

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112817\
 Data File : BF100964.D
 Acq On : 29 Nov 2017 5:45
 Operator : SJ/JU
 Sample : I6603-06
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 112217-JETT-F-D-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	505	508	518	rBV	94158	100784	4.39%	0.824%
2	5.475	572	576	589	rBV	690717	594228	25.89%	4.856%
3	6.481	744	747	755	rVB	452176	394317	17.18%	3.222%
4	6.628	768	772	775	rBV	1316090	1094331	47.68%	8.942%
5	6.857	808	811	815	rBV	482343	378997	16.51%	3.097%
6	7.016	834	838	841	rBV	1770966	1565282	68.20%	12.791%
7	7.428	902	908	911	rBV	1229674	1205813	52.54%	9.853%
8	8.139	1025	1029	1032	rBV	599013	474959	20.70%	3.881%
9	9.216	1207	1212	1215	rBV	2338378	2295014	100.00%	18.753%
10	9.892	1322	1327	1338	rVB	538444	514862	22.43%	4.207%
11	10.686	1458	1462	1473	rBV	791394	767812	33.46%	6.274%
12	11.380	1576	1580	1584	rBV	552374	456394	19.89%	3.729%
13	12.968	1845	1850	1853	rBV	1940337	1723515	75.10%	14.084%
14	14.021	2025	2029	2036	rVB	417943	355928	15.51%	2.908%
15	15.492	2274	2279	2287	rBV	253015	315560	13.75%	2.579%

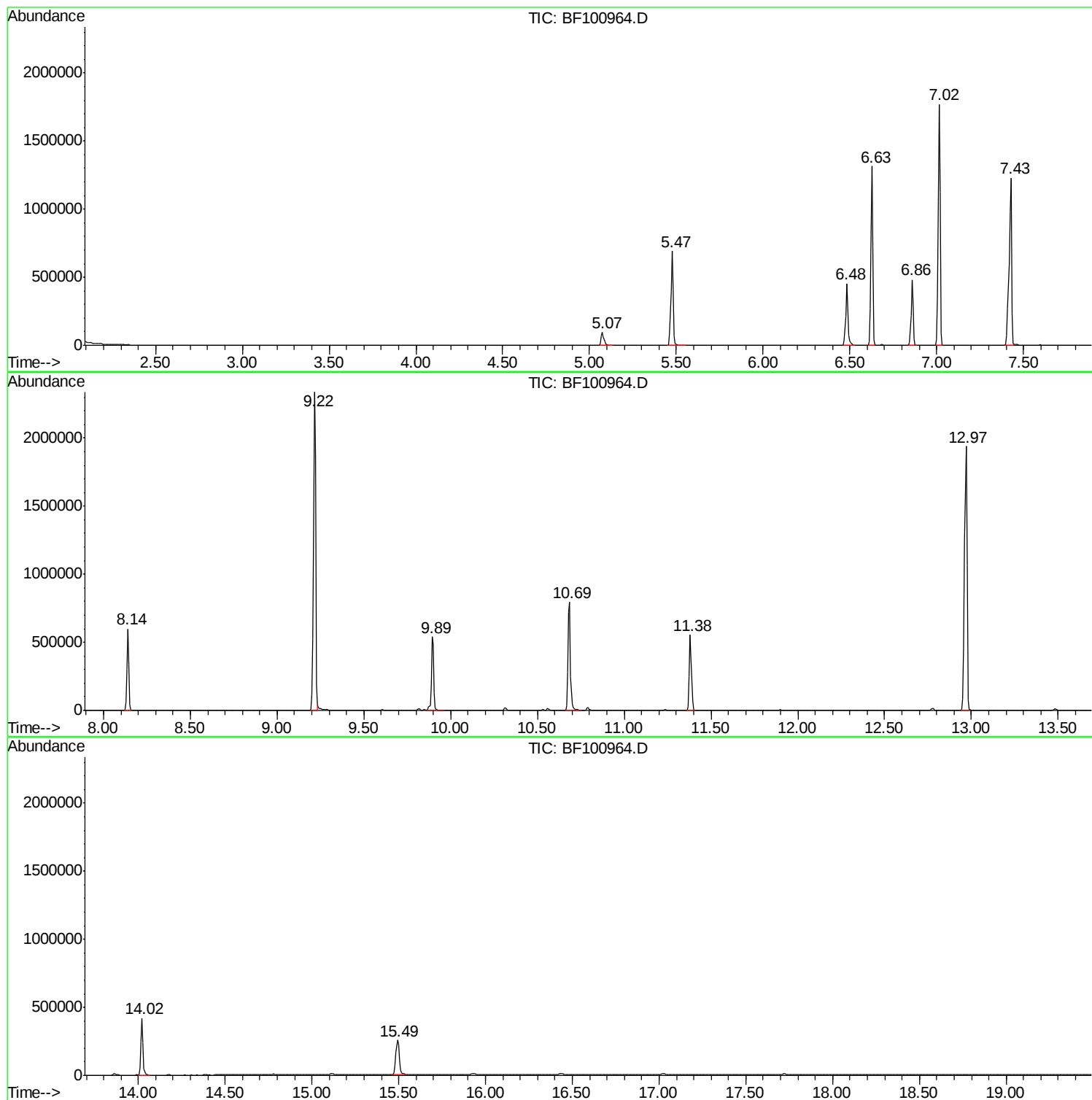
Sum of corrected areas: 12237796

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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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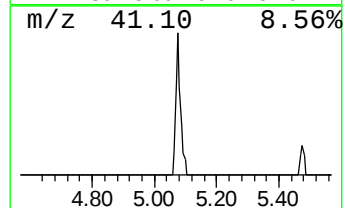
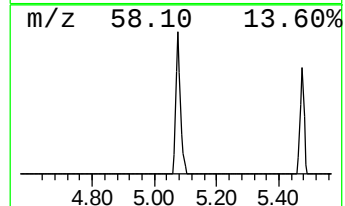
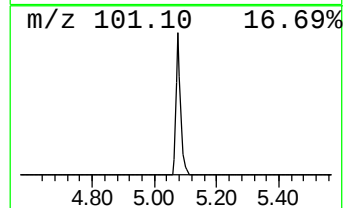
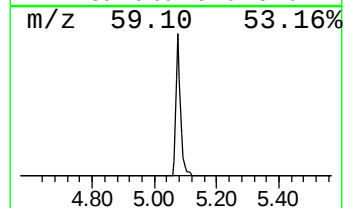
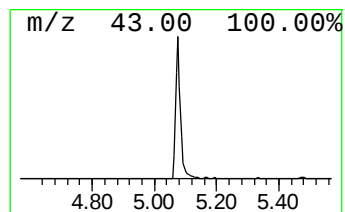
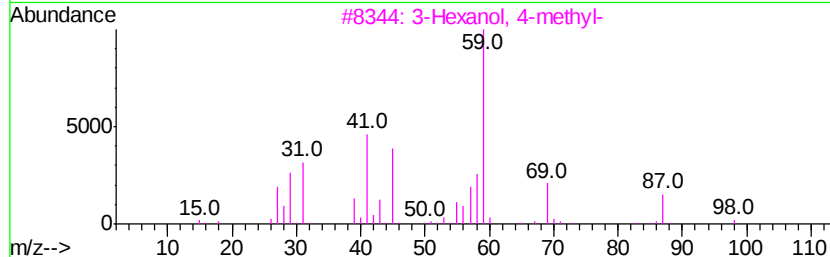
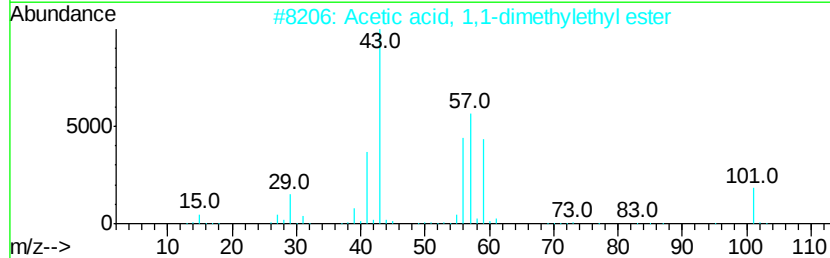
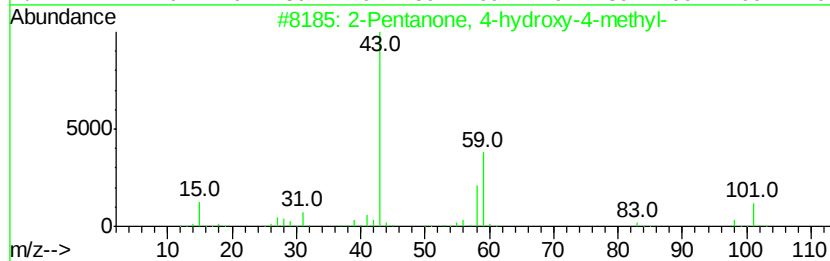
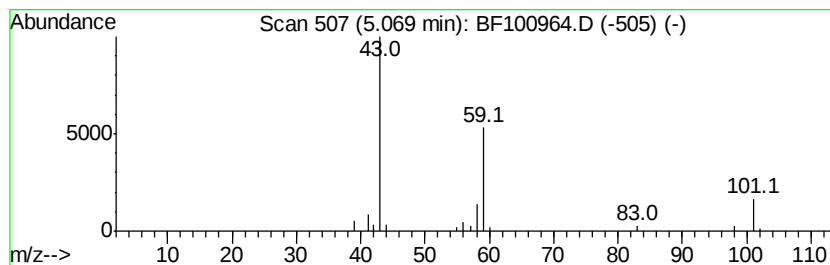
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.07	5.32 ng	100784	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			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
