

Data Path : Z:\HPCHEM1\BNA_F\Data\BF070717\
 Data File : BF096552.D
 Acq On : 7 Jul 2017 12:03
 Operator : SJ/MA
 Sample : PB100467BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB100467BL

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.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	4.887	471	476	485	rBV	337139	469127	15.68%	1.834%
2	5.316	543	549	552	rBV	2736927	2893252	96.71%	11.313%
3	6.339	717	723	726	rBV	2650558	2857040	95.50%	11.171%
4	6.469	740	745	748	rBV	2896652	2809310	93.90%	10.984%
5	6.698	780	784	787	rBV	816197	677707	22.65%	2.650%
6	6.857	806	811	814	rBV	2178083	2122980	70.96%	8.301%
7	7.257	873	879	882	rBV	1791135	1813417	60.61%	7.090%
8	7.981	997	1002	1005	rBV	1005224	935374	31.27%	3.657%
9	9.057	1179	1185	1188	rBV	3077136	2991709	100.00%	11.698%
10	9.727	1294	1299	1303	rBV	1154737	1043839	34.89%	4.081%
11	10.516	1427	1433	1436	rBV	1777883	1784713	59.66%	6.978%
12	11.210	1546	1551	1554	rBV	1111241	1023272	34.20%	4.001%
13	12.798	1815	1821	1824	rBV	2666484	2799705	93.58%	10.947%
14	13.833	1993	1997	2001	rBV	815575	762459	25.49%	2.981%
15	15.239	2231	2236	2241	rVB	483390	591573	19.77%	2.313%

