

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072117\
 Data File : BF096964.D
 Acq On : 22 Jul 2017 1:34
 Operator : SJ/JU
 Sample : PB100832BL
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB100832BL

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-BF071317.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	4.722	460	465	480	rBV	353831	498245	18.89%	2.552%
2	5.163	536	540	553	rBV	2215019	2323767	88.08%	11.902%
3	6.210	713	718	721	rBV	2078965	1911418	72.45%	9.790%
4	6.328	734	738	741	rBV	2188622	1925478	72.99%	9.862%
5	6.557	773	777	784	rVB	691729	637899	24.18%	3.267%
6	6.710	799	803	807	rBV	2195956	2105411	79.81%	10.784%
7	7.122	868	873	876	rBV	1861474	1738398	65.89%	8.904%
8	7.833	990	994	998	rBV	870432	830550	31.48%	4.254%
9	8.916	1173	1178	1181	rBV	2868160	2638152	100.00%	13.512%
10	9.580	1287	1291	1295	rBV	834875	754725	28.61%	3.866%
11	10.369	1421	1425	1434	rBV	885100	802974	30.44%	4.113%
12	11.057	1538	1542	1546	rBV	660590	560759	21.26%	2.872%
13	12.645	1807	1812	1815	rBV	1704724	1674013	63.45%	8.574%
14	13.680	1984	1988	1992	rBV	610823	558178	21.16%	2.859%
15	15.045	2215	2220	2225	rBV	483439	531825	20.16%	2.724%
16	16.386	2443	2448	2454	rBV8	13227	32482	1.23%	0.166%

