

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF061419\
 Data File : BF115039.D
 Acq On : 14 Jun 2019 14:55
 Operator : HP/JU
 Sample : PB120599BL
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB120599BL

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061219.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	4.834	463	467	480	rBV	262993	351848	5.94%	0.817%
2	5.257	535	539	543	rBV	3099612	3176646	53.59%	7.379%
3	6.287	709	714	717	rBV	3414558	3073097	51.84%	7.139%
4	6.416	731	736	739	rBV	3763330	3453018	58.25%	8.021%
5	6.645	771	775	779	rBV	1715756	1409172	23.77%	3.274%
6	6.804	797	802	805	rBV	4394151	4071109	68.68%	9.457%
7	7.210	865	871	874	rBV	2878745	3092048	52.16%	7.183%
8	7.928	988	993	1005	rBV	1808611	1845281	31.13%	4.287%
9	9.004	1170	1176	1179	rBV	5765578	5927747	100.00%	13.770%
10	9.675	1284	1290	1305	rBV	2151764	2116042	35.70%	4.916%
11	10.457	1419	1423	1449	rBV	2456460	2448229	41.30%	5.687%
12	11.145	1536	1540	1548	rBV	2460818	2262681	38.17%	5.256%
13	12.733	1805	1810	1813	rBV	5583055	5755984	97.10%	13.371%
14	13.768	1982	1986	2001	rBV	2157594	2095766	35.36%	4.868%
15	15.145	2215	2220	2226	rBV	1551053	1885729	31.81%	4.381%
16	15.198	2226	2229	2238	rVB	56065	83234	1.40%	0.193%

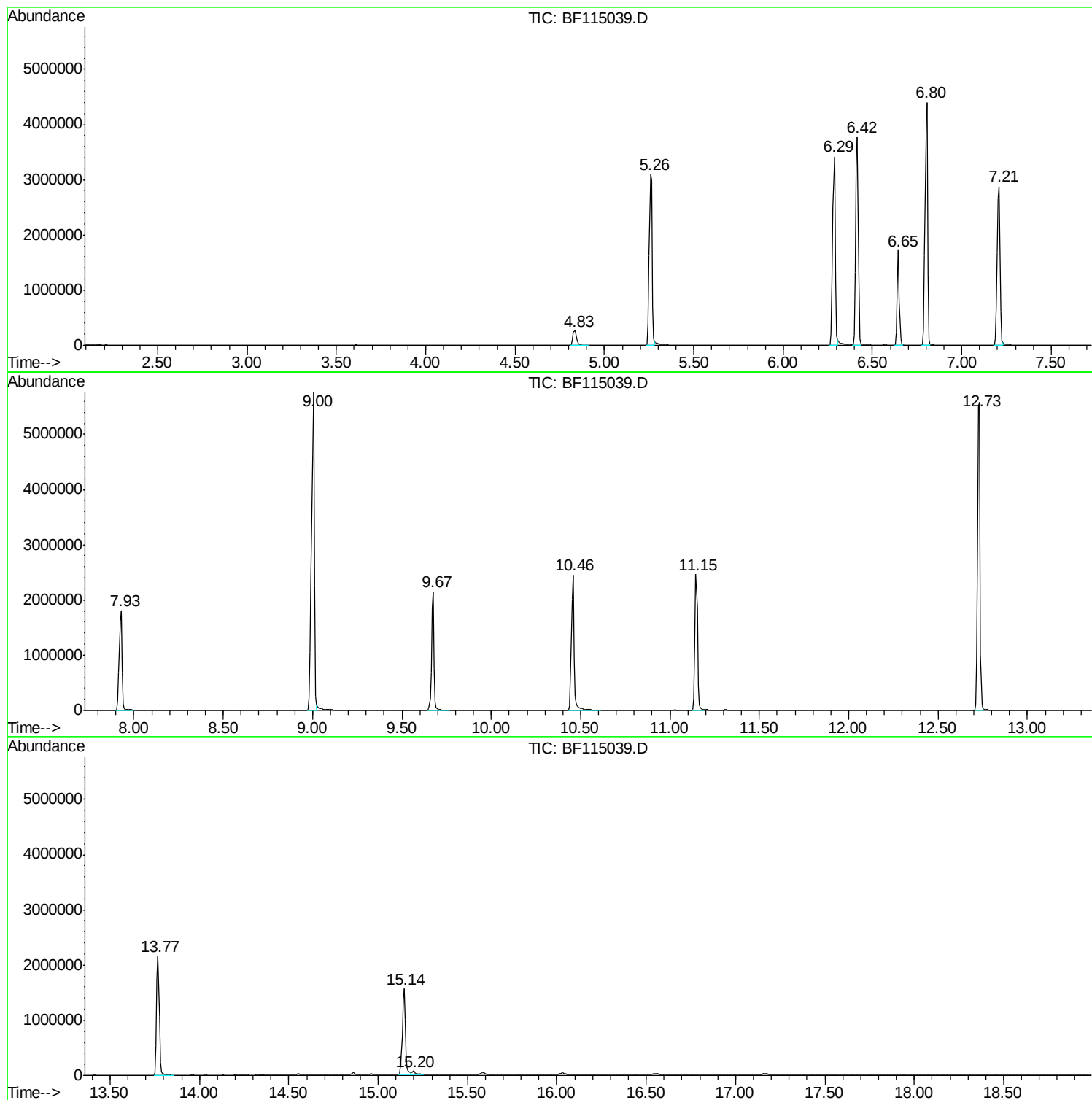
Sum of corrected areas: 43047631

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061219.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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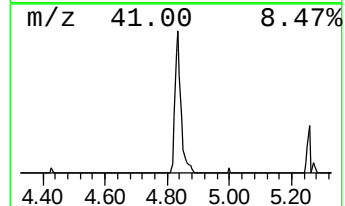
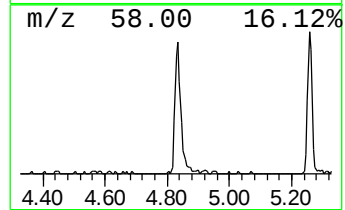
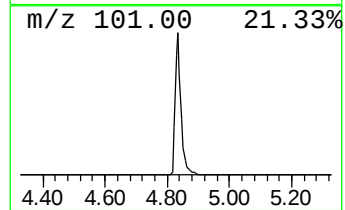
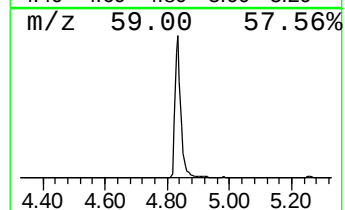
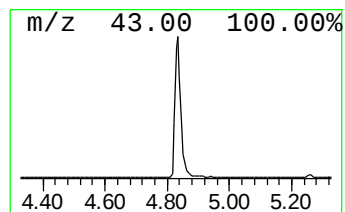
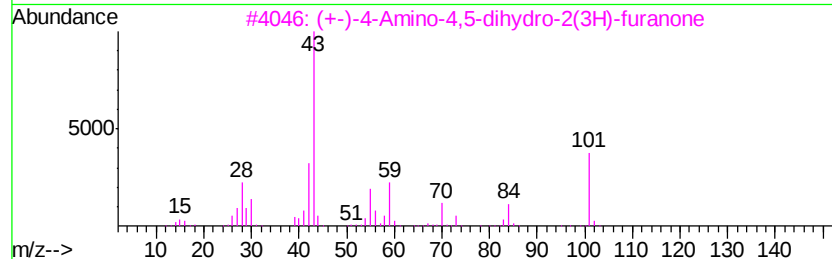
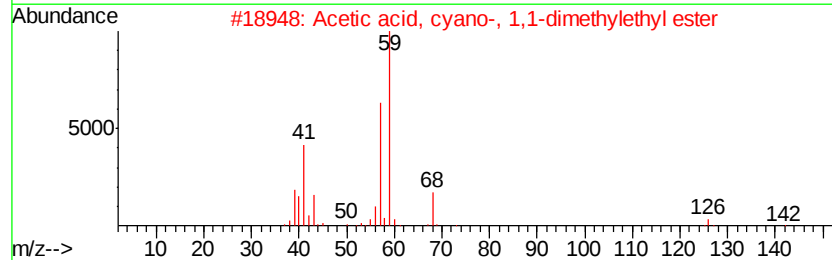
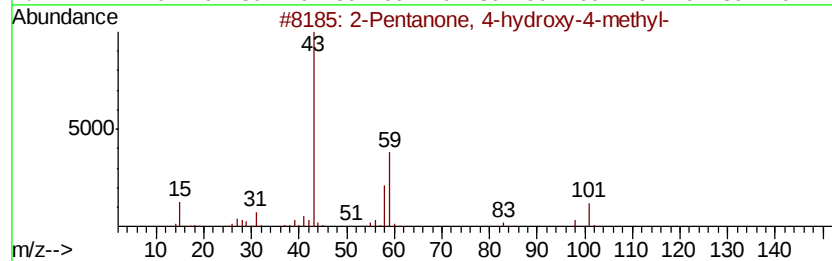
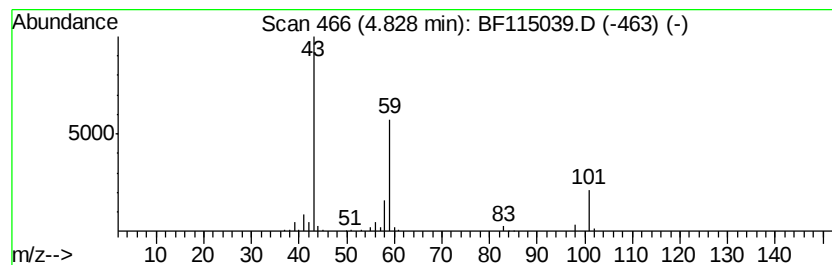
