

Data Path : Z:\HPCHEM1\BNA F\DATA\BF112217\
 Data File : BF100872.D
 Acq On : 23 Nov 2017 13:45
 Operator : SJ/JU
 Sample : I6567-04
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 112017-TIDO-F-U-S

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-BF110817.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	5.075	504	508	515	rBV	136920	143413	5.89%	1.078%
2	5.469	572	575	579	rBV	640331	614645	25.24%	4.618%
3	6.481	743	747	755	rVB	421225	422921	17.37%	3.178%
4	6.628	767	772	775	rBV	1212388	1140402	46.83%	8.569%
5	6.857	807	811	814	rBV	516673	414371	17.02%	3.114%
6	7.016	833	838	841	rBV	1817563	1613619	66.26%	12.125%
7	7.422	901	907	910	rBV	1227625	1322282	54.30%	9.936%
8	8.139	1024	1029	1032	rBV	664354	540199	22.18%	4.059%
9	9.216	1206	1212	1215	rBV	2613314	2435142	100.00%	18.298%
10	9.604	1275	1278	1281	rBV	65524	48931	2.01%	0.368%
11	9.892	1321	1327	1331	rVB2	680405	605582	24.87%	4.550%
12	10.316	1391	1399	1402	rBV2	31643	30515	1.25%	0.229%
13	10.563	1438	1441	1444	rVB	35499	26392	1.08%	0.198%
14	10.680	1457	1461	1465	rBV	935117	855708	35.14%	6.430%
15	10.786	1476	1479	1487	rVB	32007	29522	1.21%	0.222%
16	11.380	1576	1580	1583	rBV	670941	542403	22.27%	4.076%
17	12.969	1845	1850	1853	rBV	1899944	1763261	72.41%	13.249%
18	14.021	2025	2029	2032	rBV	388639	359337	14.76%	2.700%
19	15.492	2273	2279	2287	rBV	320412	399711	16.41%	3.003%

