

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061317\
 Data File : BF095897.D
 Acq On : 13 Jun 2017 16:55
 Operator : SJ/MA
 Sample : PB99721BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB99721BL

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	4.587	506	510	530	rBV2	51491	168841	7.36%	1.012%
2	5.251	619	623	649	rBV	952401	1761245	76.78%	10.554%
3	6.916	901	906	917	rBV	994454	1640260	71.51%	9.829%
4	7.004	917	921	938	rVB	1288423	2026587	88.35%	12.144%
5	7.351	976	980	995	rVB	262626	374001	16.30%	2.241%
6	7.586	1015	1020	1040	rBV	1048469	1461401	63.71%	8.758%
7	8.269	1131	1136	1162	rBV	718145	1094096	47.70%	6.556%
8	9.380	1321	1325	1345	rBV	350863	523116	22.81%	3.135%
9	11.163	1620	1628	1644	rBV	1570930	2293791	100.00%	13.746%
10	12.198	1799	1804	1820	rVB2	429877	602150	26.25%	3.608%
11	13.498	2020	2025	2049	rBV	829050	1205977	52.58%	7.227%
12	14.598	2207	2212	2226	rBV	389149	601446	26.22%	3.604%
13	16.974	2610	2616	2620	rBV2	1681071	2120157	92.43%	12.705%
14	18.315	2840	2844	2856	rBV	353368	481626	21.00%	2.886%
15	19.986	3125	3128	3142	rBV	204030	332603	14.50%	1.993%

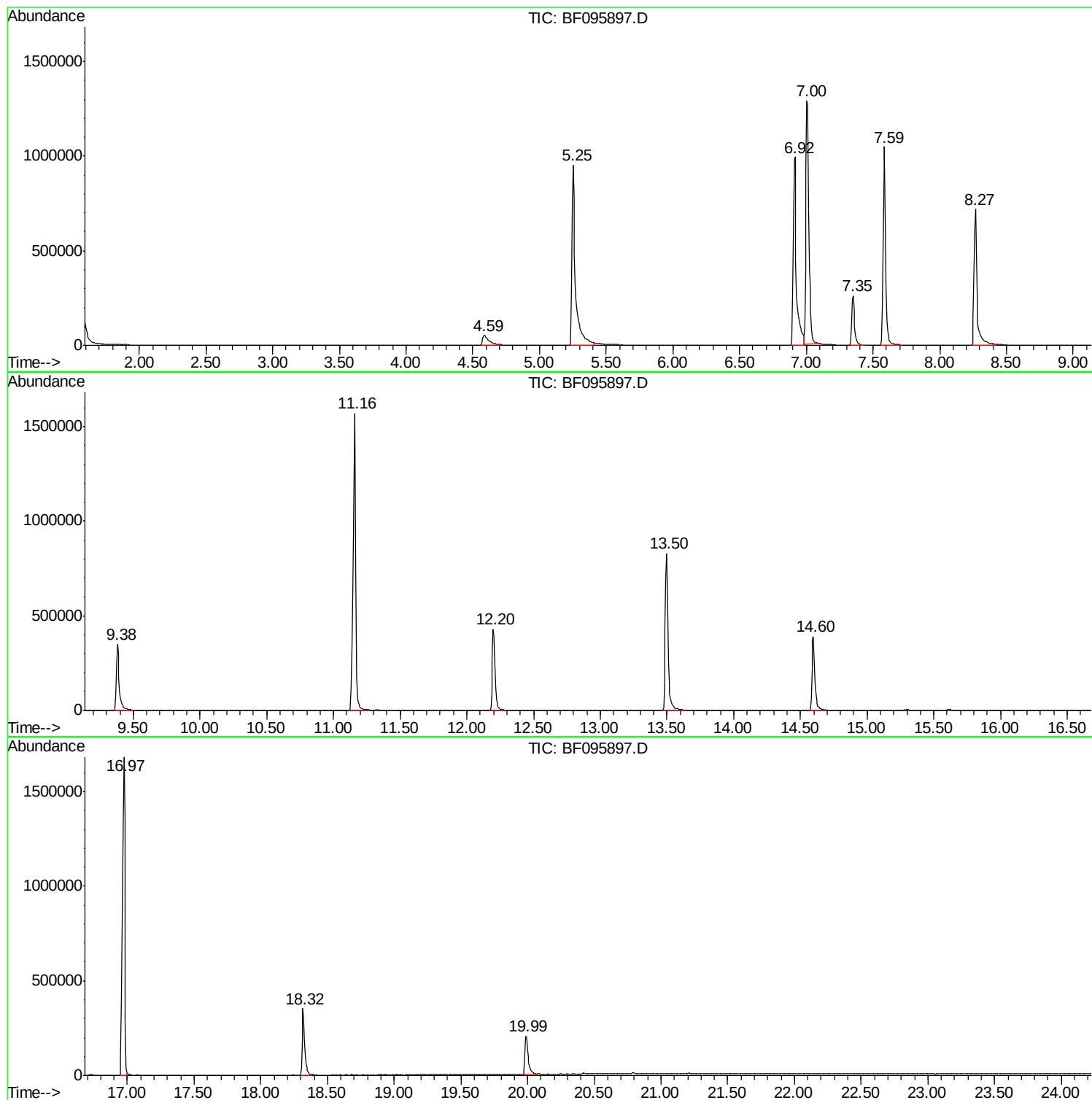
Sum of corrected areas: 16687297

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060517.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 :
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ClientSampled :
PB99721BL