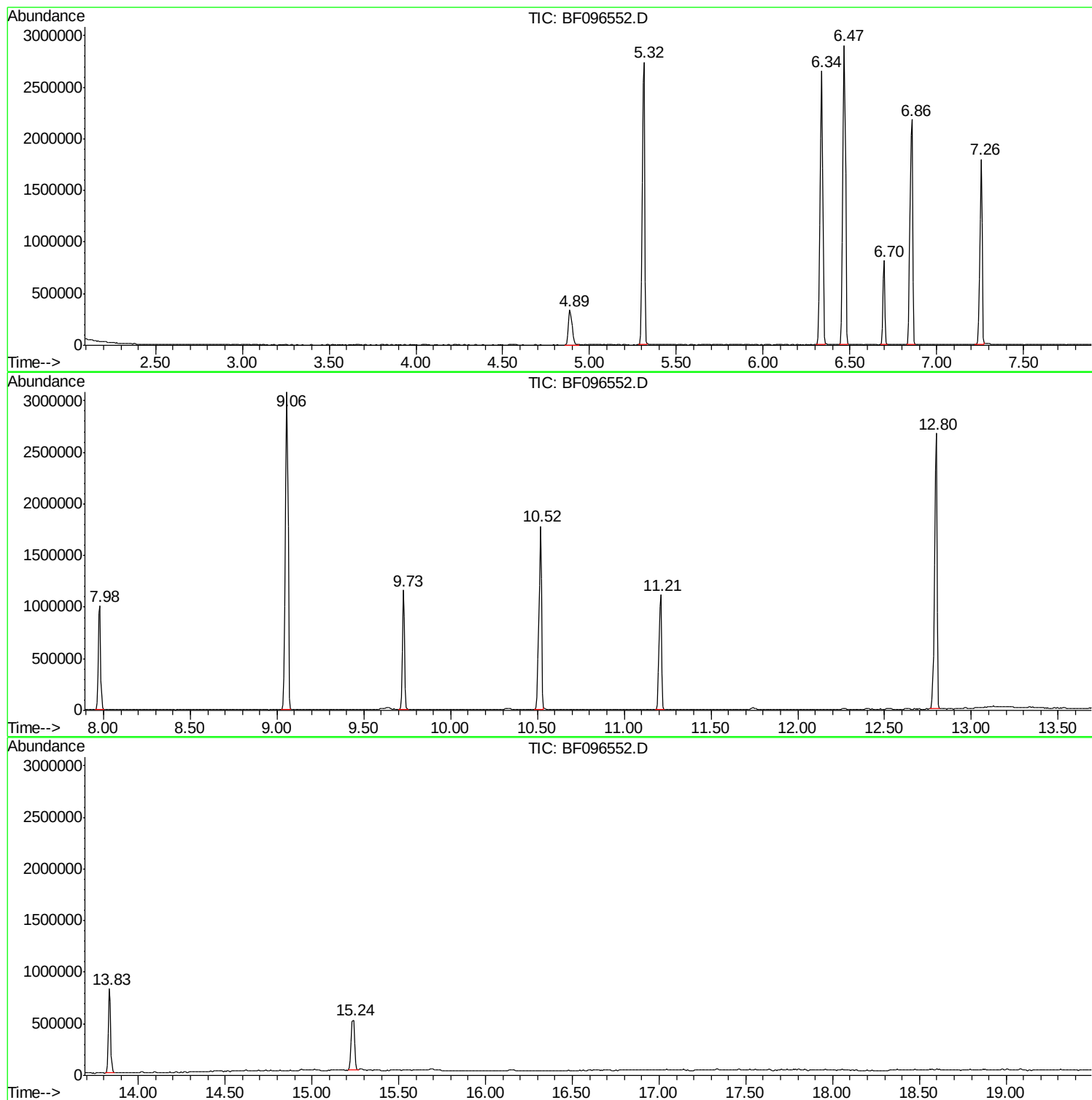
Sum of corrected areas: 25575477

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.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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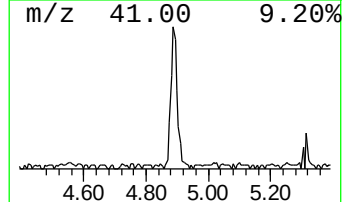
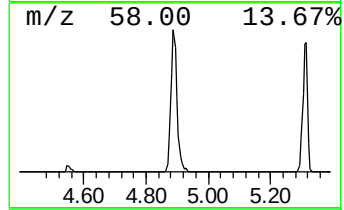
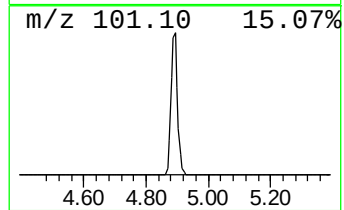
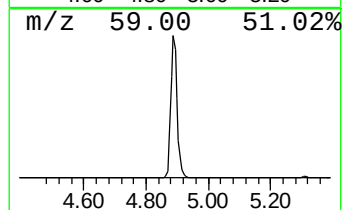
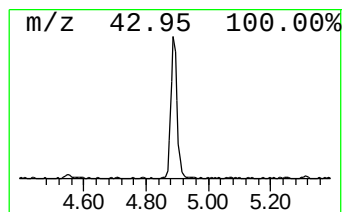
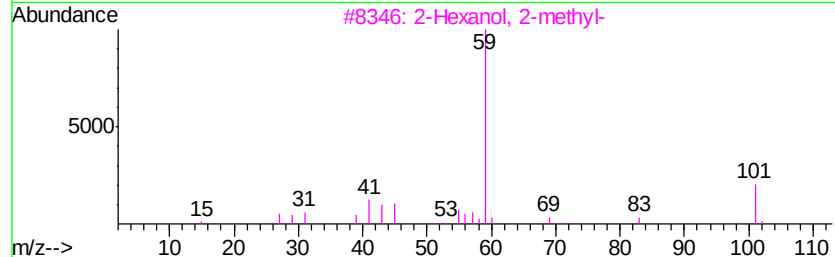
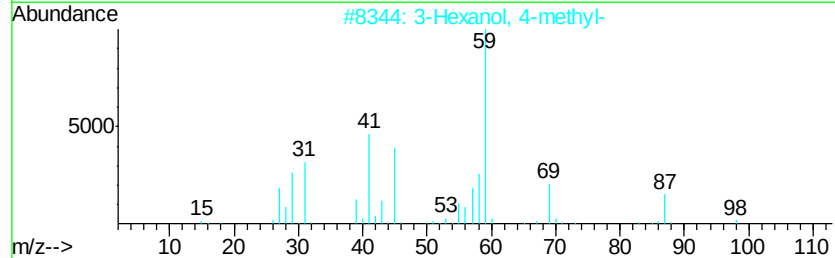
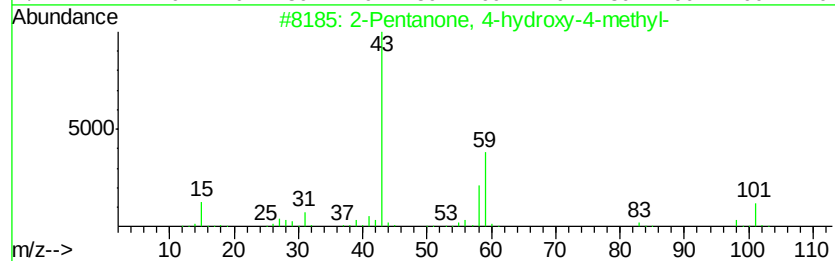
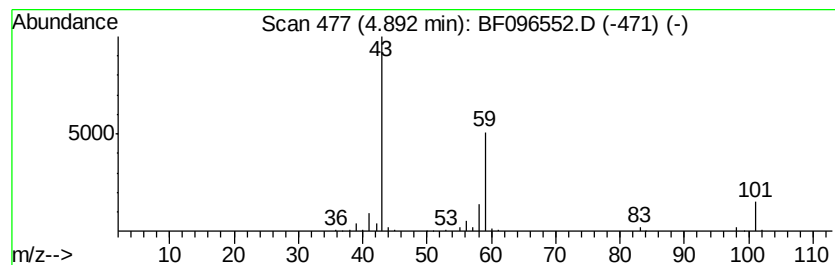
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.89	13.84 ng	469127	1,4-Dichlorobenzene-d4	6.70

