

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060217\
 Data File : BF095656.D
 Acq On : 3 Jun 2017 4:21
 Operator : SJ/MA
 Sample : PB99424BL
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB99424BL

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-BF050217.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	4.787	523	527	545	rBV	92185	220422	8.48%	1.367%
2	5.457	638	641	674	rBV	733431	1522277	58.58%	9.438%
3	7.086	915	918	932	rVV	772045	1395821	53.72%	8.654%
4	7.198	932	937	965	rVB	1091339	1792187	68.97%	11.111%
5	7.533	990	994	1007	rBV	345608	472056	18.17%	2.927%
6	7.775	1030	1035	1050	rBV	1401353	1858746	71.53%	11.524%
7	8.451	1145	1150	1177	rBV	923584	1340961	51.60%	8.313%
8	9.569	1336	1340	1357	rBV	454808	635538	24.46%	3.940%
9	11.351	1636	1643	1662	rBV	1925821	2598533	100.00%	16.110%
10	12.398	1816	1821	1835	rBV2	474010	624627	24.04%	3.872%
11	13.698	2037	2042	2060	rBV	421256	677756	26.08%	4.202%
12	14.798	2225	2229	2246	rBV2	339893	480457	18.49%	2.979%
13	17.150	2623	2629	2637	rBV	1548666	1752056	67.42%	10.862%
14	18.497	2854	2858	2867	rBV	354375	436161	16.78%	2.704%
15	20.174	3139	3143	3154	rBV	236428	322411	12.41%	1.999%

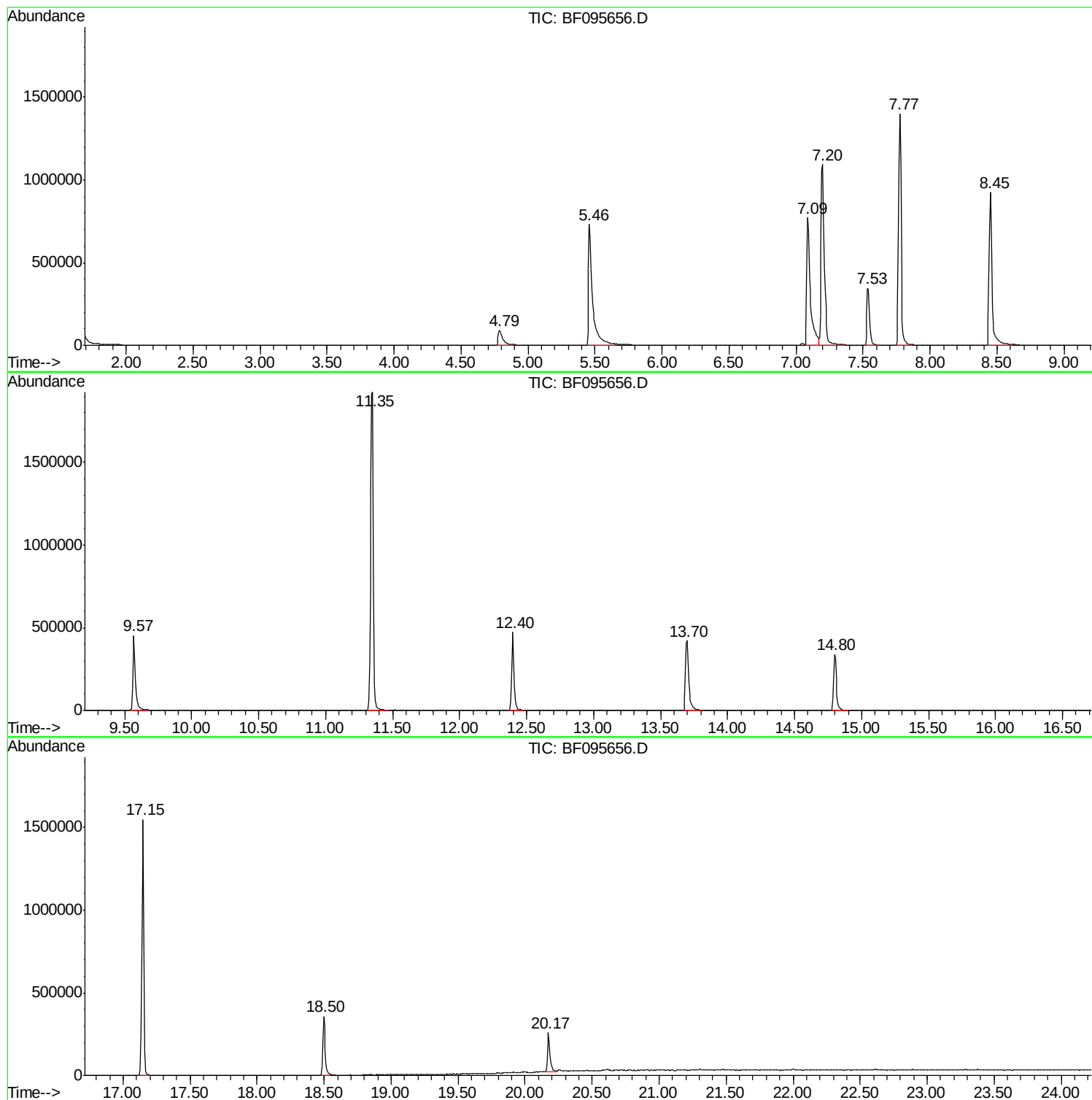
Sum of corrected areas: 16130009

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.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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Instrument :
BNA_F
ClientSampled :
PB99424BL