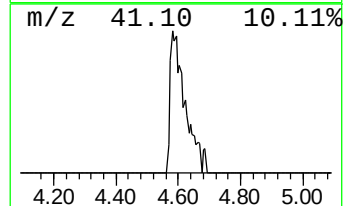
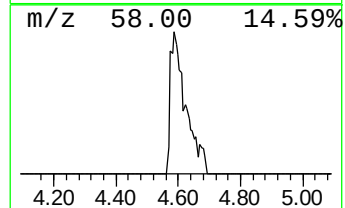
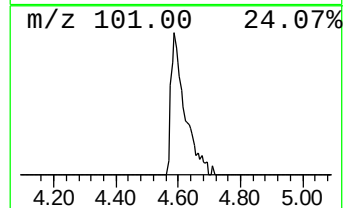
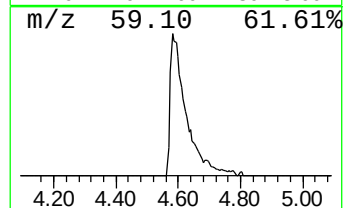
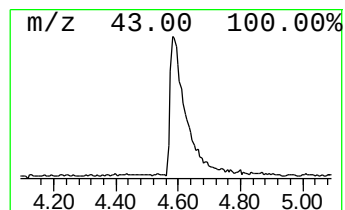
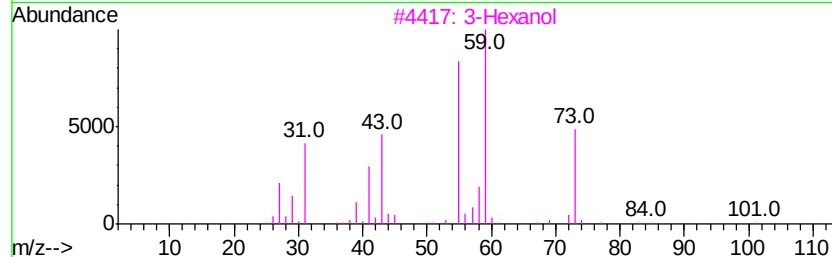
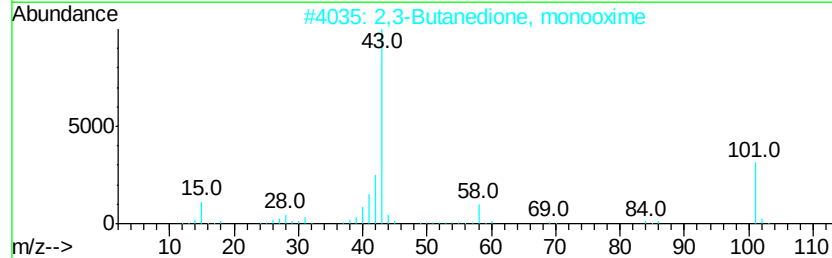
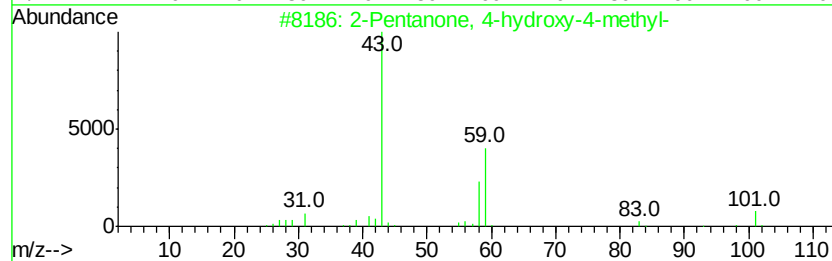
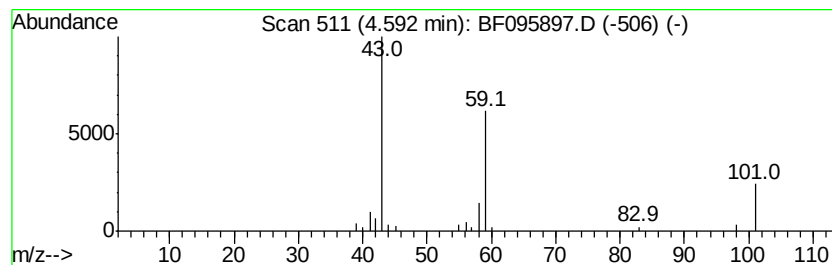
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.59	9.03 ng	168841	1,4-Dichlorobenzene-d4	7.35

