

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090817\
 Data File : BF098374.D
 Acq On : 9 Sep 2017 4:36
 Operator : SJ/JU
 Sample : PB102179BL
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB102179BL

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	4.881	471	475	486	rBV	123621	166533	7.52%	0.933%
2	5.304	543	547	562	rBV	2053997	2100832	94.87%	11.773%
3	6.340	718	723	726	rBV	2178784	2183934	98.62%	12.239%
4	6.457	739	743	746	rBV	2209996	2062207	93.13%	11.557%
5	6.681	778	781	785	rBV	671290	556473	25.13%	3.118%
6	6.840	803	808	811	rBV	1984123	1803441	81.44%	10.106%
7	7.251	873	878	881	rBV	1473850	1365171	61.65%	7.650%
8	7.963	995	999	1003	rBV	720688	648309	29.28%	3.633%
9	9.045	1177	1183	1186	rBV	2515914	2214424	100.00%	12.410%
10	9.716	1293	1297	1301	rBV2	694789	551380	24.90%	3.090%
11	10.504	1428	1431	1437	rBV	726510	703575	31.77%	3.943%
12	11.198	1545	1549	1553	rBV	614348	503242	22.73%	2.820%
13	12.610	1785	1789	1792	rBV2	34167	33354	1.51%	0.187%
14	12.674	1796	1800	1804	rVV5	15370	27013	1.22%	0.151%
15	12.786	1814	1819	1822	rBV	2053465	1890625	85.38%	10.595%
16	13.316	1906	1909	1912	rBV3	28733	31673	1.43%	0.177%
17	13.833	1993	1997	2001	rBV	697958	590629	26.67%	3.310%
18	15.239	2232	2236	2242	rBV	341457	411726	18.59%	2.307%