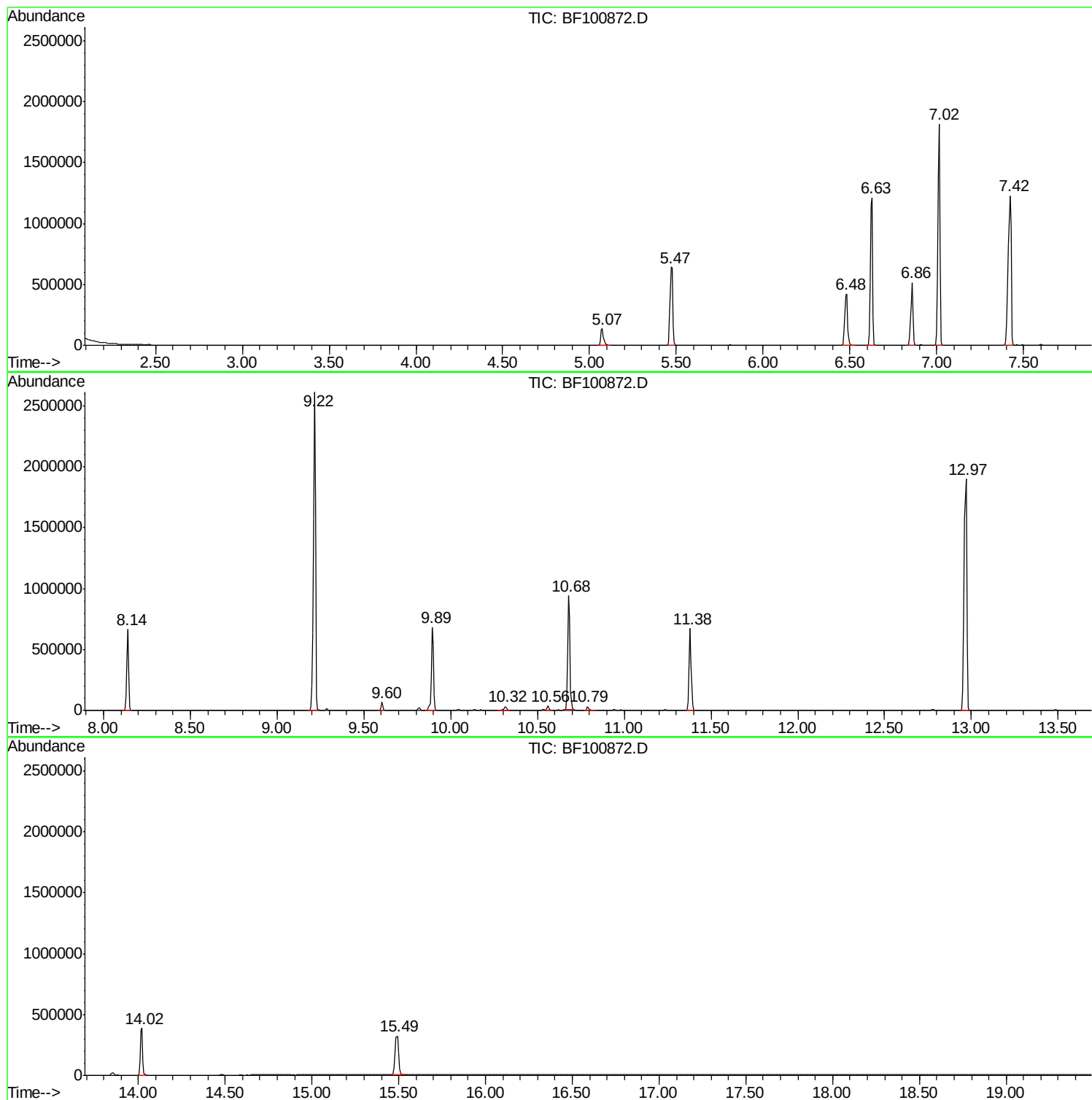
Sum of corrected areas: 13308356

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110817.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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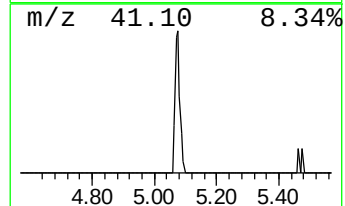
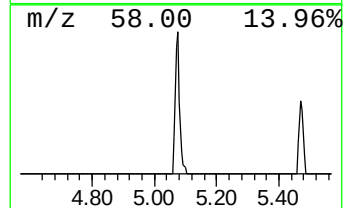
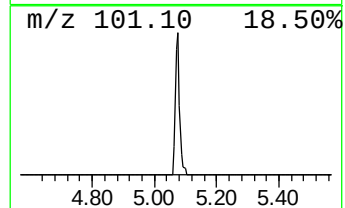
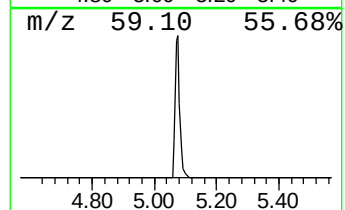
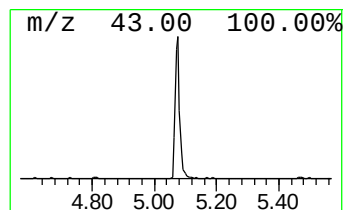
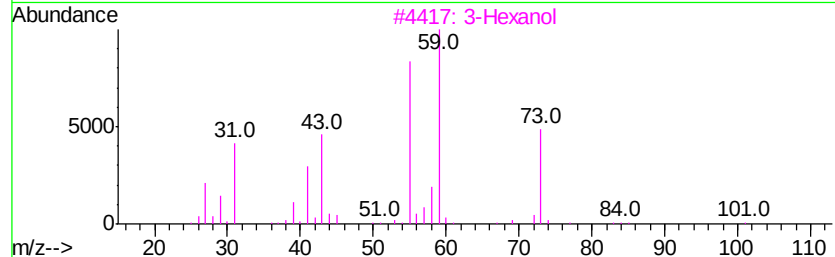
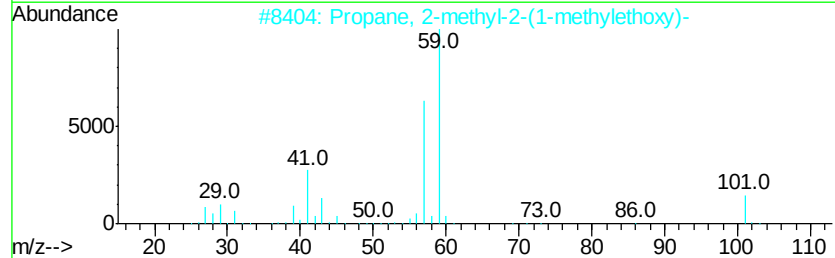
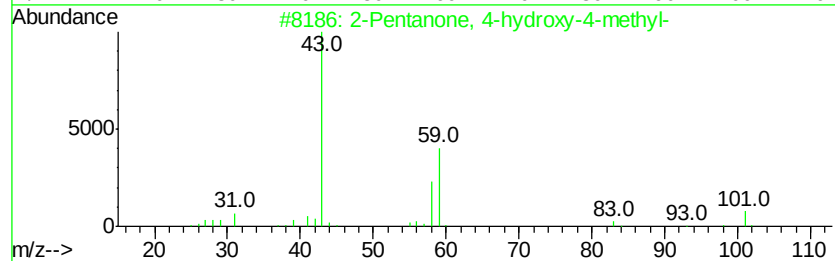
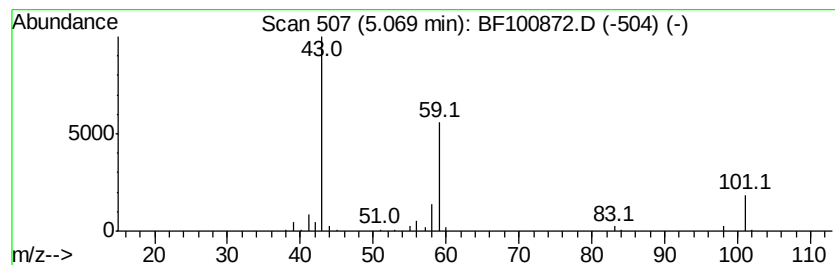
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.07	6.92 ng	143413	1,4-Dichlorobenzene-d4	6.86

