

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062615\
 Data File : BF080192.D
 Acq On : 26 Jun 2015 16:33
 Operator : TP/IZ
 Sample : PB84223BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB84223BL

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-BF061715.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	5.510	209	212	218	rBV	89862	126518	10.79%	1.536%
2	5.899	244	246	249	rBV	805309	767434	65.46%	9.318%
3	6.893	331	333	336	rBV	870629	778330	66.39%	9.451%
4	7.064	345	348	350	rBV	663269	820001	69.95%	9.956%
5	7.293	366	368	371	rBV	244936	197916	16.88%	2.403%
6	7.453	380	382	384	rBV	857163	678007	57.84%	8.232%
7	7.853	415	417	419	rBV	539606	492553	42.02%	5.981%
8	8.585	479	481	484	rBV	330061	266087	22.70%	3.231%
9	9.659	573	575	578	rBV	1265927	1025108	87.45%	12.447%
10	10.356	634	636	639	rVB	365539	306023	26.10%	3.716%
11	11.145	703	705	708	rBV2	554943	575474	49.09%	6.987%
12	11.854	765	767	770	rBV	239468	311979	26.61%	3.788%
13	13.511	910	912	915	rBV	1241427	1172284	100.00%	14.234%
14	14.745	1018	1020	1023	rBV	359519	349288	29.80%	4.241%
15	16.563	1176	1179	1187	rVB	203696	334847	28.56%	4.066%
16	18.026	1302	1307	1313	rBV5	7674	33998	2.90%	0.413%

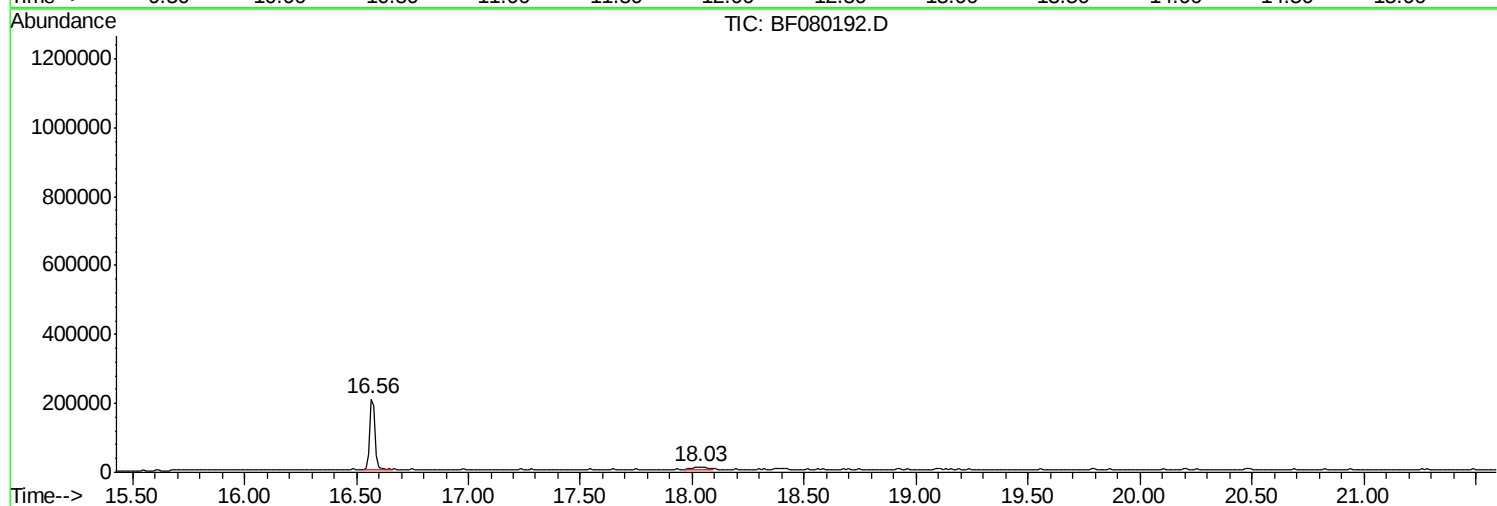
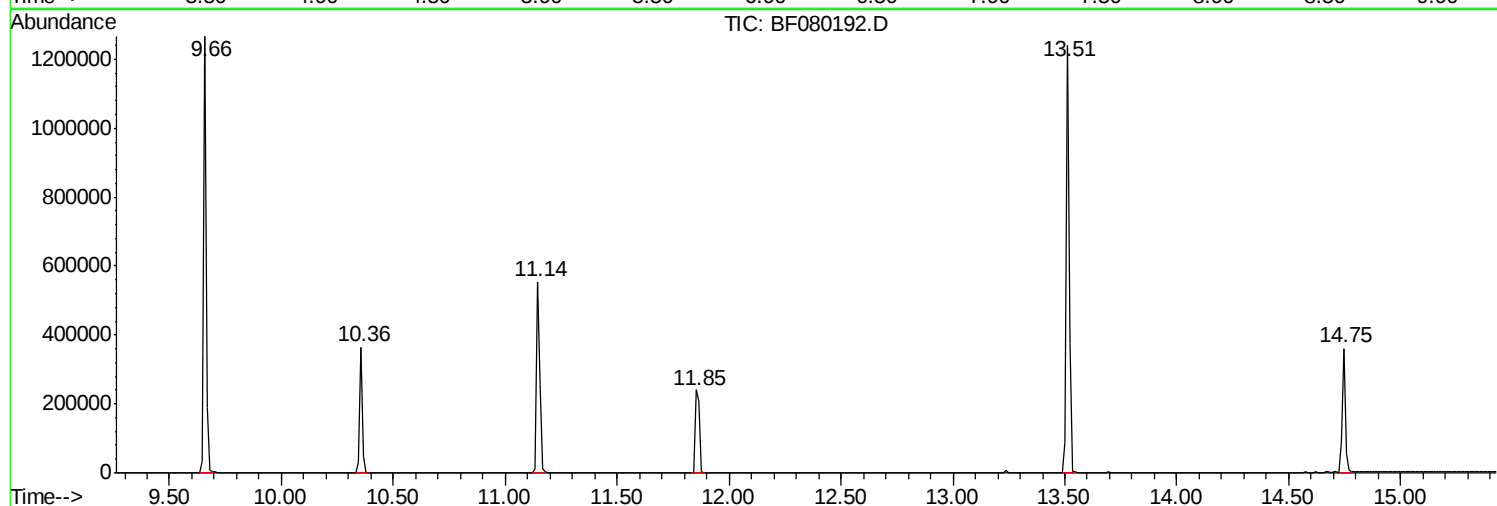
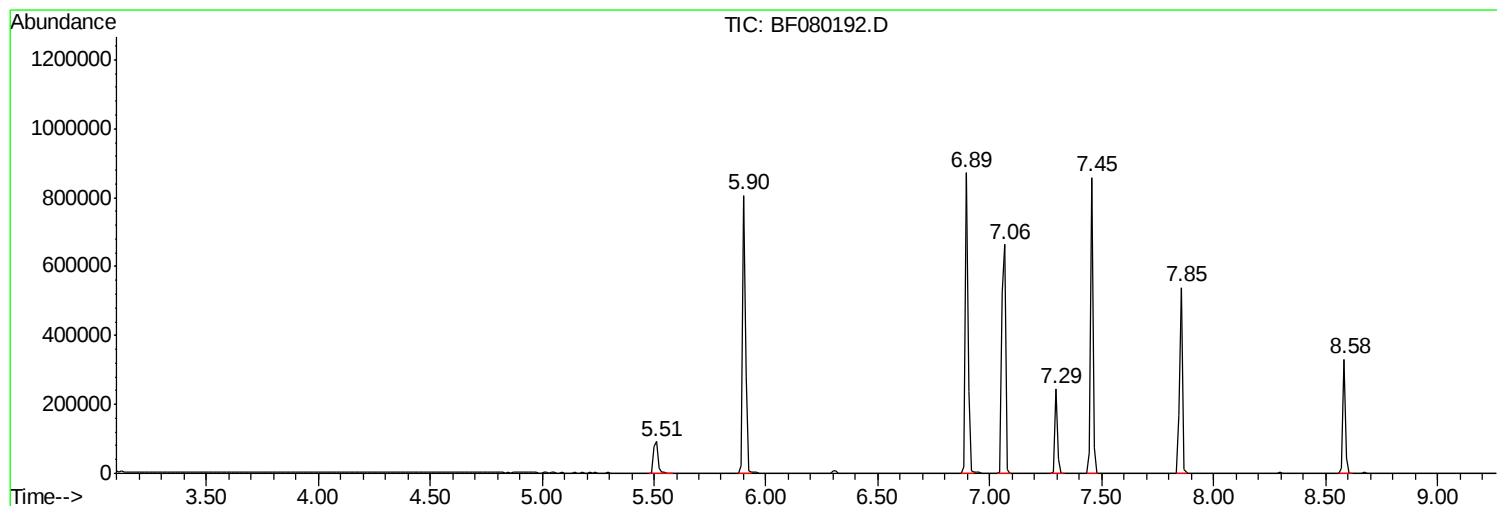
Sum of corrected areas: 8235847

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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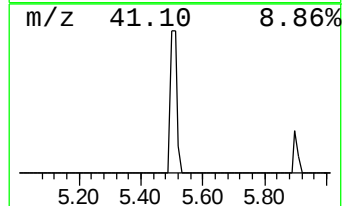