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	38
3			3-Hexanol	102	C6H14O	000623-37-0	28
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



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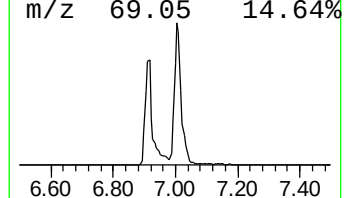
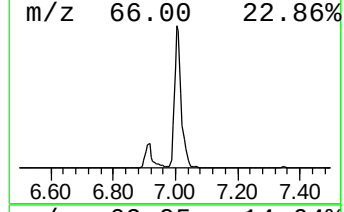
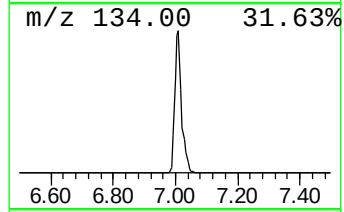
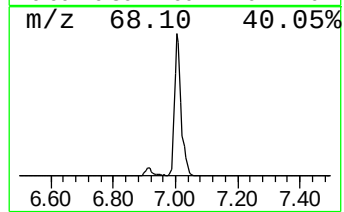
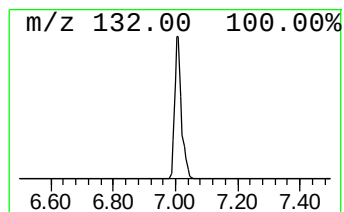
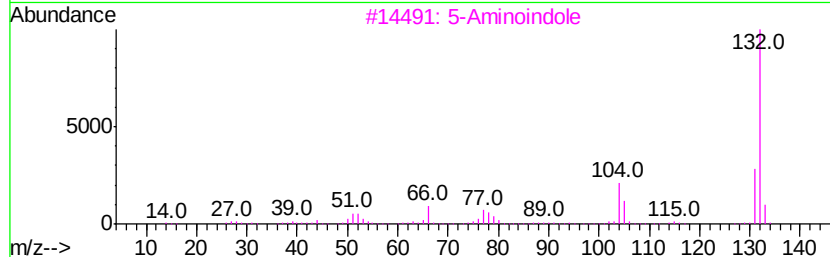
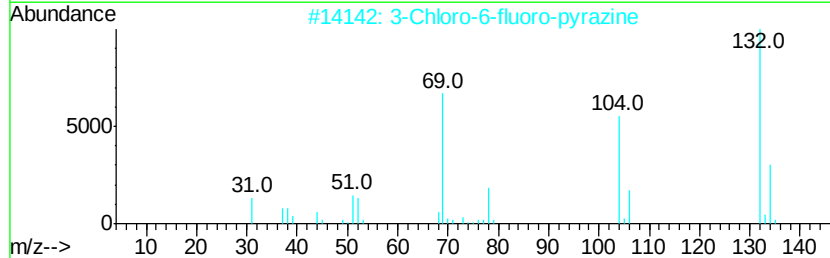
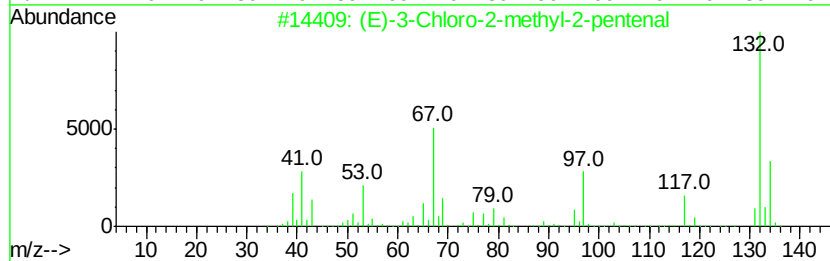
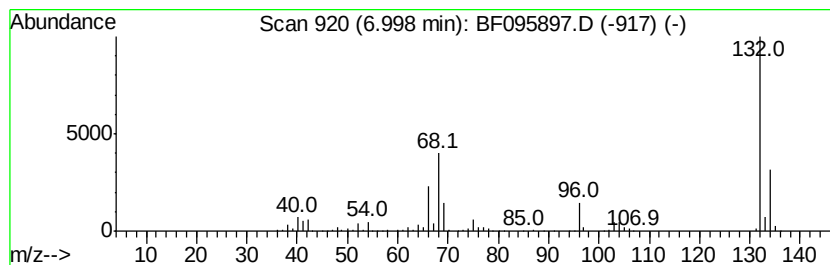
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Peak Number 2 unknown7.00 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.00	108.37 ng	2026590	1,4-Dichlorobenzene-d4	7.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.59	9.0	ng	168841	1	7.35	374001	20.0
unknown7.00	7.00	108.4	ng	2026590	1	7.35	374001	20.0