

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112717\
 Data File : BF100909.D
 Acq On : 27 Nov 2017 23:46
 Operator : SJ/JU
 Sample : PB104511BL
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB104511BL

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	5.087	506	510	518	rBV	265795	303534	16.21%	2.502%
2	5.481	573	577	581	rBV	1040878	999311	53.38%	8.238%
3	6.486	743	748	751	rBV	1035815	973558	52.01%	8.026%
4	6.628	767	772	775	rBV	1199747	1043255	55.73%	8.600%
5	6.857	807	811	814	rBV	455968	364422	19.47%	3.004%
6	7.016	833	838	841	rBV	1343782	1247157	66.62%	10.281%
7	7.422	901	907	910	rBV	957608	945706	50.52%	7.796%
8	8.139	1024	1029	1032	rBV	546534	466953	24.94%	3.849%
9	9.216	1206	1212	1215	rBV	2014148	1872036	100.00%	15.433%
10	9.892	1322	1327	1331	rBV	555402	517388	27.64%	4.265%
11	10.680	1457	1461	1476	rBV	691114	640431	34.21%	5.280%
12	11.380	1576	1580	1584	rBV	603047	496039	26.50%	4.089%
13	12.968	1845	1850	1853	rBV	1711787	1559903	83.33%	12.860%
14	14.021	2025	2029	2036	rBV	425477	373568	19.96%	3.080%
15	15.492	2273	2279	2290	rBV	253788	326974	17.47%	2.696%

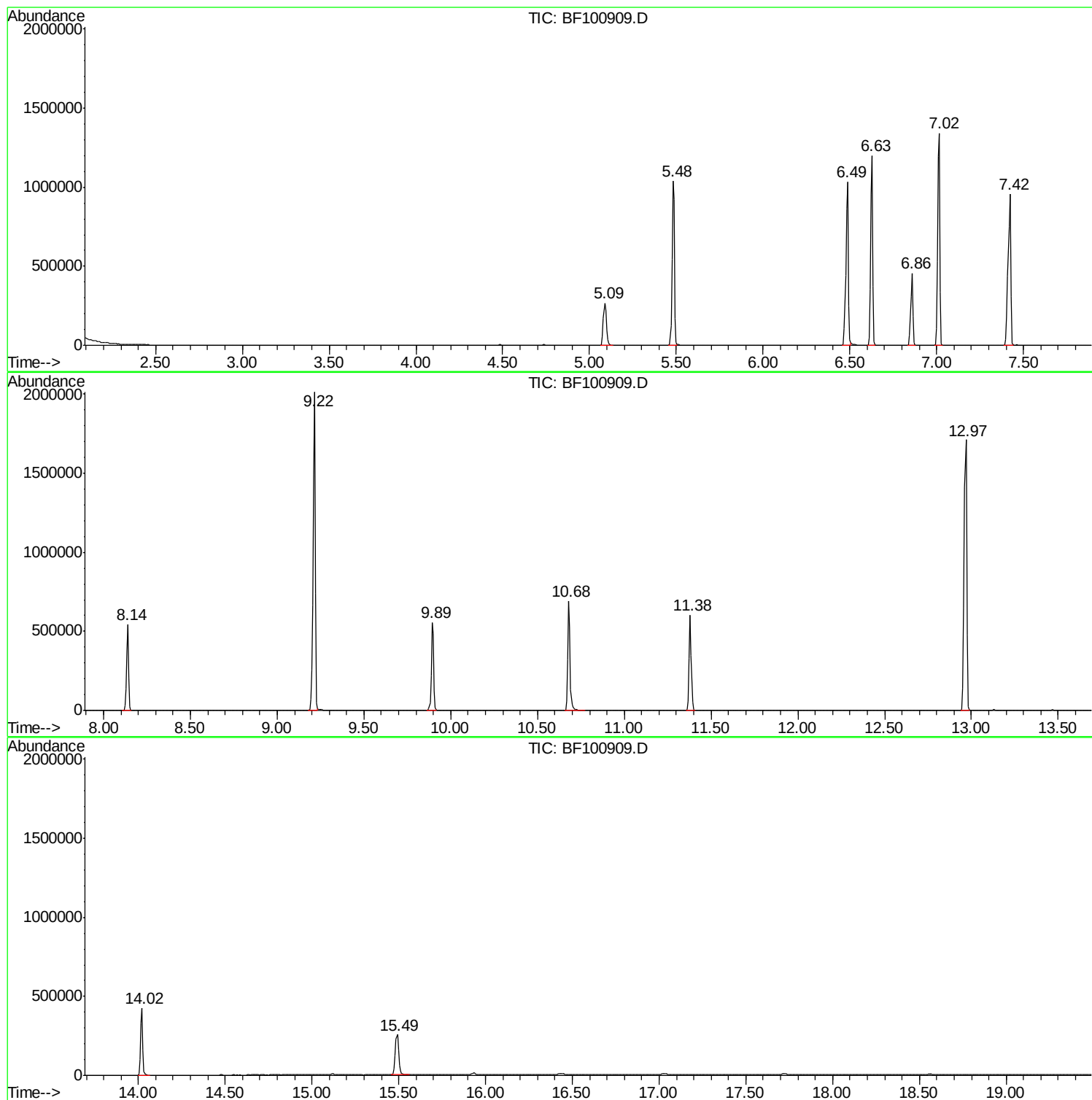
Sum of corrected areas: 12130235

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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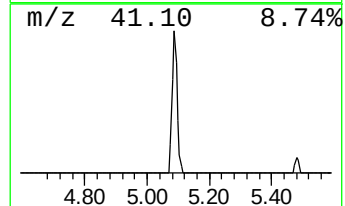
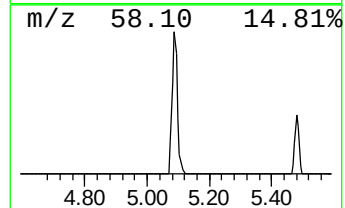
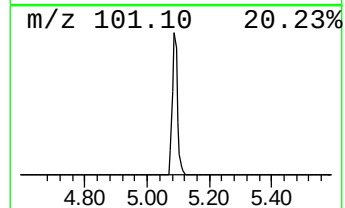
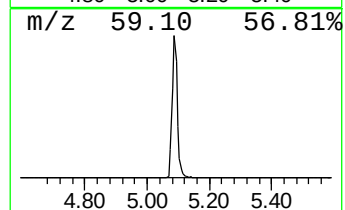
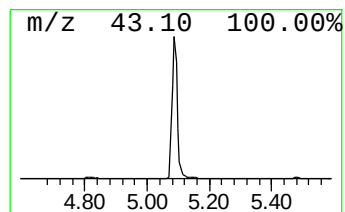
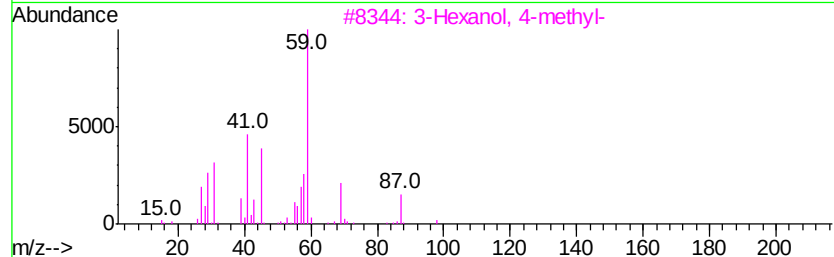
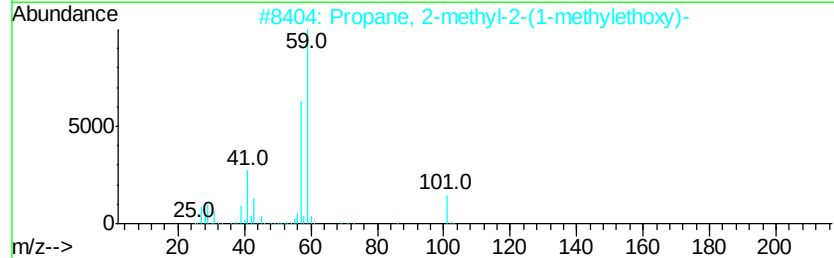
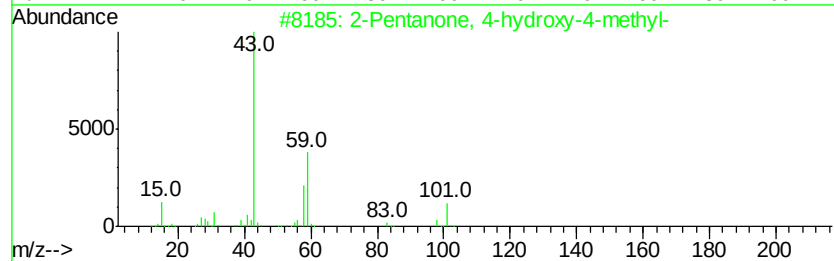
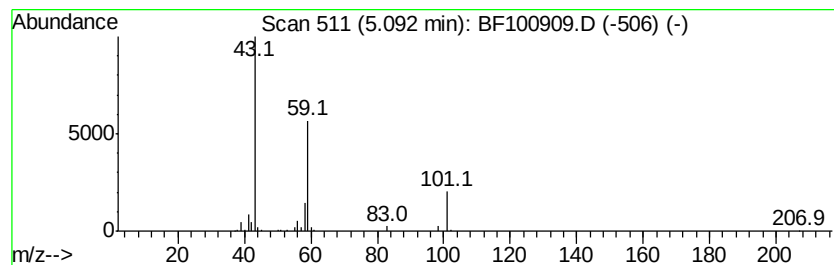
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.09	16.66 ng	303534	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	64
2			Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	33