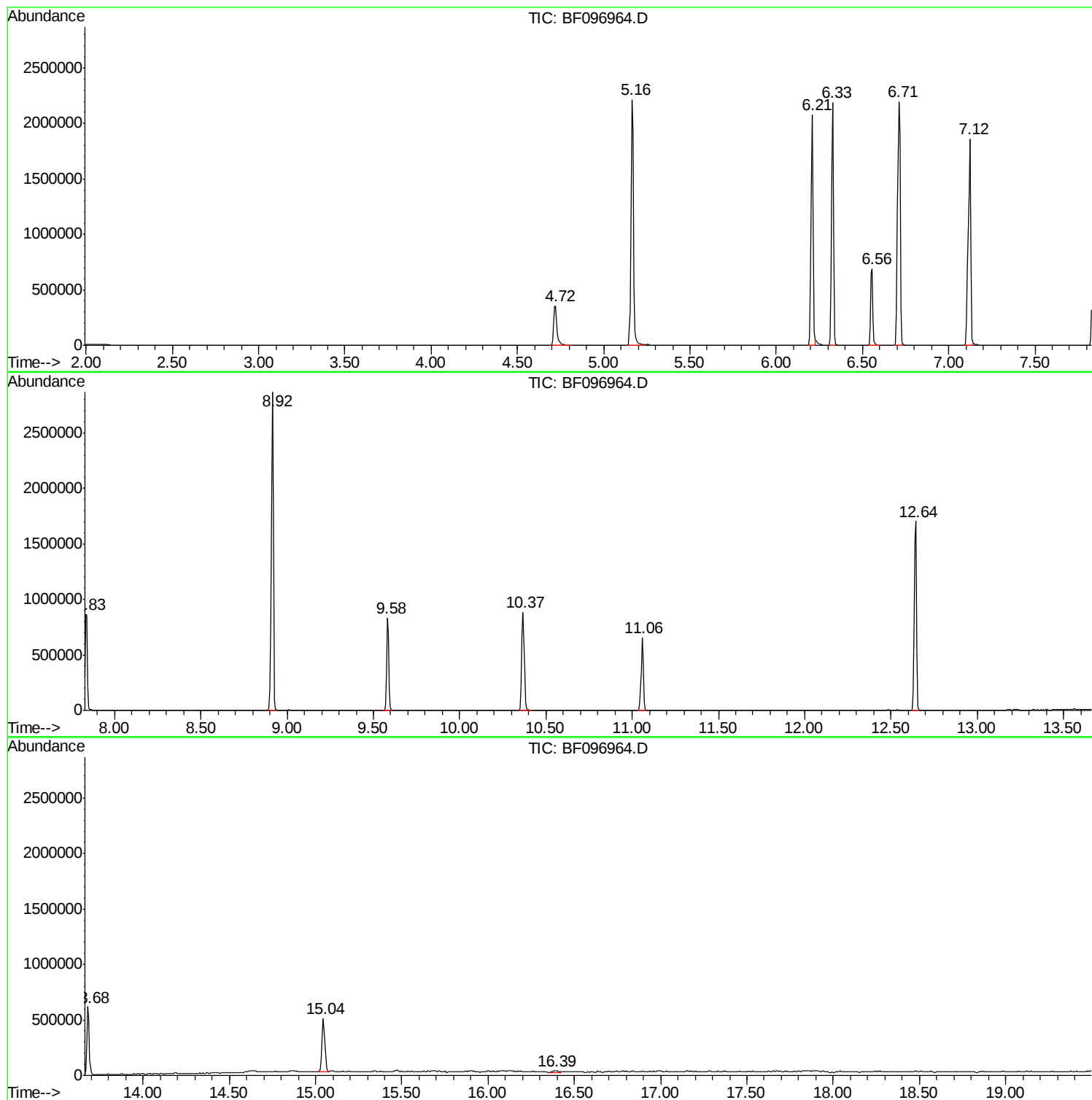
Sum of corrected areas: 19524274

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071317.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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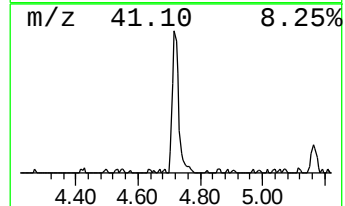
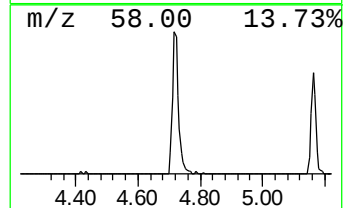
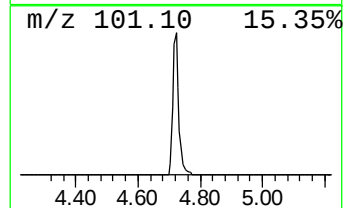
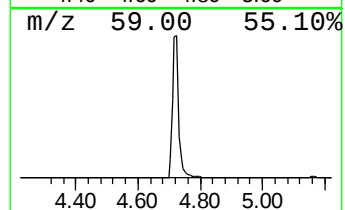
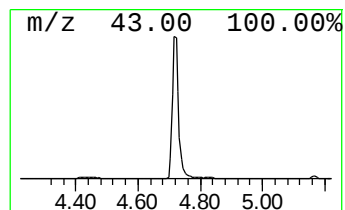
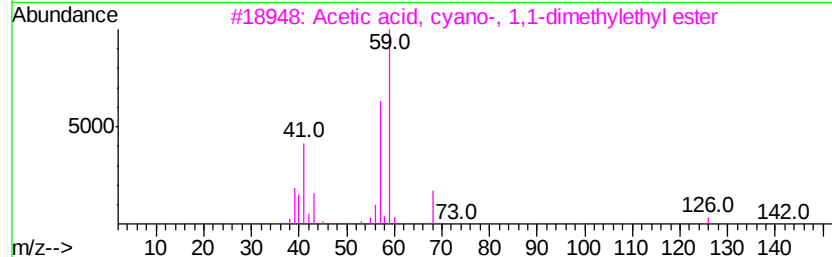
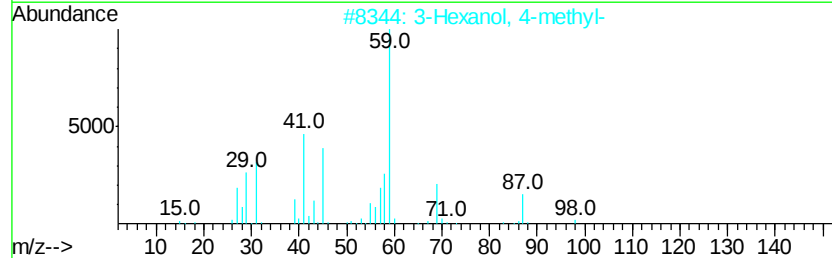
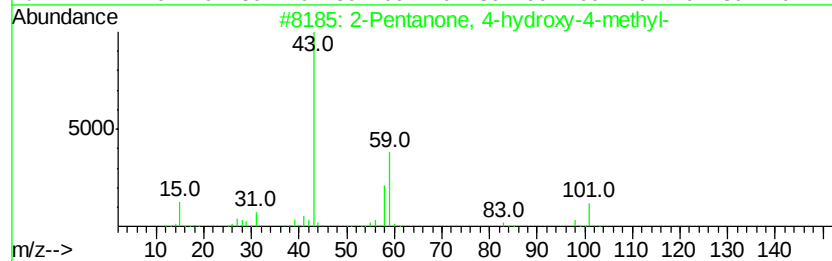
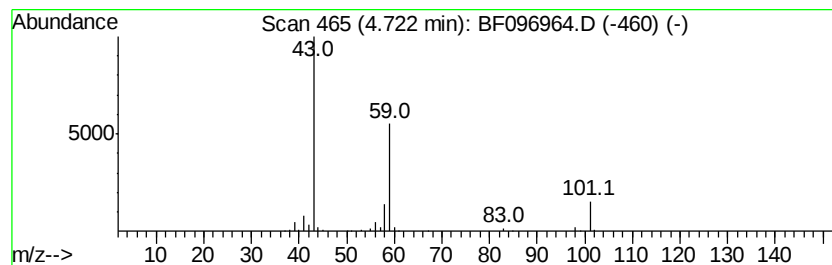
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.72	15.62 ng	498245	1,4-Dichlorobenzene-d4	6.56

