

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050517\
 Data File : BF094930.D
 Acq On : 5 May 2017 22:19
 Operator : SJ/MA
 Sample : I2949-08
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TWP-07-DUP

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	5.019	520	524	537	rBV	88300	178060	7.00%	1.269%
2	5.678	632	636	656	rBV	269460	571502	22.46%	4.071%
3	7.266	903	906	921	rBV	170765	390995	15.37%	2.785%
4	7.384	921	926	939	rVB	693981	1057916	41.58%	7.537%
5	7.719	979	983	996	rVB	341091	403196	15.85%	2.872%
6	7.960	1019	1024	1043	rBV	1333043	1571670	61.77%	11.197%
7	8.619	1131	1136	1161	rBV	887171	1203992	47.32%	8.577%
8	9.742	1323	1327	1342	rBV	384448	531247	20.88%	3.785%
9	11.513	1623	1628	1647	rBV	1883246	2524759	99.23%	17.987%
10	12.571	1803	1808	1821	rBV2	388105	598059	23.51%	4.261%
11	13.865	2023	2028	2043	rBV	483768	778022	30.58%	5.543%
12	14.977	2212	2217	2231	rBV	376350	611223	24.02%	4.354%
13	17.301	2606	2612	2621	rBV	2072113	2544261	100.00%	18.126%
14	18.642	2835	2840	2853	rBV	443247	621592	24.43%	4.428%
15	20.306	3118	3123	3135	rBV2	274969	450402	17.70%	3.209%