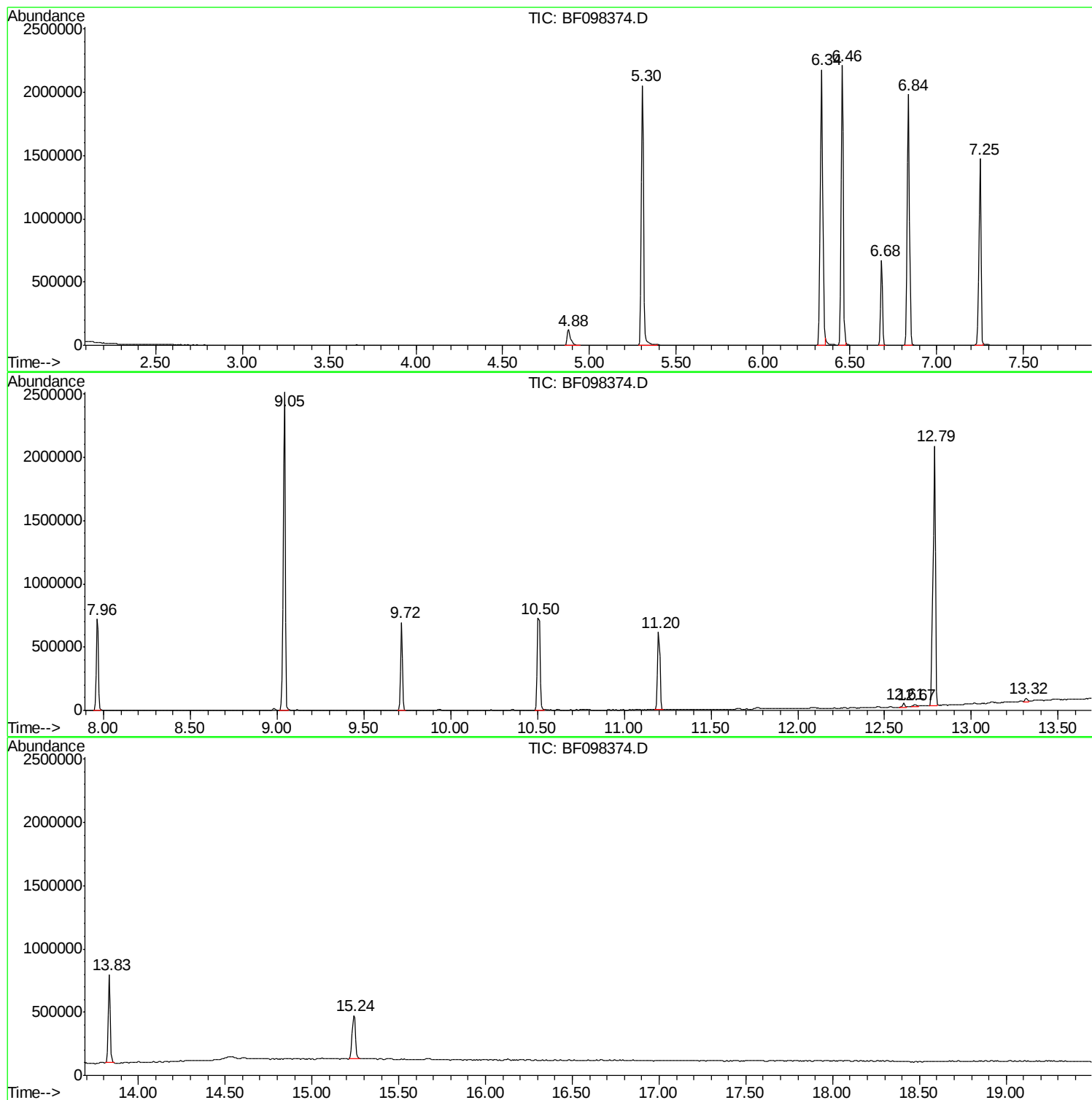
Sum of corrected areas: 17844541

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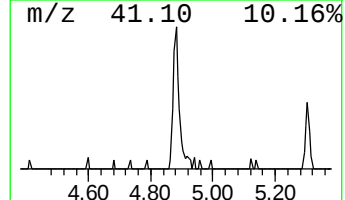
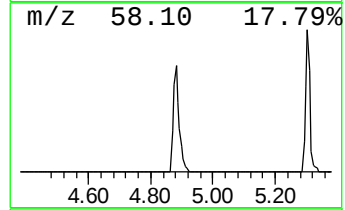
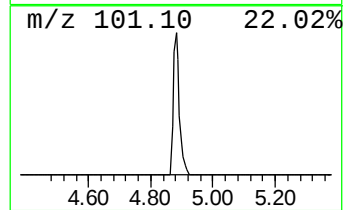
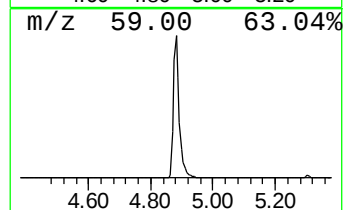
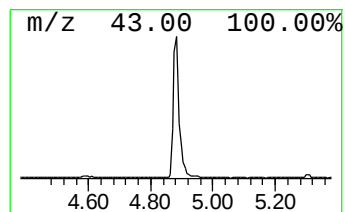
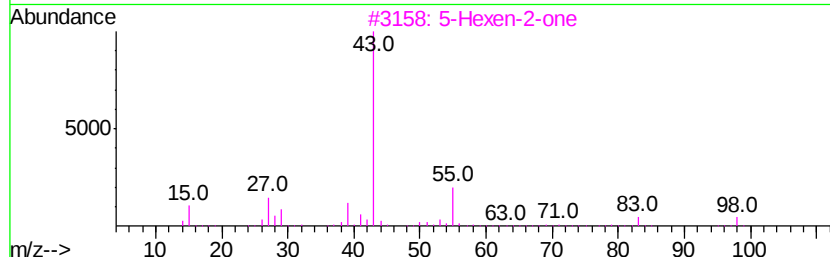
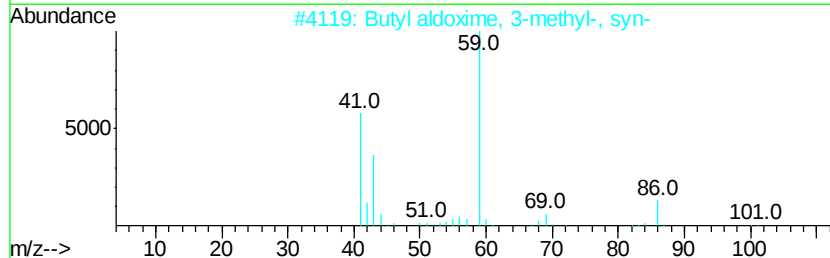
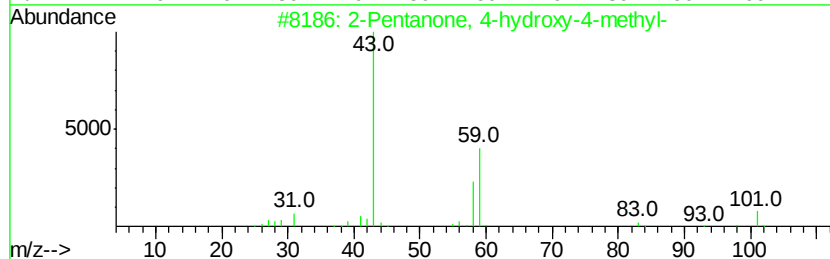
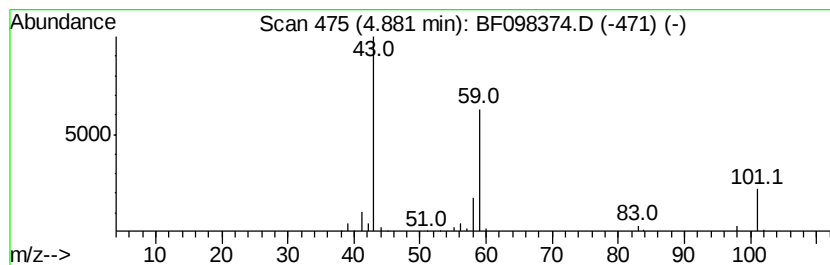
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.88	5.99 ng	166533	1,4-Dichlorobenzene-d4	6.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	25
3		5-Hexen-2-one	98	C6H10O	000109-49-9	9
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		Acetone	58	C3H6O	000067-64-1	9