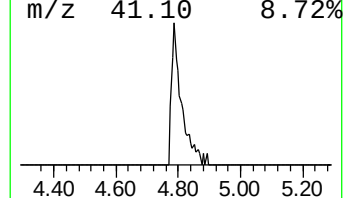
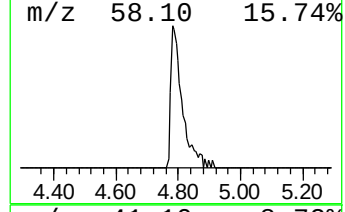
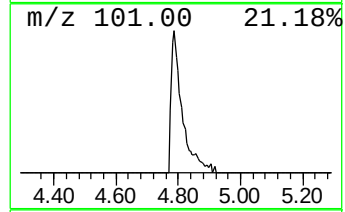
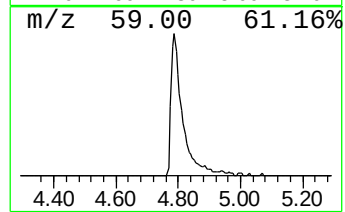
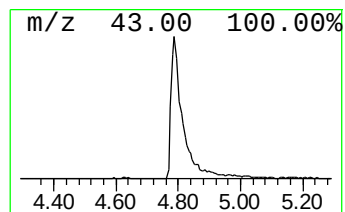
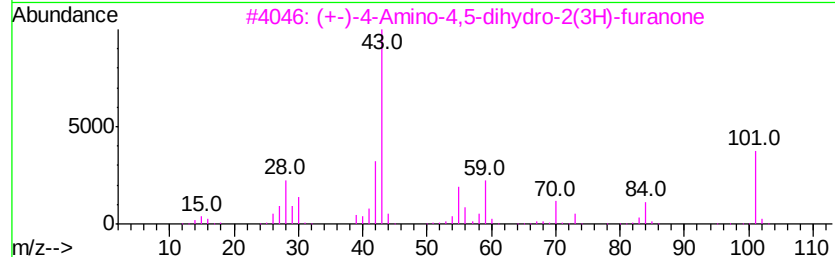
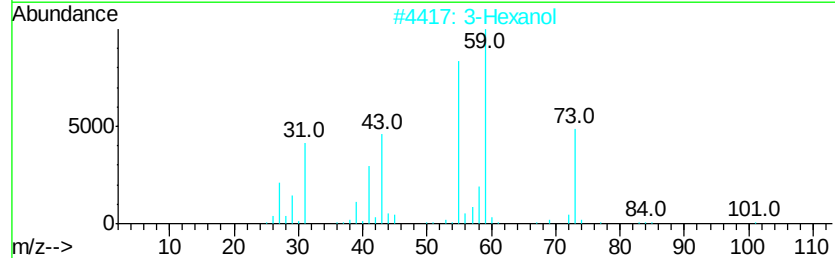
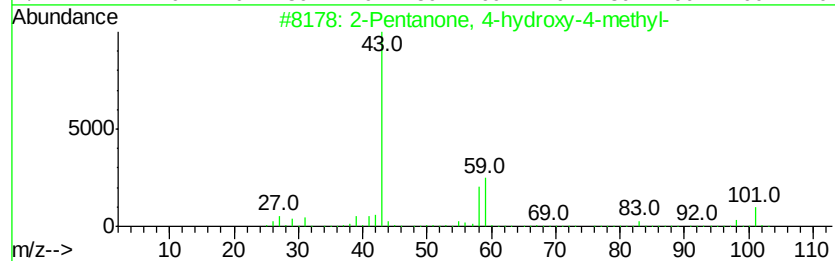
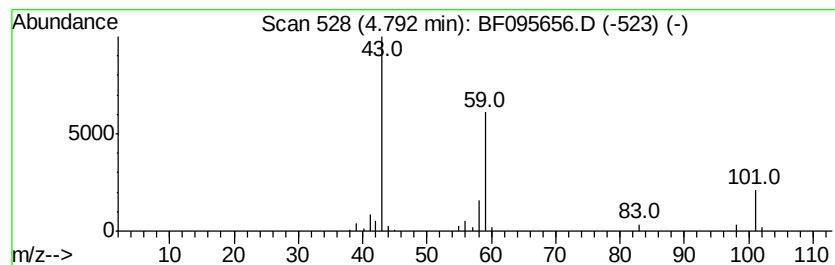
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.79	9.34 ng	220422	1,4-Dichlorobenzene-d4	7.54