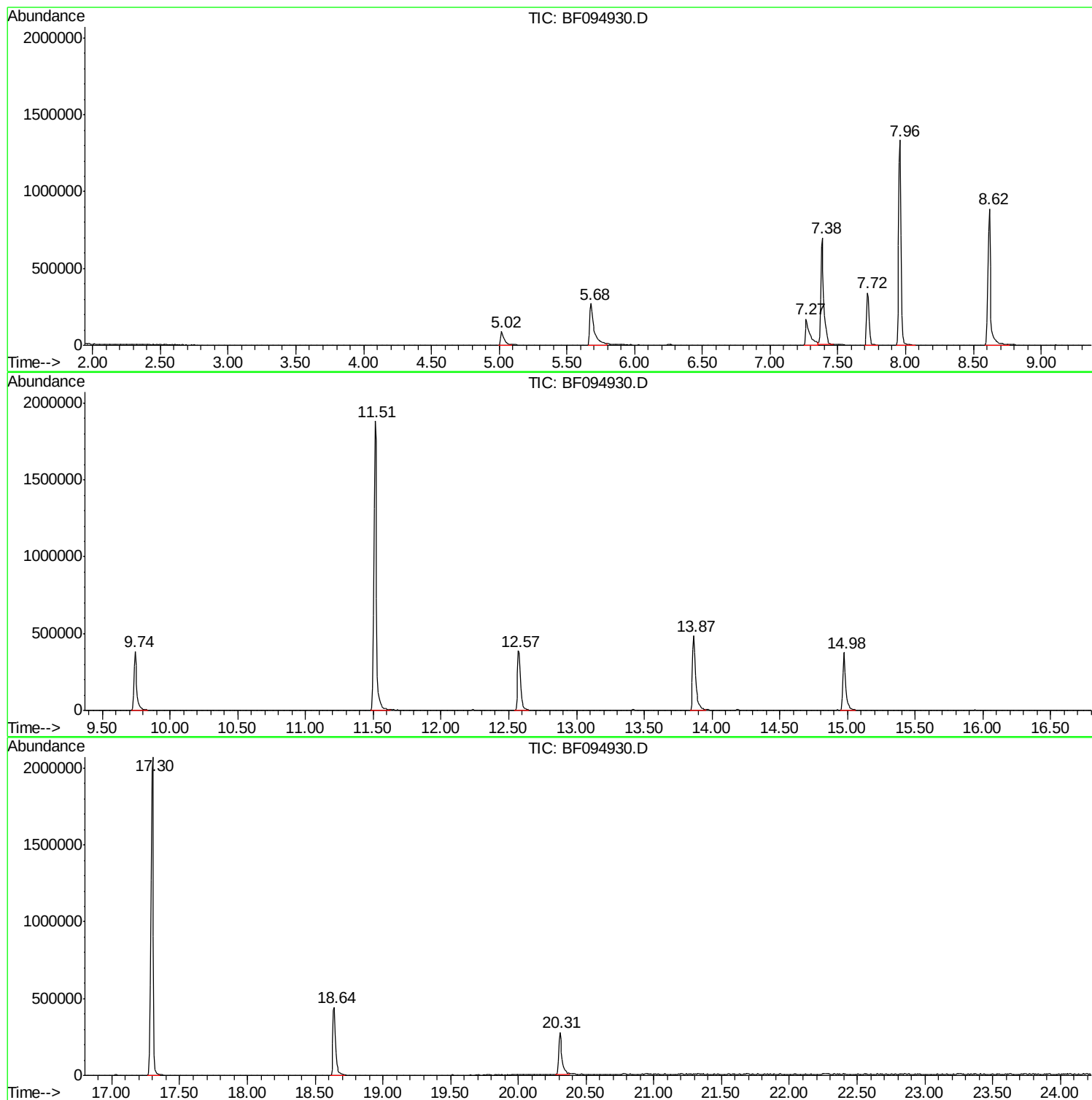
Sum of corrected areas: 14036896

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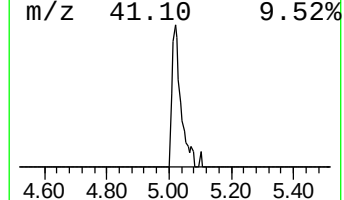
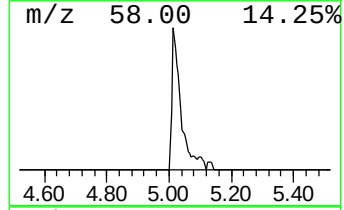
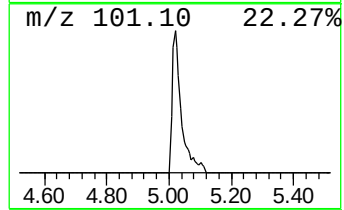
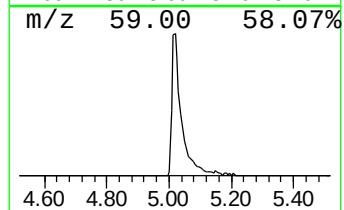
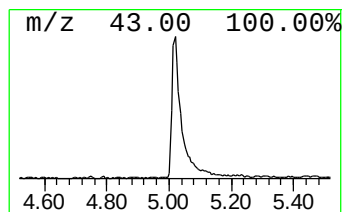
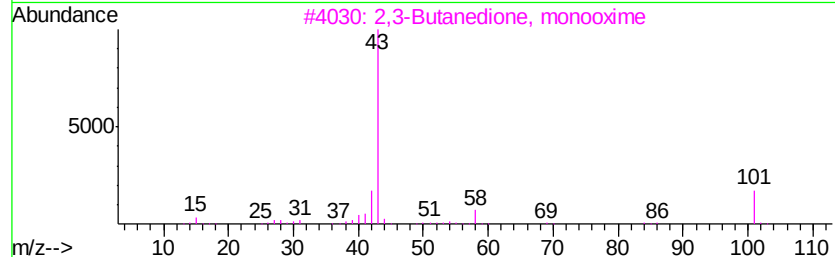
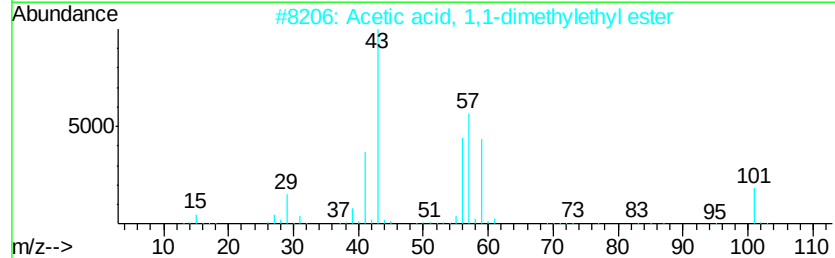
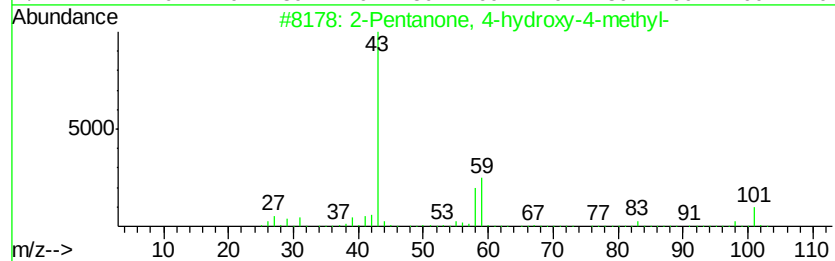
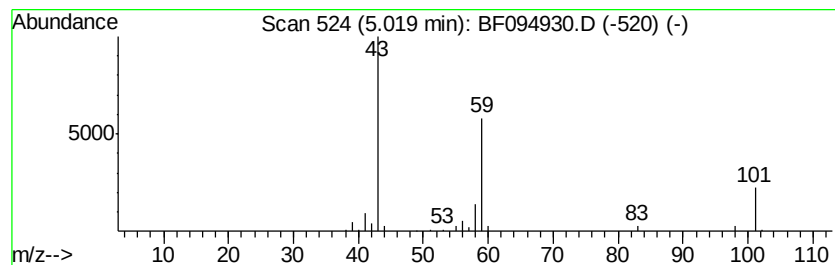
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.02	8.83 ng	178060	1,4-Dichlorobenzene-d4	7.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
5			3-Pentanol, 2-methyl-	102	C6H14O	000565-67-3	9