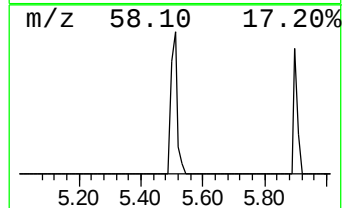
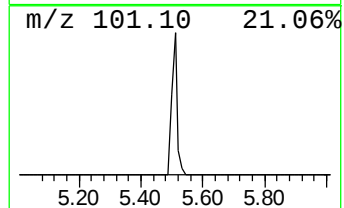
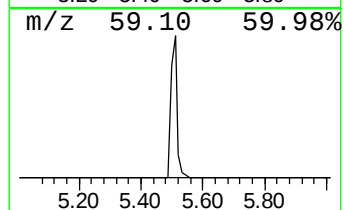
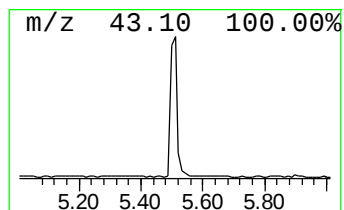
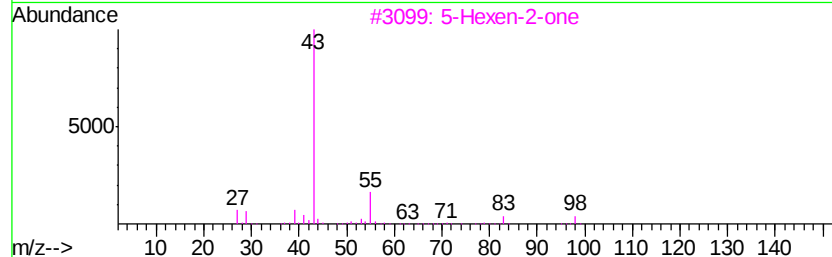
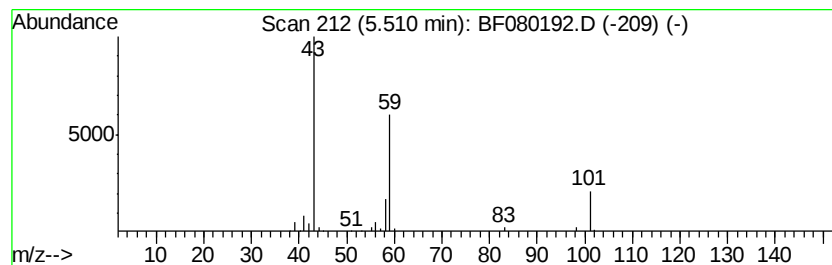
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.51	12.79 ng	126518	1,4-Dichlorobenzene-d4	7.29

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53	
2	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17	
3	5-Hexen-2-one	98	C6H10O	000109-49-9	9	
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9	
5	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9	



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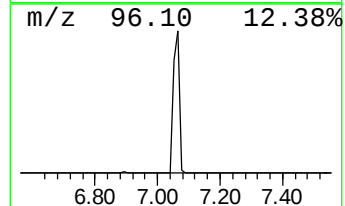
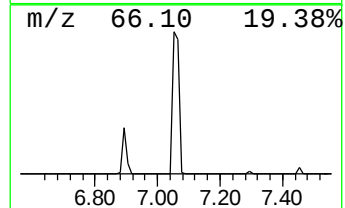