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4			Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	25
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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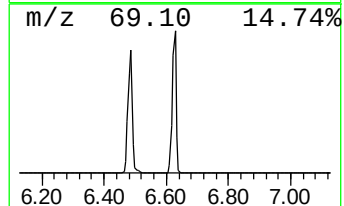
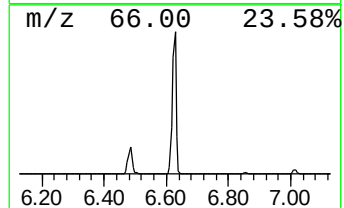
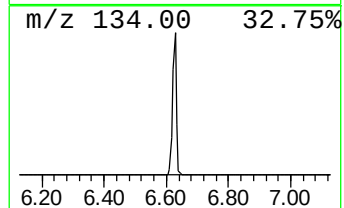
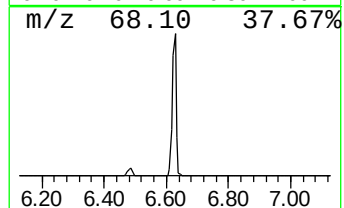
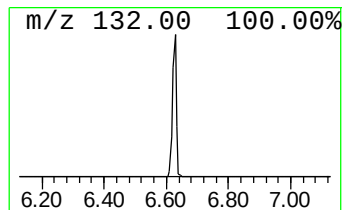
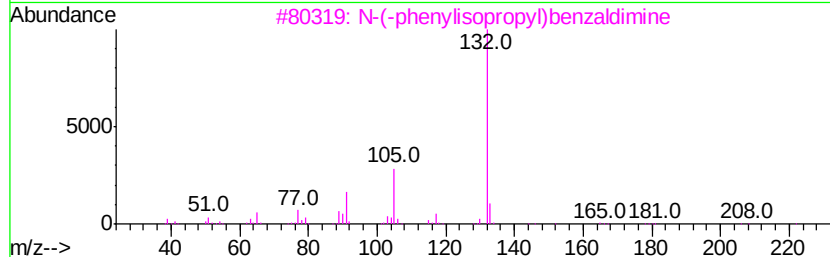
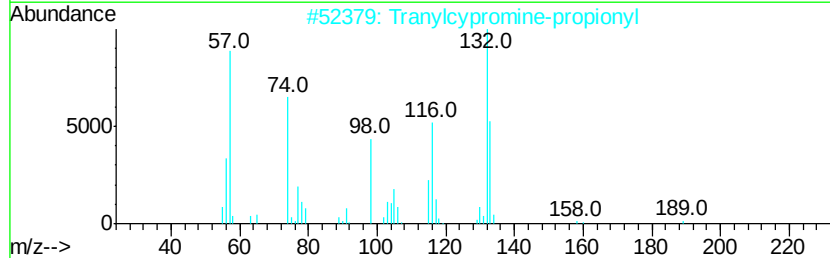
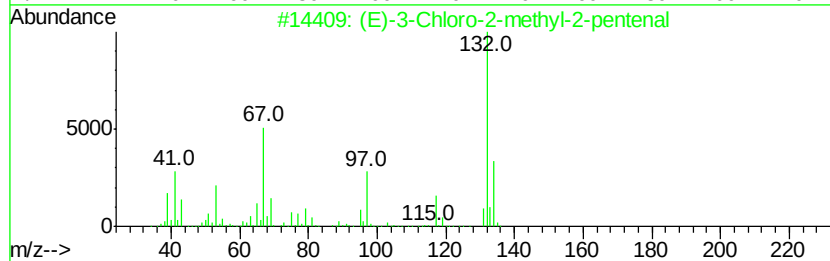
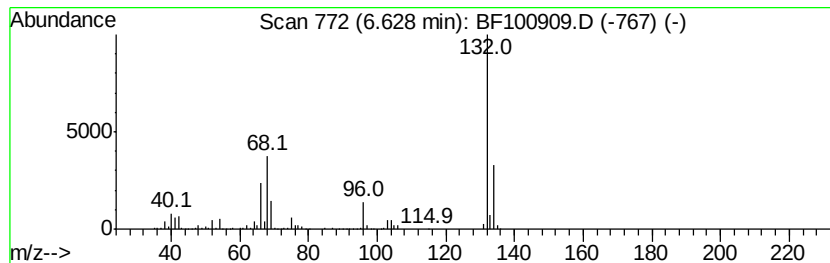
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Peak Number 2 unknown6.63 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	57.26 ng	1043260	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		N-(-phenylisopropyl)benzalimine	223	C16H17N	1000379-01-3	12
4		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.09	16.7	ng	303534	1	6.86	364422	20.0
unknown6.63	6.63	57.3	ng	1043260	1	6.86	364422	20.0