

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061917\
 Data File : BF096011.D
 Acq On : 19 Jun 2017 17:50
 Operator : SJ/MA
 Sample : PB99828BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB99828BL

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-BF060517.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	1.557	10	12	61	rVB	1903725	3966737	100.00%	18.804%
2	4.510	510	514	541	rBV	49374	180845	4.56%	0.857%
3	5.192	626	630	662	rBV	984624	1938738	48.87%	9.191%
4	6.875	912	916	926	rBV	1095754	1758617	44.33%	8.337%
5	6.963	926	931	951	rVB	1429334	2299061	57.96%	10.899%
6	7.304	985	989	1009	rVB	274956	390884	9.85%	1.853%
7	7.539	1024	1029	1046	rBV	1165874	1601172	40.36%	7.590%
8	8.227	1141	1146	1173	rBV	809830	1216644	30.67%	5.768%
9	9.339	1331	1335	1357	rBV	351029	548485	13.83%	2.600%
10	11.116	1631	1637	1654	rBV	1834670	2439425	61.50%	11.564%
11	12.157	1809	1814	1828	rBV2	419548	587984	14.82%	2.787%
12	13.451	2029	2034	2062	rBV	746731	1154381	29.10%	5.472%
13	14.551	2216	2221	2243	rBV2	324225	515052	12.98%	2.442%
14	16.933	2620	2626	2646	rBB	1441652	1806940	45.55%	8.566%
15	18.280	2851	2855	2872	rBB	247239	337356	8.50%	1.599%
16	19.950	3135	3139	3151	rBV	219217	352309	8.88%	1.670%

