

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092816\
 Data File : BF090278.D
 Acq On : 28 Sep 2016 14:13
 Operator : UM/SJ
 Sample : PB93644BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB93644BL

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-BF092016.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	1.644	4	5	49	rVB	1535986	4050055	100.00%	13.314%
2	4.558	257	260	272	rBV	335233	589124	14.55%	1.937%
3	5.039	299	302	304	rBV	2404356	2700250	66.67%	8.877%
4	6.136	395	398	400	rBV	2483335	2752127	67.95%	9.047%
5	6.239	405	407	410	rVV	2430737	2700143	66.67%	8.876%
6	6.479	425	428	430	rBV	435946	626488	15.47%	2.060%
7	6.627	439	441	444	rBV	2130240	2194803	54.19%	7.215%
8	7.050	475	478	480	rBV	1797914	1907618	47.10%	6.271%
9	7.759	538	540	543	rBV	810696	832849	20.56%	2.738%
10	8.845	632	635	637	rBV	3441276	3754258	92.70%	12.342%
11	9.496	690	692	695	rBV	973943	968913	23.92%	3.185%
12	10.285	758	761	764	rBV	1862580	1902611	46.98%	6.255%
13	10.971	818	821	823	rBV	856992	939617	23.20%	3.089%
14	12.559	957	960	962	rBV	2181038	3167822	78.22%	10.414%
15	13.576	1046	1049	1055	rBV	728761	704629	17.40%	2.316%
16	14.902	1162	1165	1167	rBV	419776	627927	15.50%	2.064%

