

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072417\
 Data File : BF096995.D
 Acq On : 24 Jul 2017 19:57
 Operator : SJ/JU
 Sample : PB100872BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB100872BL

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	4.675	470	474	487	rBV	178331	282098	9.99%	1.198%
2	5.128	547	551	567	rBV	2688674	2777046	98.31%	11.793%
3	6.181	725	730	744	rBV	2769840	2824773	100.00%	11.996%
4	6.287	744	748	752	rBV	2885581	2728755	96.60%	11.588%
5	6.510	783	786	790	rBV	854052	745871	26.40%	3.167%
6	6.669	809	813	817	rBV	2144324	1998215	70.74%	8.486%
7	7.081	878	883	886	rBV	1579835	1597987	56.57%	6.786%
8	7.792	1000	1004	1008	rBV	909916	891533	31.56%	3.786%
9	8.875	1182	1188	1191	rBV	2533172	2620888	92.78%	11.130%
10	9.539	1296	1301	1305	rBV	987001	923588	32.70%	3.922%
11	9.957	1369	1372	1373	rBV	34227	28869	1.02%	0.123%
12	10.327	1430	1435	1439	rBV	1495149	1409944	49.91%	5.988%
13	11.016	1547	1552	1555	rBV	1036843	904664	32.03%	3.842%
14	12.604	1817	1822	1825	rBV	2303648	2374448	84.06%	10.084%
15	13.639	1994	1998	2002	rBV	957323	874827	30.97%	3.715%
16	14.992	2223	2228	2234	rBV	448589	534058	18.91%	2.268%
17	15.821	2365	2369	2374	rBV3	17559	30125	1.07%	0.128%