5			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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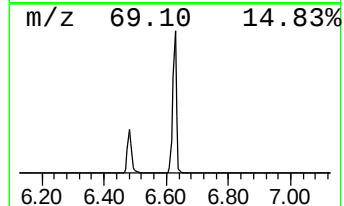
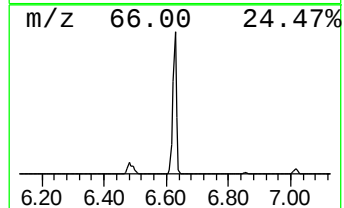
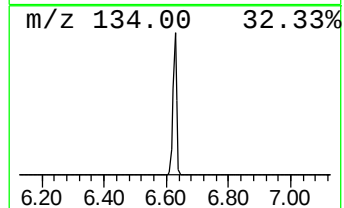
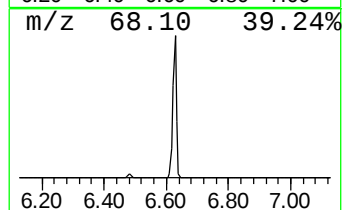
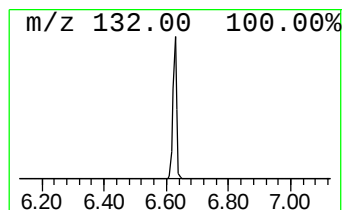
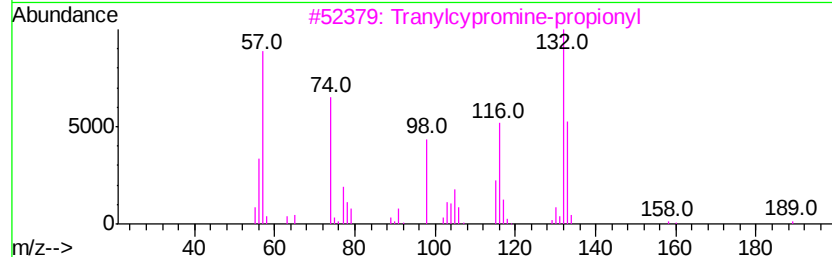
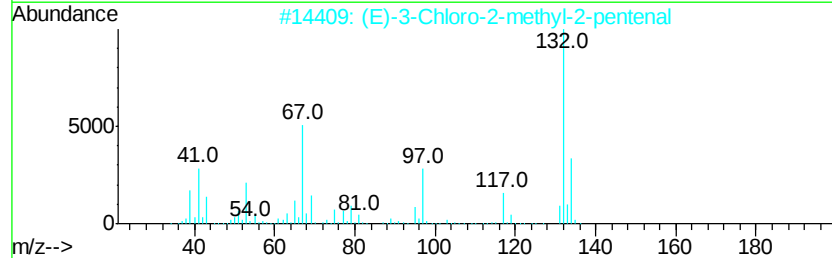
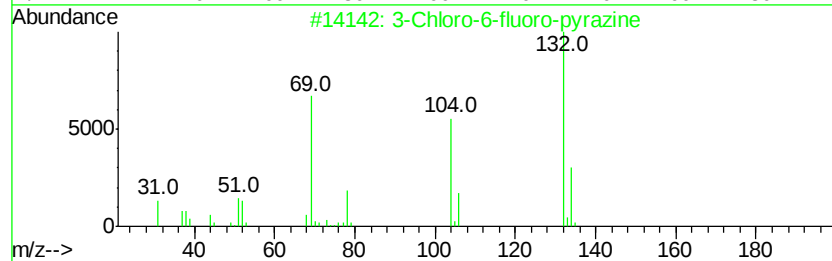
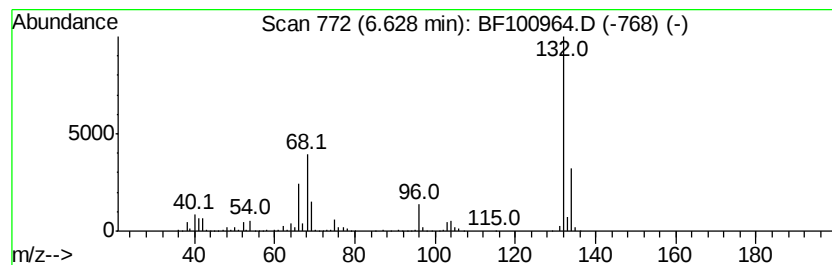
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Peak Number 2 unknown6.63 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	57.75 ng	1094330	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	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
5		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.07	5.3	ng	100784	1	6.86	378997	20.0
unknown6.63	6.63	57.8	ng	1094330	1	6.86	378997	20.0