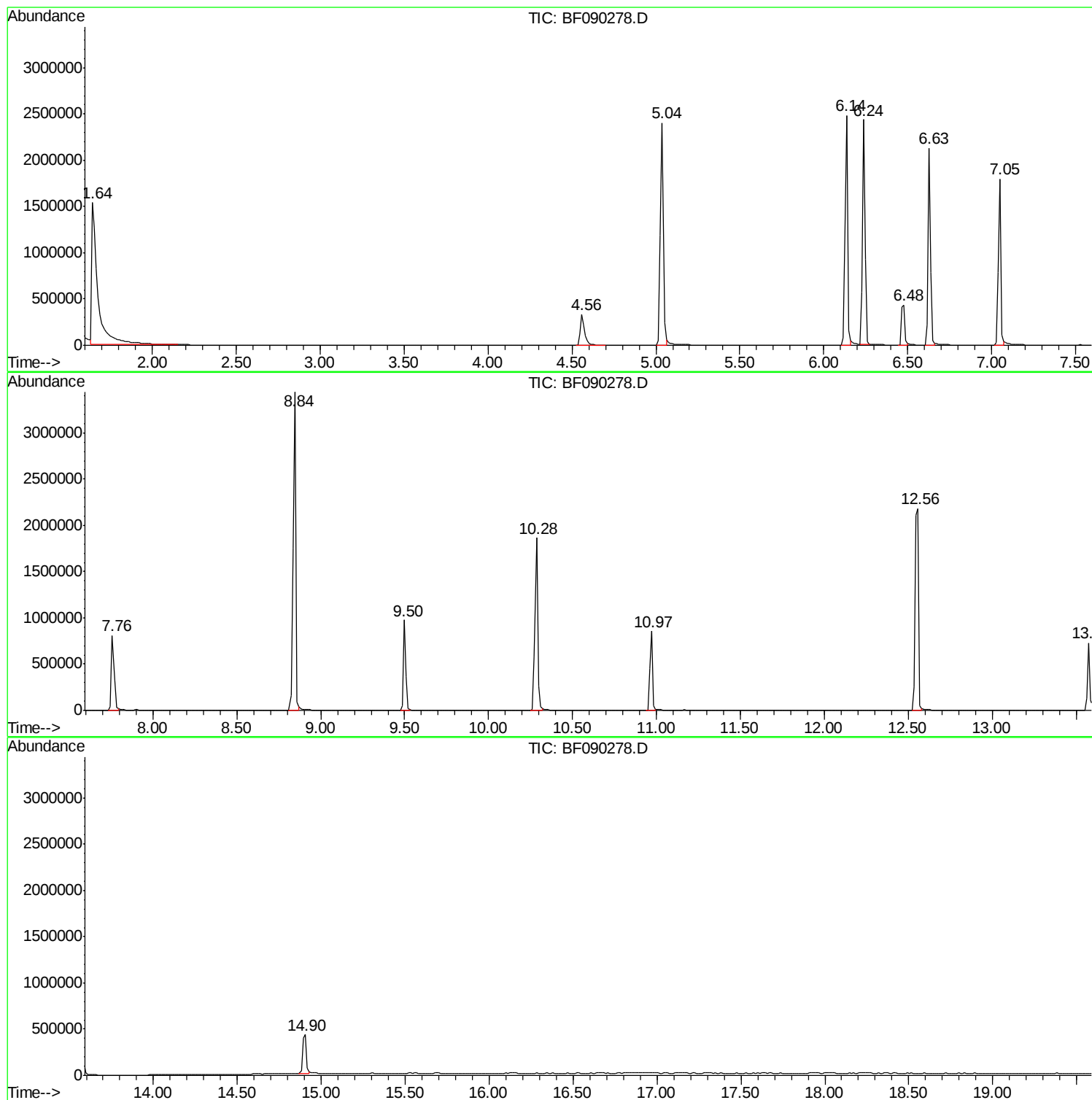
Sum of corrected areas: 30419234

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092016.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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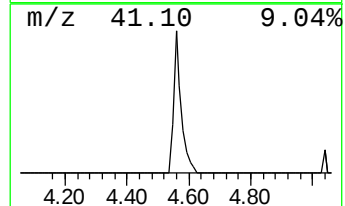
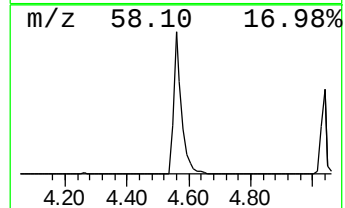
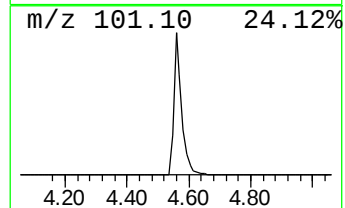
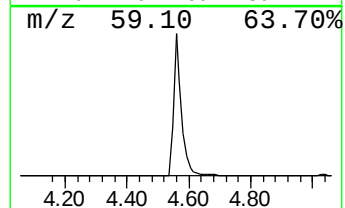
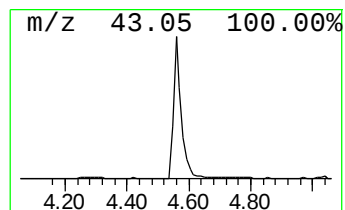
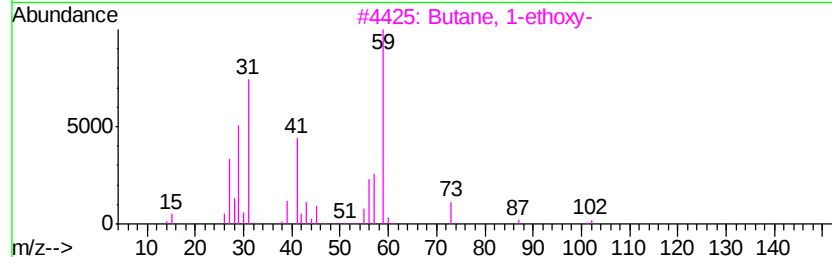
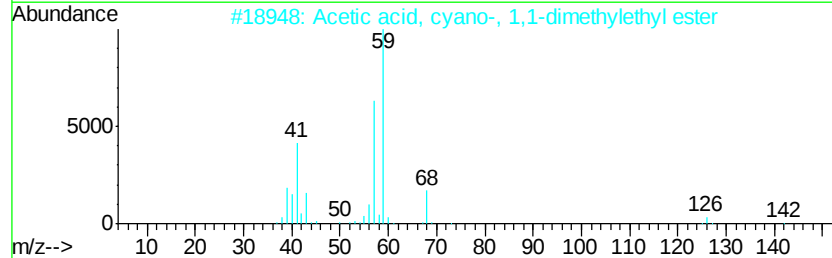
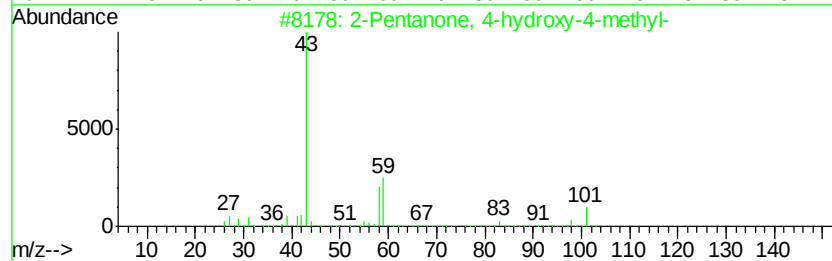
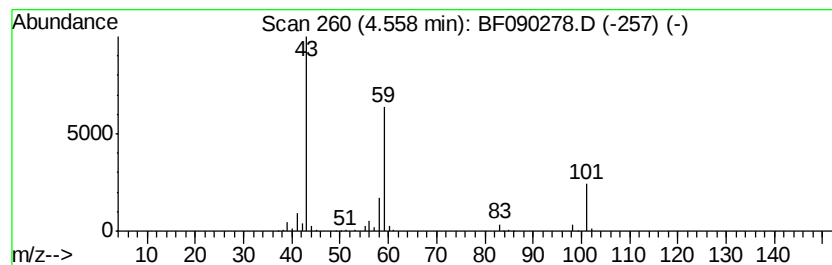
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.56	18.81 ng	589124	1,4-Dichlorobenzene-d4	6.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		5-Hexen-2-one	98	C6H10O	000109-49-9	9