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.83	4.99 ng	351848	1,4-Dichlorobenzene-d4	6.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
4		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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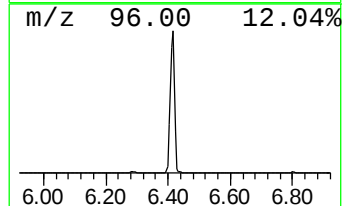
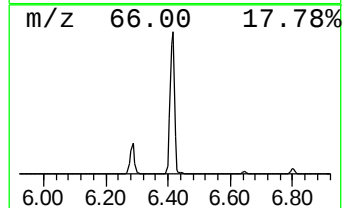
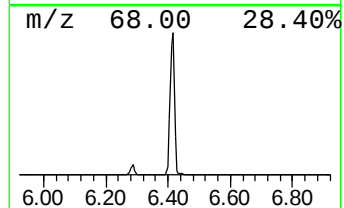
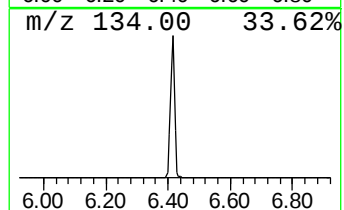
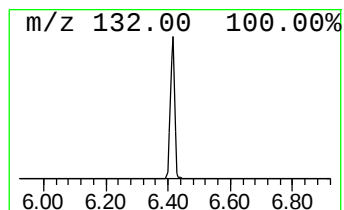
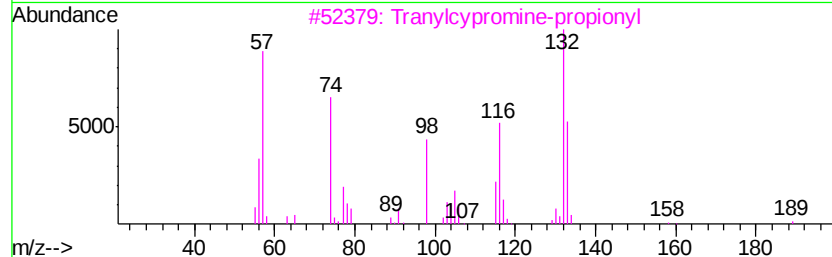
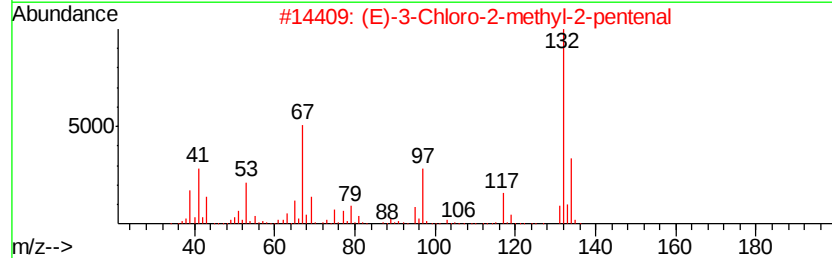
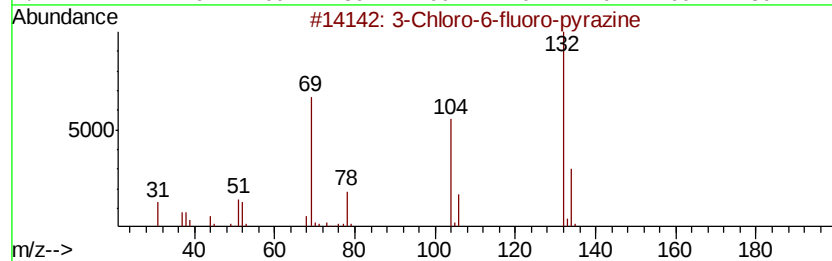
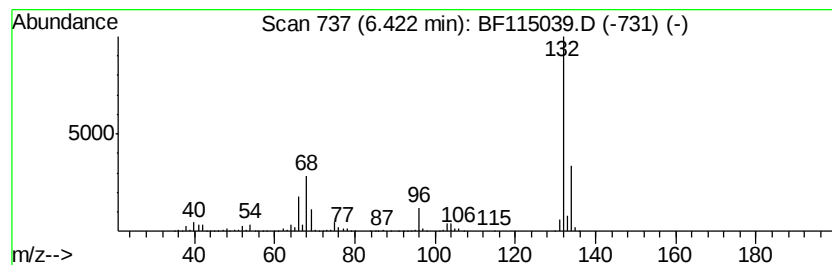
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Peak Number 2 unknown6.42 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.42	49.01 ng	3453020	1,4-Dichlorobenzene-d4	6.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.83	5.0	ng	351848	1	6.65	1409170	20.0
unknown6.42	6.42	49.0	ng	3453020	1	6.65	1409170	20.0