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
2			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
3			3-Hexanol	102	C6H14O	000623-37-0	28
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23



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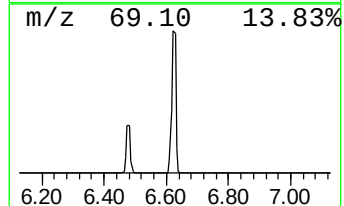
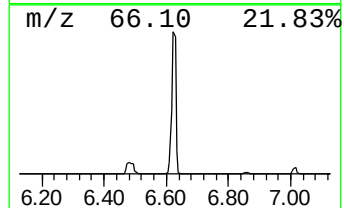
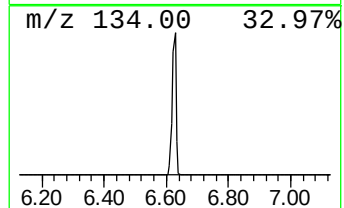
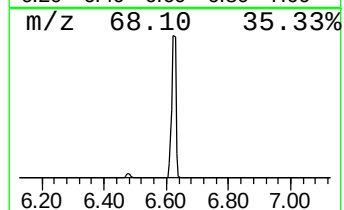
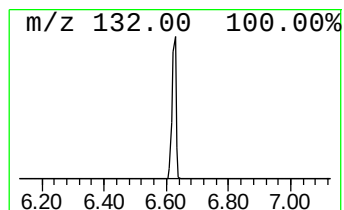
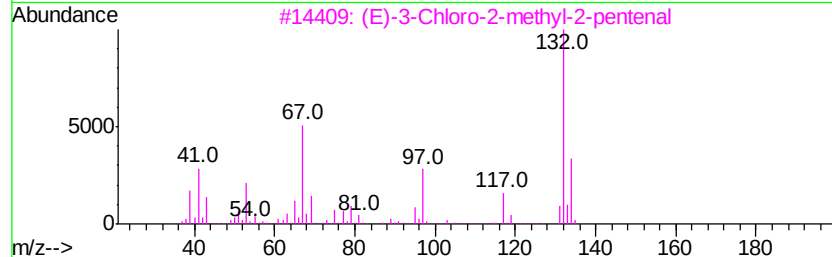
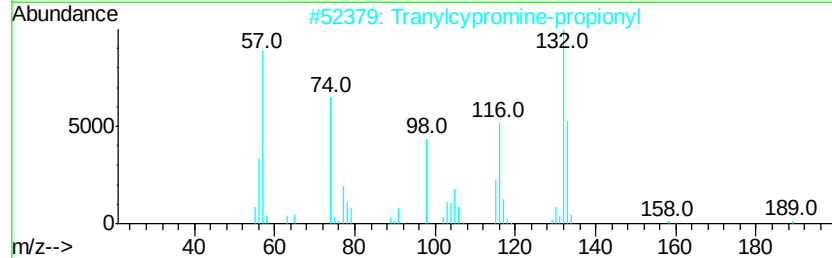
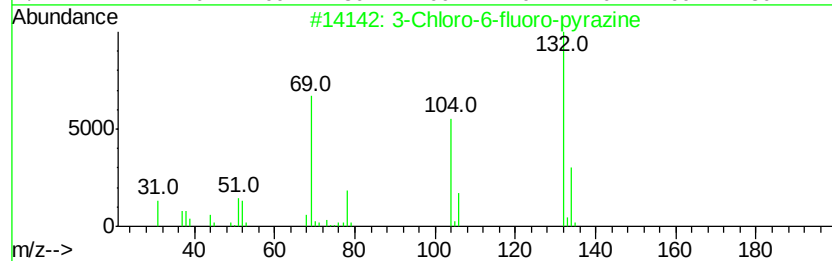
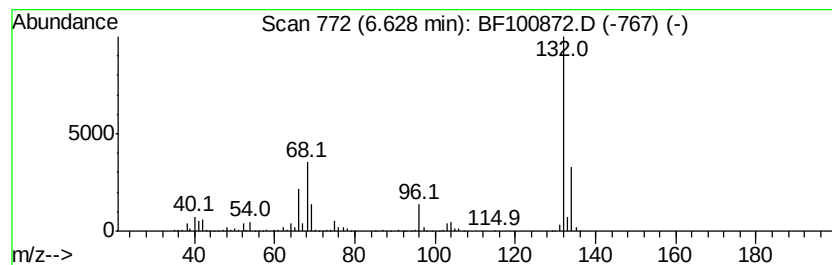
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Peak Number 2 unknown6.63 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	55.04 ng	1140400	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.07	6.9	ng	143413	1	6.86	414371	20.0
unknown6.63	6.63	55.0	ng	1140400	1	6.86	414371	20.0