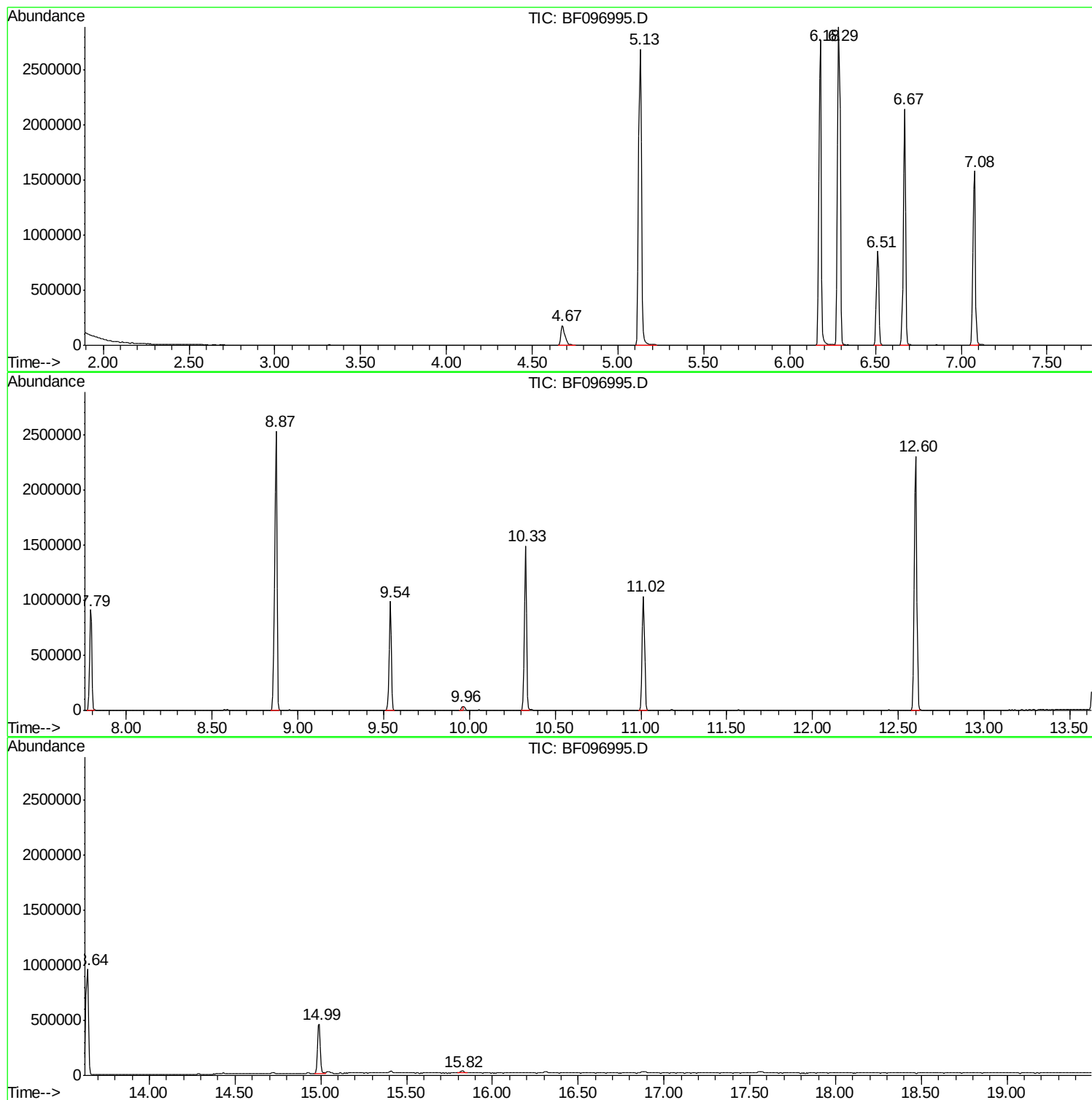
Sum of corrected areas: 23547689

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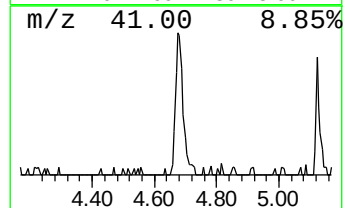
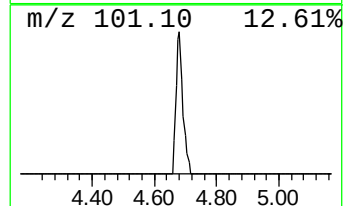
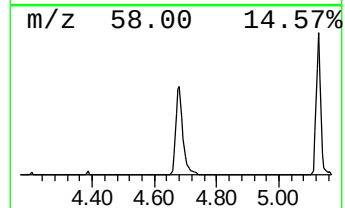
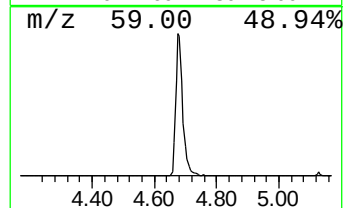
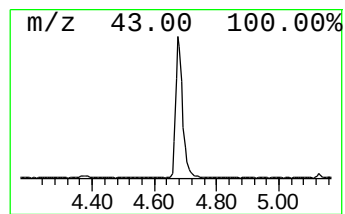
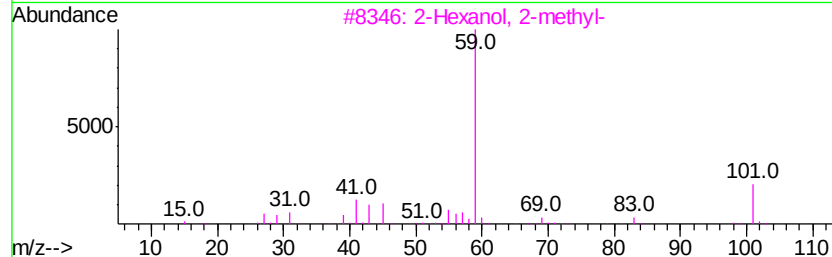
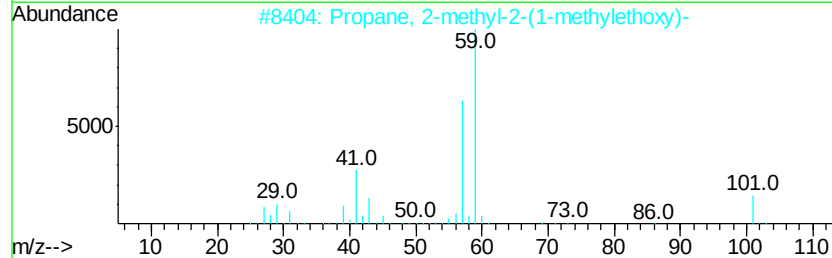
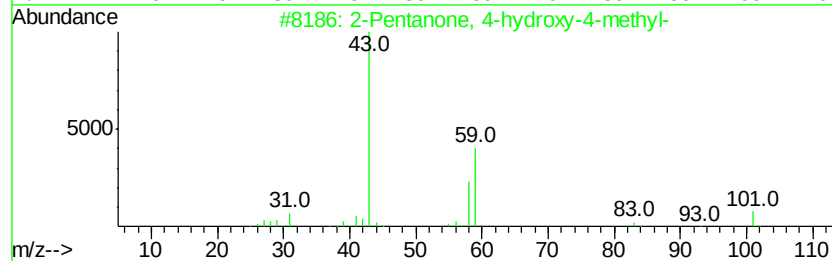
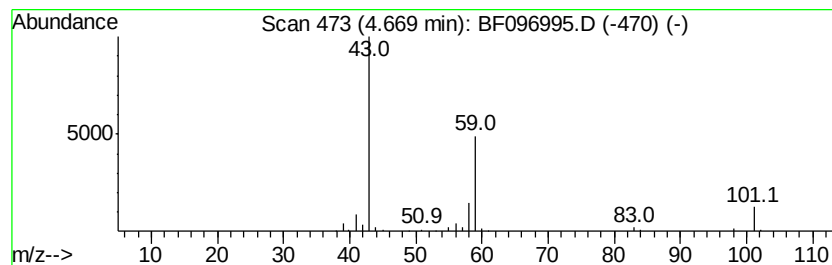
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.67	7.56 ng	282098	1,4-Dichlorobenzene-d4	6.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
3		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
4		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
5		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23