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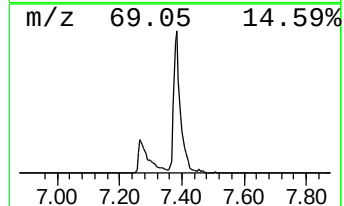
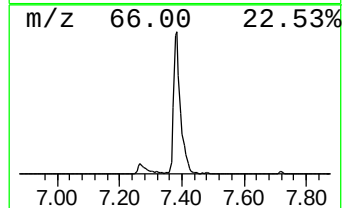
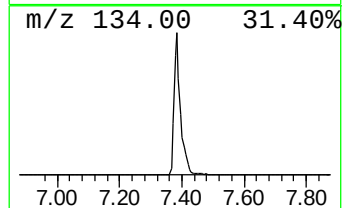
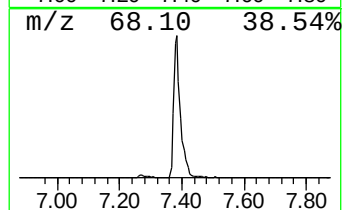
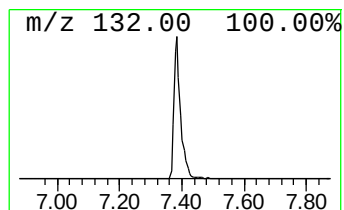
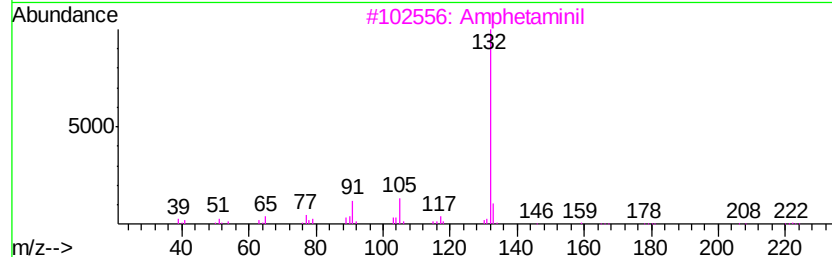
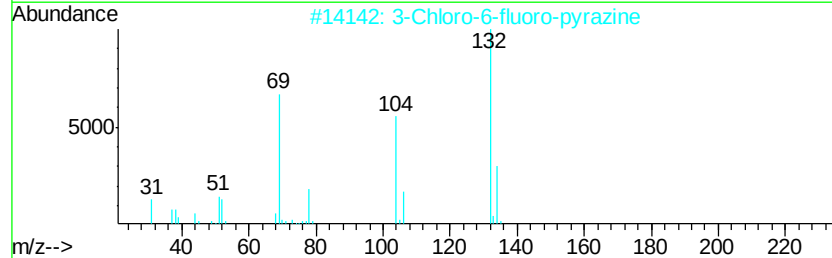
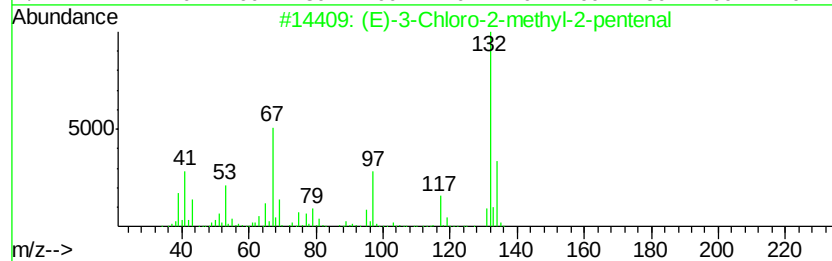
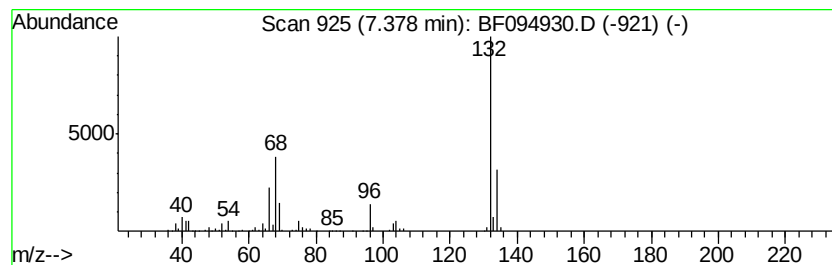
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Peak Number 2 unknown7.38 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.38	52.48 ng	1057920	1,4-Dichlorobenzene-d4	7.72

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		Amphetaminil	250	C17H18N2	017590-01-1	12
4		Tranlylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5		5-Aminoindole	132	C8H8N2	005192-03-0	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.02	8.8	ng	178060	1	7.72	403196	20.0
unknown7.38	7.38	52.5	ng	1057920	1	7.72	403196	20.0