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			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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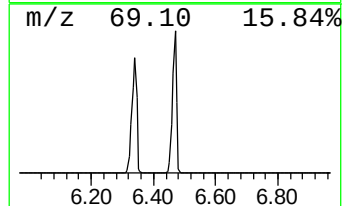
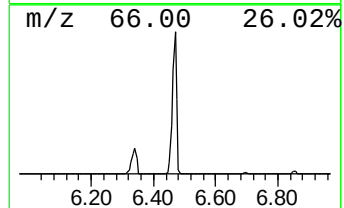
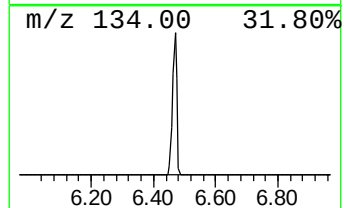
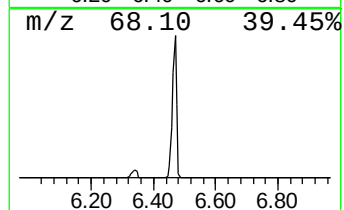
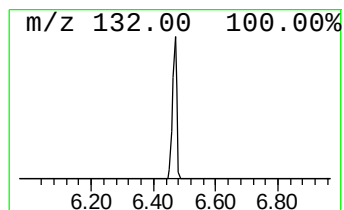
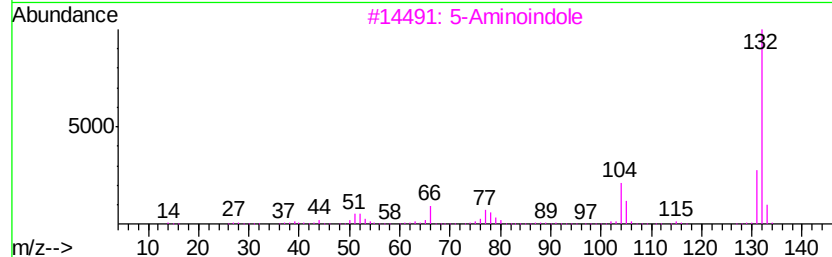
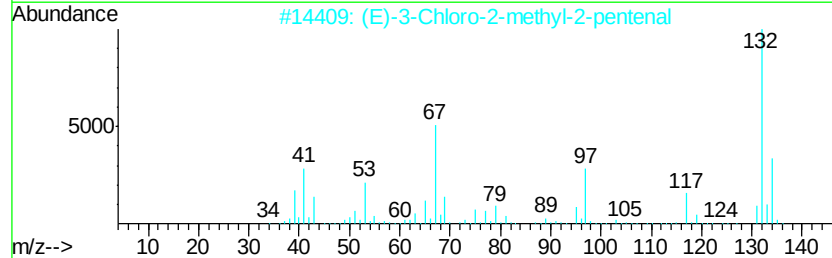
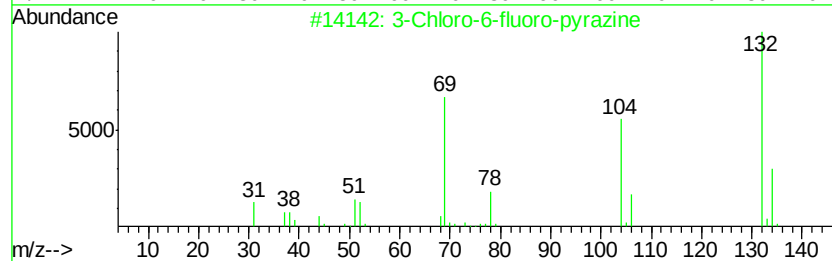
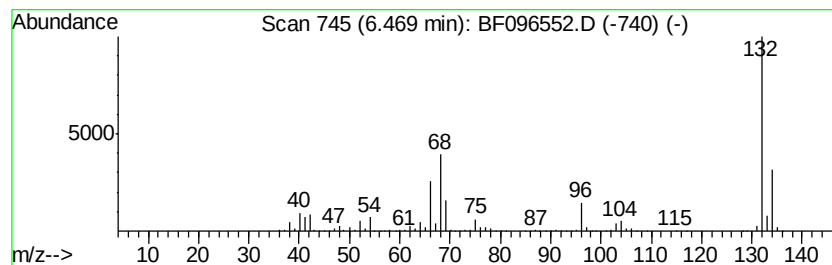
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Peak Number 2 unknown6.47 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.47	82.91 ng	2809310	1,4-Dichlorobenzene-d4	6.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.89	13.8	ng	469127	1	6.70	677707	20.0
unknown6.47	6.47	82.9	ng	2809310	1	6.70	677707	20.0