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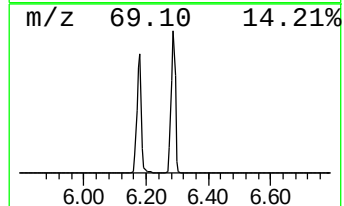
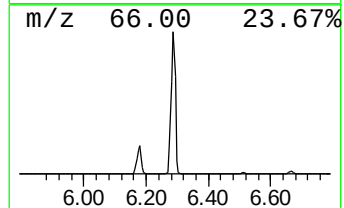
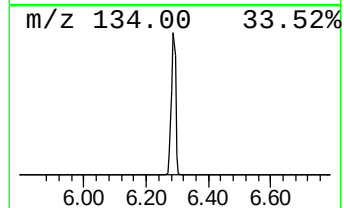
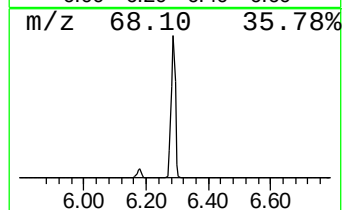
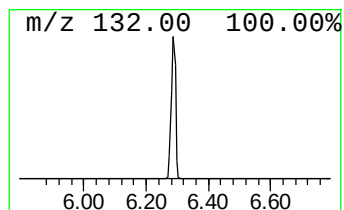
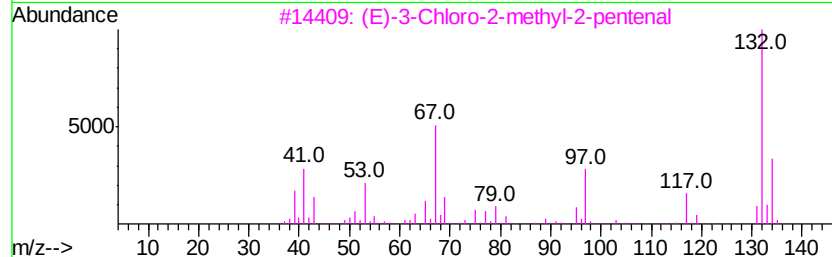
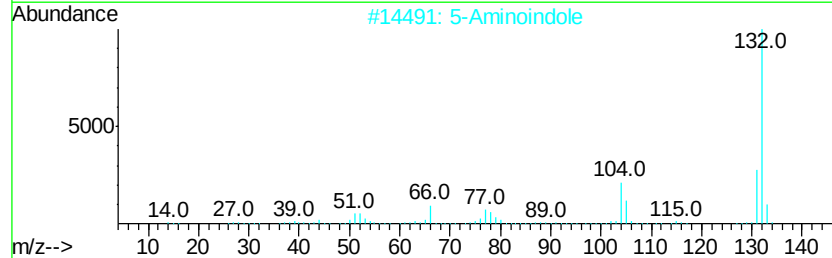
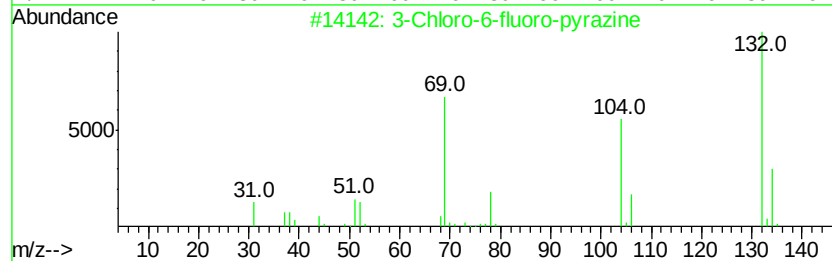
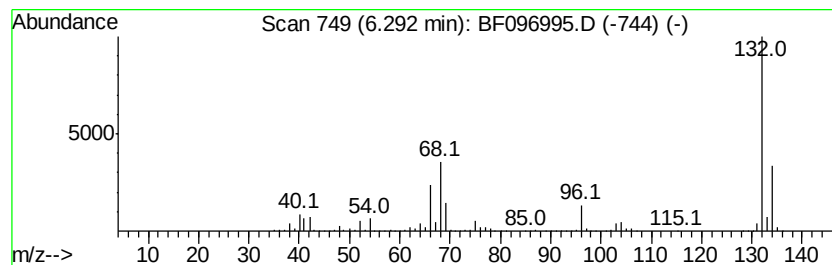
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Peak Number 2 unknown6.29 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.29	73.17 ng	2728760	1,4-Dichlorobenzene-d4	6.51

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine			132	C4H2ClFN2	1000146-10-7	27
2	5-Aminoindole			132	C8H8N2	005192-03-0	12
3	(E)-3-Chloro-2-methyl-2-pentenal			132	C6H9ClO	031357-76-3	9
4	Bicyclo[4.2.0]octa-1,3,5-triene, ...			132	C10H12	028749-81-7	9
5	Benzene, 1-ethenyl-3,5-dimethyl-			132	C10H12	005379-20-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.67	7.6	ng	282098	1	6.51	745871	20.0
unknown6.29	6.29	73.2	ng	2728760	1	6.51	745871	20.0