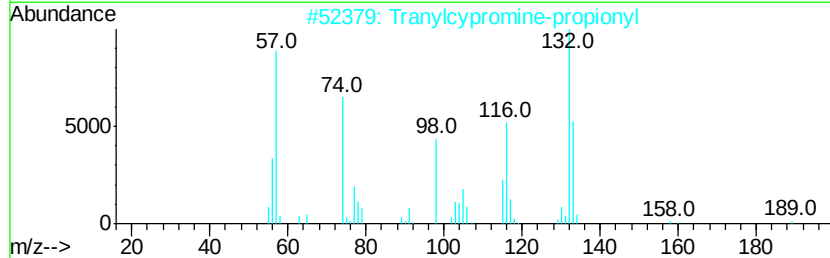
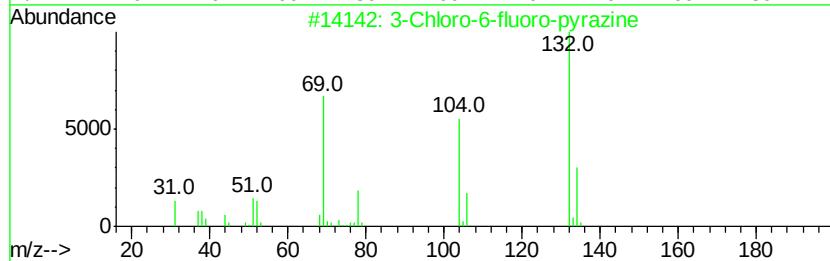
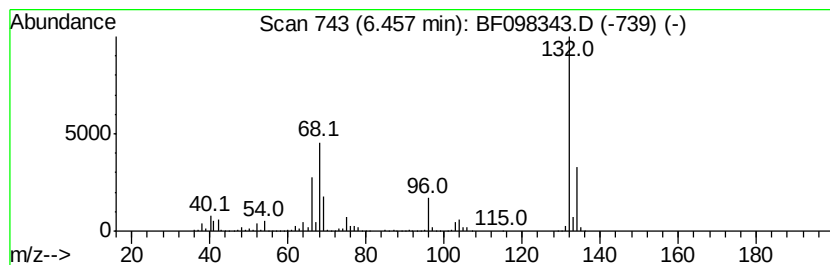
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Peak Number 2 unknown6.46 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.46	56.21 ng	1773730	1,4-Dichlorobenzene-d4	6.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
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		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.86	5.3	ng	168189	1	6.69	631067	20.0
unknown6.46	6.46	56.2	ng	1773730	1	6.69	631067	20.0