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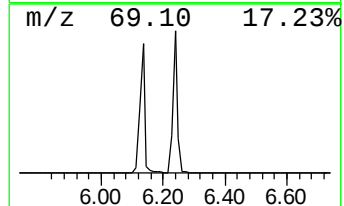
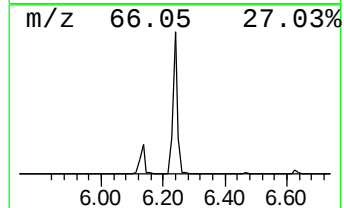
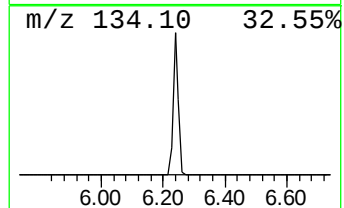
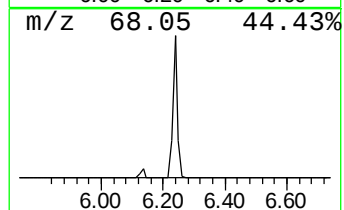
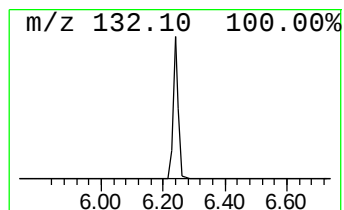
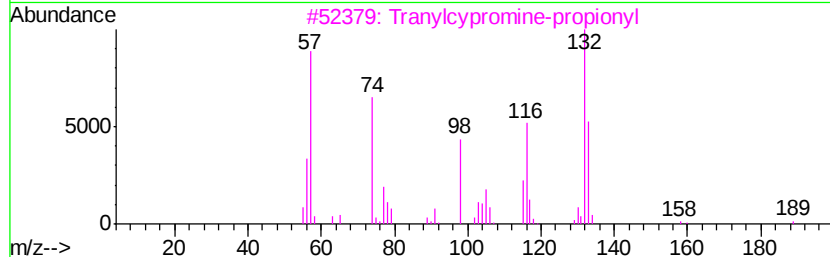
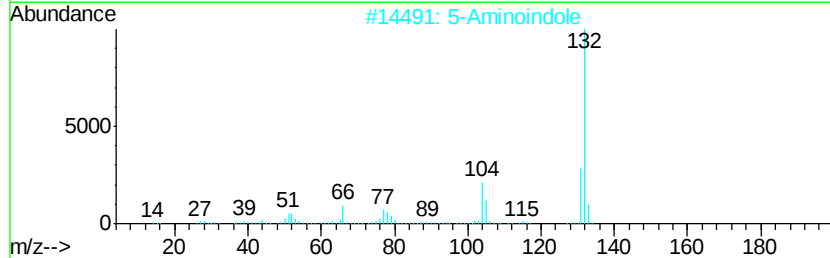
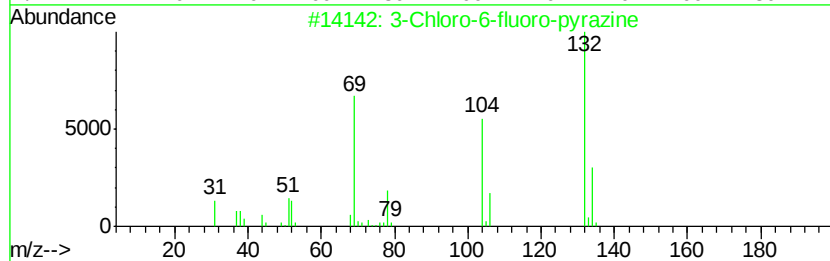
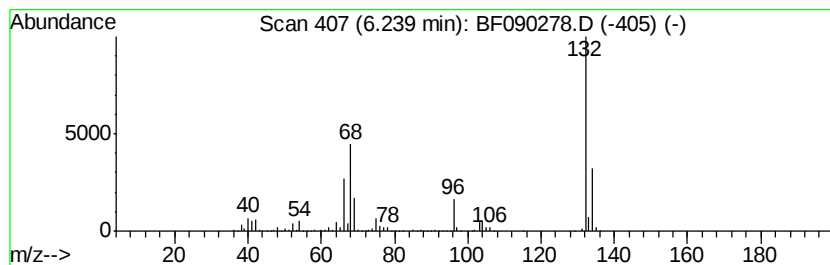
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Peak Number 3 unknown6.24 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.24	86.20 ng	2700140	1,4-Dichlorobenzene-d4	6.48

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10
4			2H-Cyclopentadlpvridazine, 2-me...	132	C8H8N2	022291-85-6	9
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...	4.56	18.8	ng	589124	1	6.48	626488	20.0
unknown6.24	6.24	86.2	ng	2700140	1	6.48	626488	20.0