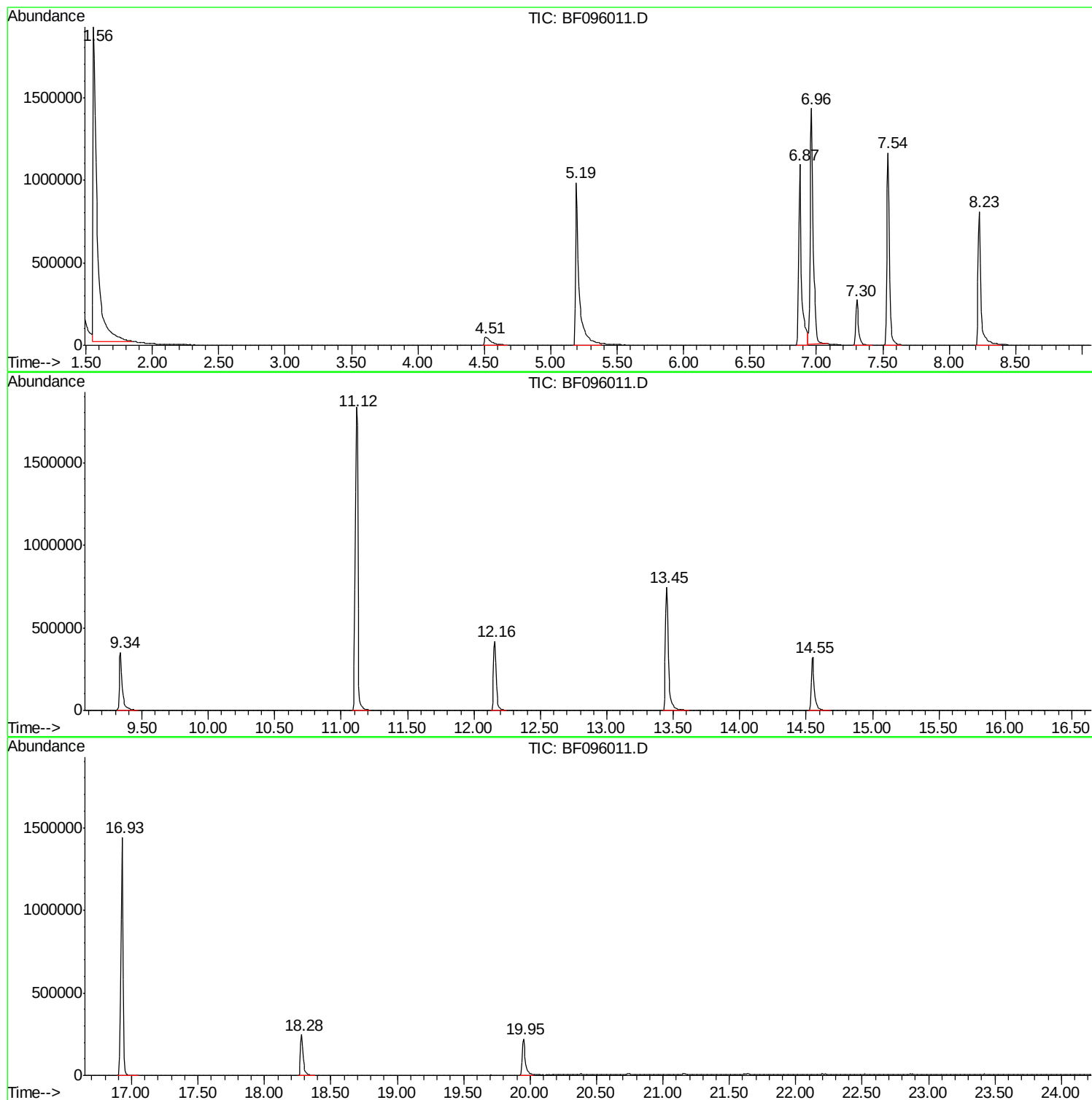
Sum of corrected areas: 21094630

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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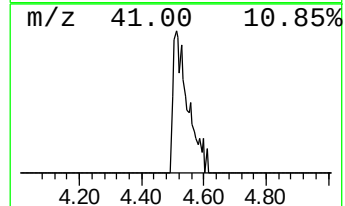
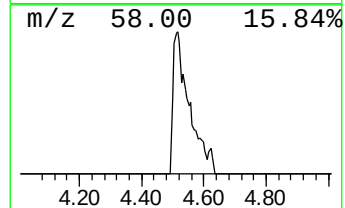
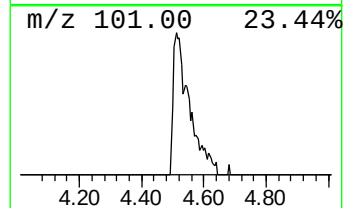
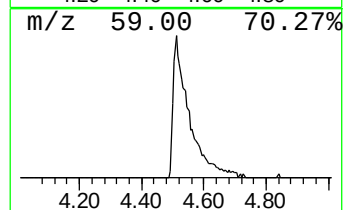
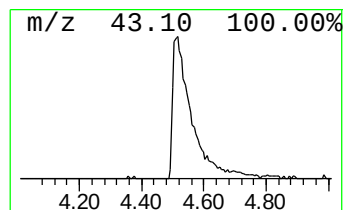
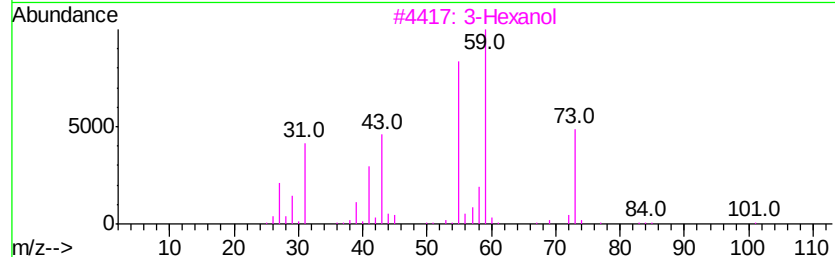
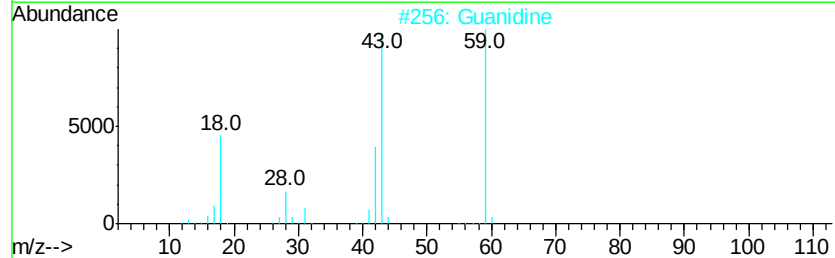
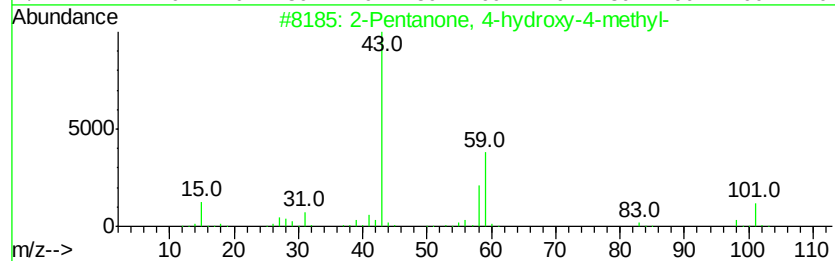
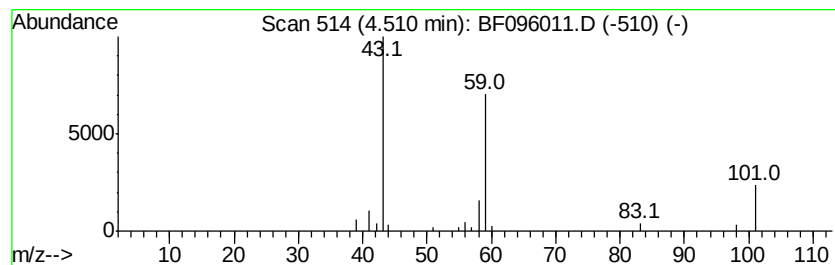
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.51	9.25 ng	180845	1,4-Dichlorobenzene-d4	7.30