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			3-Hexanol	102	C6H14O	000623-37-0	28
3			(+)-4-Amino-4,5-dihydro-2(3H)-furanone	101	C4H7NO2	016504-58-8	9
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			Propanamide, N-ethyl-	101	C5H11NO	005129-72-6	9



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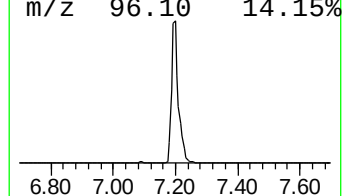
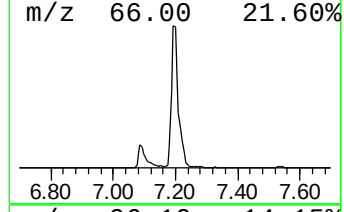
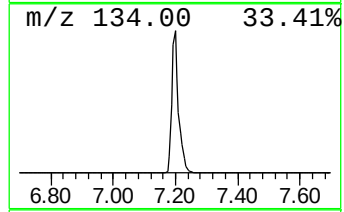
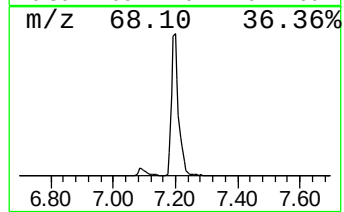
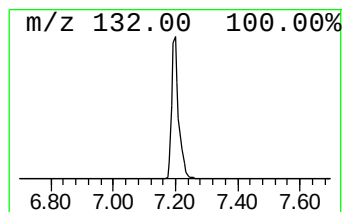
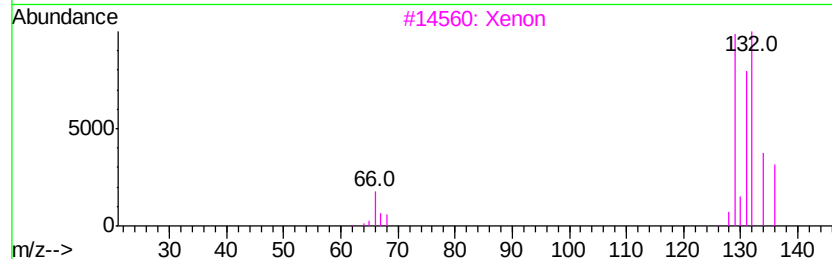
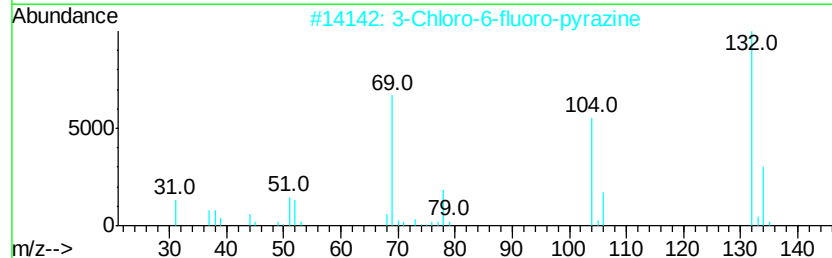
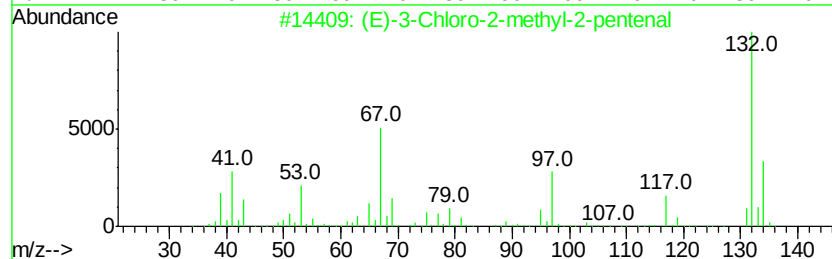
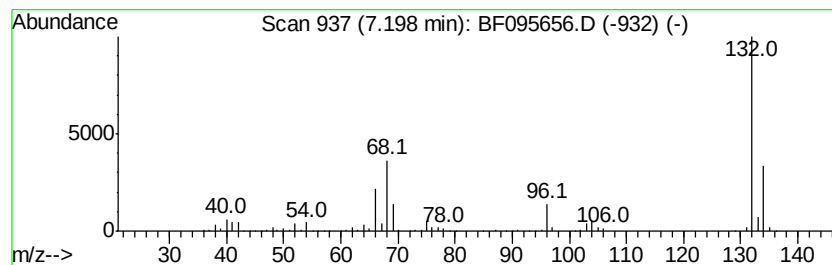
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Peak Number 2 unknown7.20 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.20	75.93 ng	1792190	1,4-Dichlorobenzene-d4	7.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
3		Xenon	132	Xe	007440-63-3	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...	4.79	9.3	ng	220422	1	7.54	472056	20.0
unknown7.20	7.20	75.9	ng	1792190	1	7.54	472056	20.0