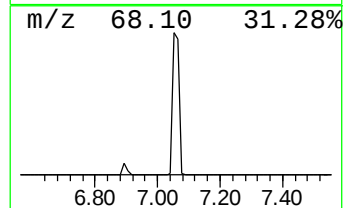
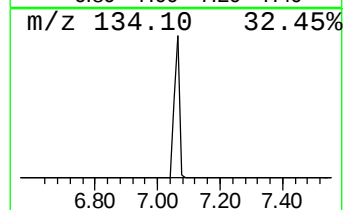
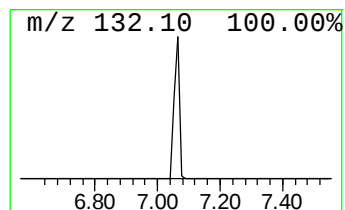
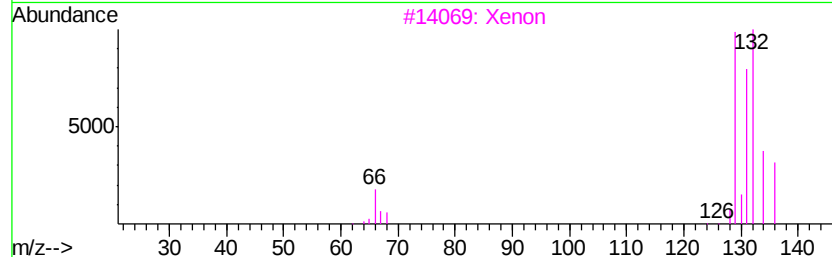
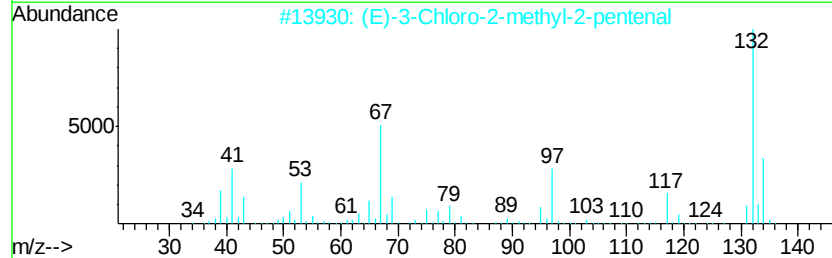
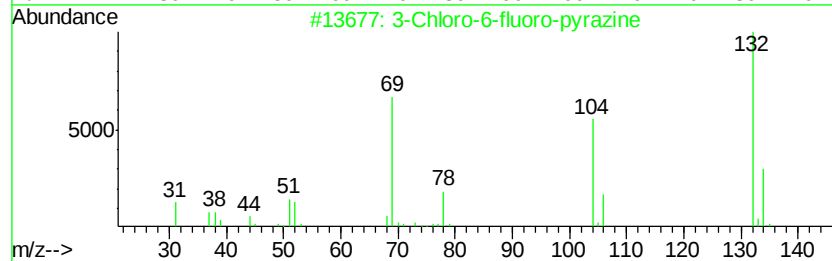
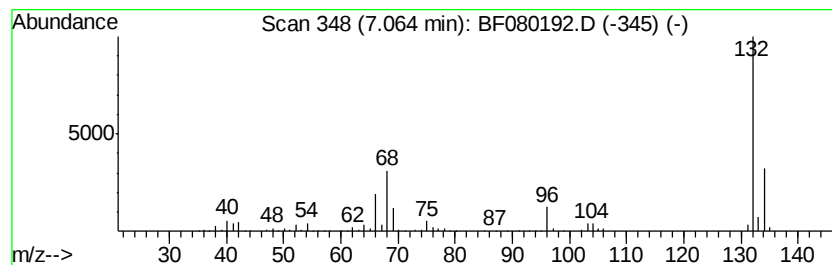
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Peak Number 2 unknown7.06 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.06	82.86 ng	820001	1,4-Dichlorobenzene-d4	7.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Xenon	132	Xe	007440-63-3	9
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	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...	5.51	12.8	ng	126518	1	7.29	197916	20.0
unknown7.06	7.06	82.9	ng	820001	1	7.29	197916	20.0