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Guanidine	59	CH5N3	000113-00-8	43
3		3-Hexanol	102	C6H14O	000623-37-0	28
4		5-Hexen-2-one	98	C6H10O	000109-49-9	10
5		Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	9



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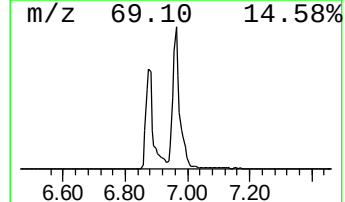
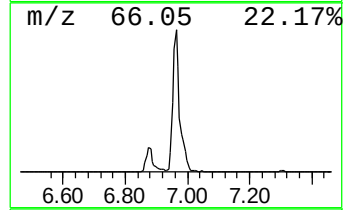
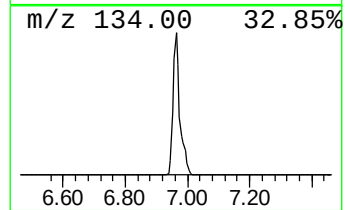
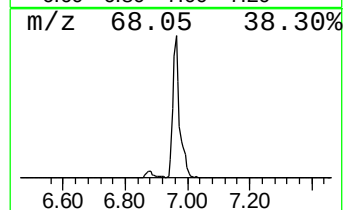
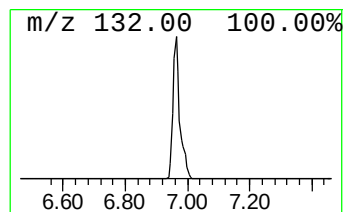
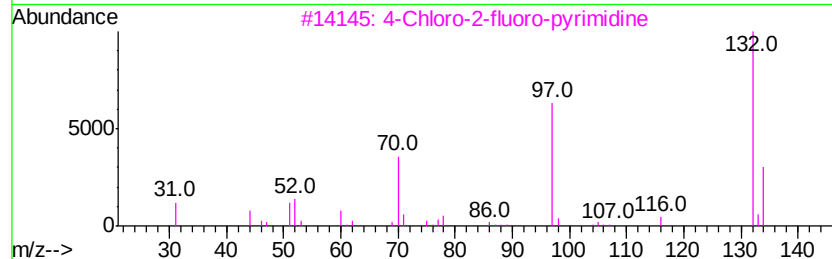
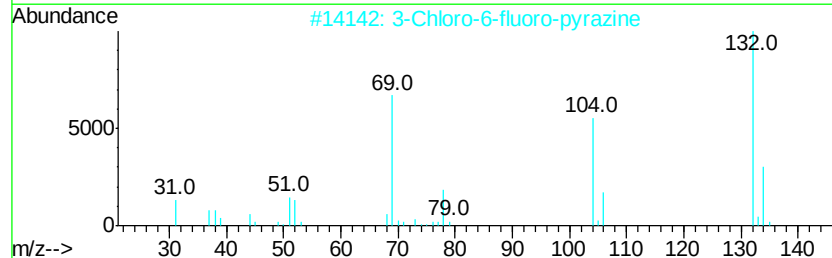
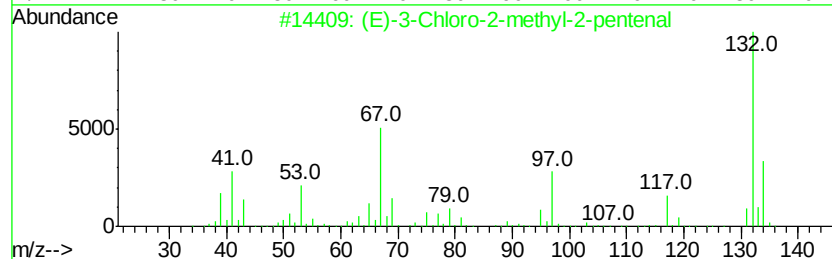
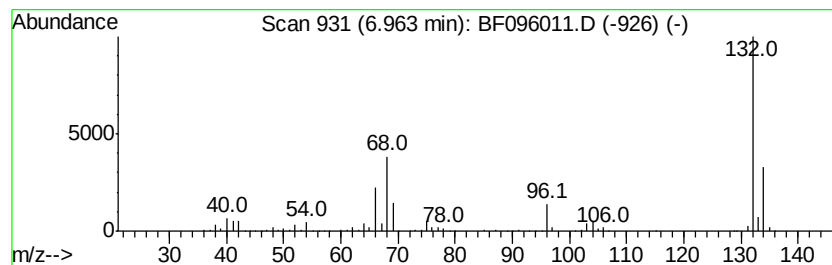
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Peak Number 3 unknown6.96 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.96	117.63 ng	2299060	1,4-Dichlorobenzene-d4	7.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
3		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
4		Xenon	132	Xe	007440-63-3	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...	4.51	9.3	ng	180845	1	7.30	390884	20.0
unknown6.96	6.96	117.6	ng	2299060	1	7.30	390884	20.0