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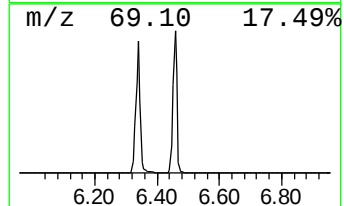
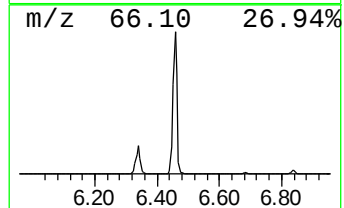
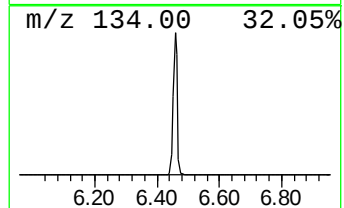
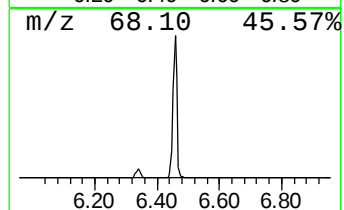
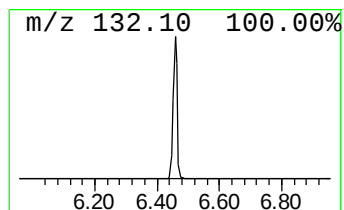
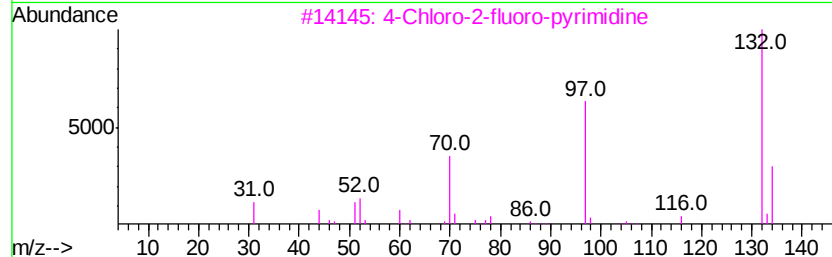
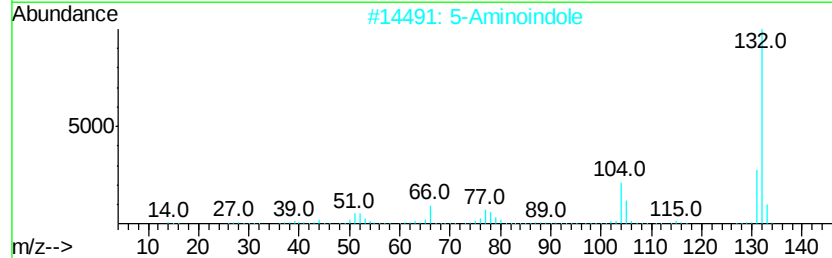
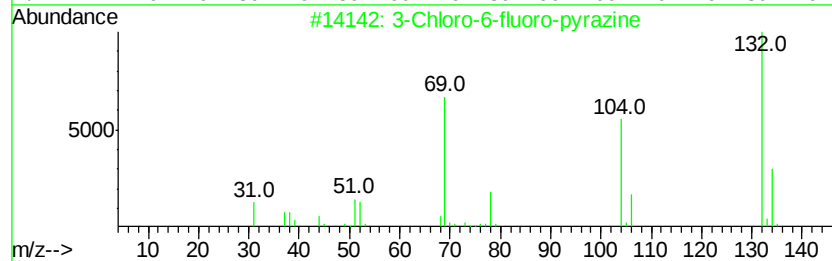
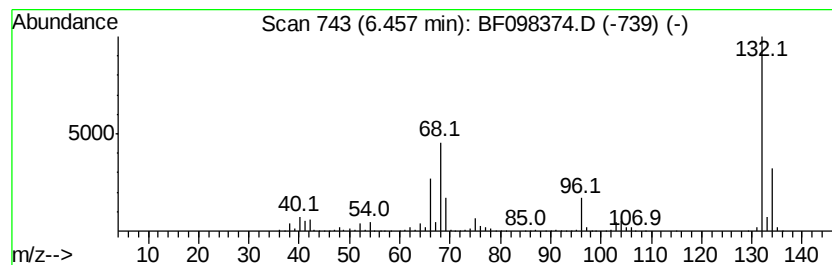
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Peak Number 2 unknown6.46 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.46	74.12 ng	2062210	1,4-Dichlorobenzene-d4	6.68

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
4			1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	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...	4.88	6.0	ng	166533	1	6.68	556473	20.0
unknown6.46	6.46	74.1	ng	2062210	1	6.68	556473	20.0