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, 4-methyl-	116	C7H16O	000615-29-2	33
3		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	9



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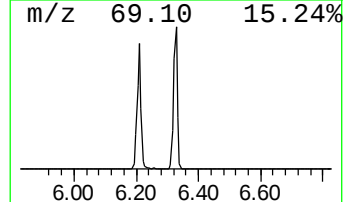
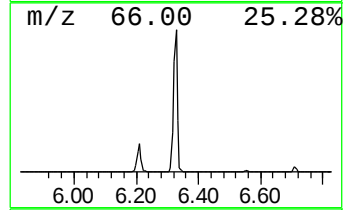
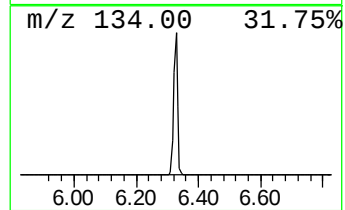
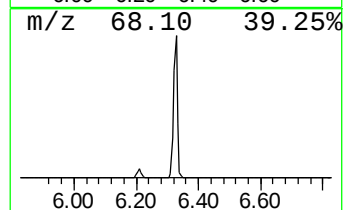
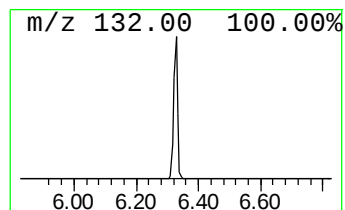
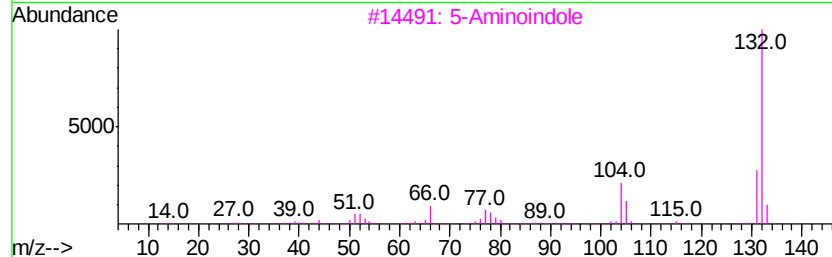
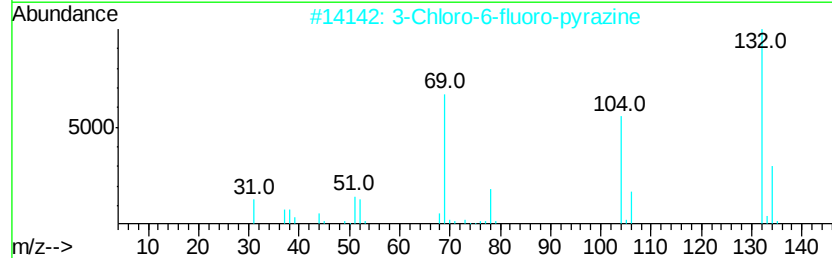
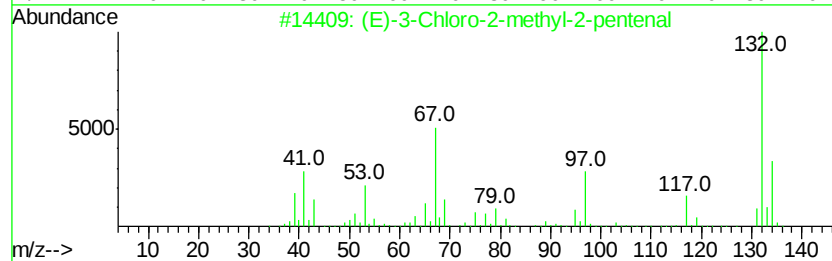
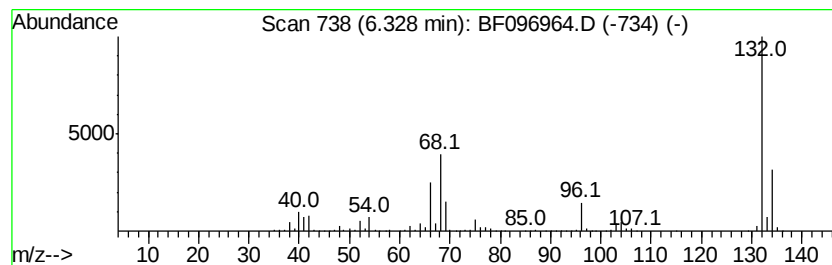
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Peak Number 2 unknown6.33 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.33	60.37 ng	1925480	1,4-Dichlorobenzene-d4	6.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.72	15.6	ng	498245	1	6.56	637899	20.0
unknown6.33	6.33	60.4	ng	1925480	1	6.56	637899	20.0