

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF080918\
 Data File : BF108074.D
 Acq On : 10 Aug 2018 2:02
 Operator : JU/SJ
 Sample : J4380-06
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FIELD-BLANK

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-BF080818.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	5.248	490	495	501	rBV	716954	763741	8.35%	1.281%
2	5.636	556	561	564	rBV	3698337	3576278	39.08%	5.999%
3	6.625	724	729	732	rBV	2743151	2288621	25.01%	3.839%
4	6.783	751	756	759	rBV	7324882	6803691	74.35%	11.412%
5	7.013	792	795	799	rVB	2131701	1792213	19.58%	3.006%
6	7.172	818	822	826	rVB	6737034	6219831	67.97%	10.433%
7	7.577	885	891	894	rBV	5593409	5738482	62.71%	9.626%
8	8.301	1009	1014	1017	rBV	2317138	2175310	23.77%	3.649%
9	9.377	1191	1197	1200	rBV	9419427	9151483	100.00%	15.350%
10	10.060	1309	1313	1317	rVB	2362738	2237773	24.45%	3.754%
11	10.854	1443	1448	1463	rBV	5060805	4988557	54.51%	8.368%
12	11.560	1564	1568	1571	rBV	2310129	1986167	21.70%	3.332%
13	13.148	1833	1838	1842	rBV	8148744	8472688	92.58%	14.212%
14	14.212	2015	2019	2023	rBV	2246596	1961071	21.43%	3.289%
15	15.771	2279	2284	2295	rVB	1056523	1461212	15.97%	2.451%

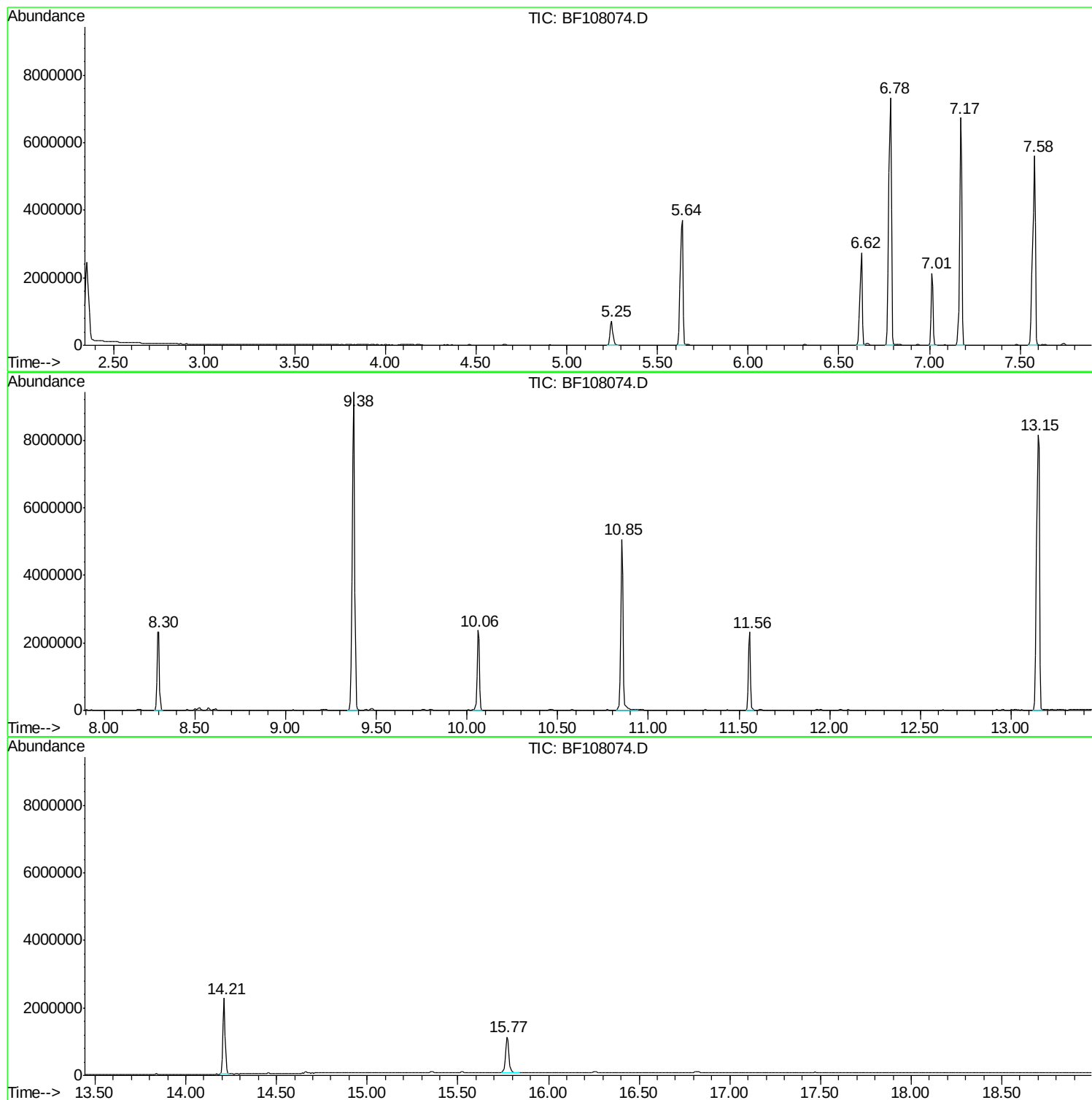
Sum of corrected areas: 59617118

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



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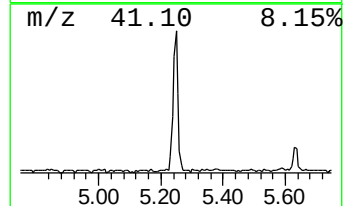
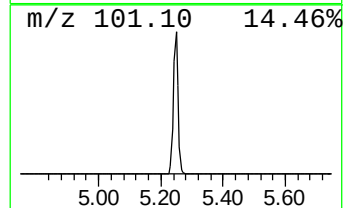
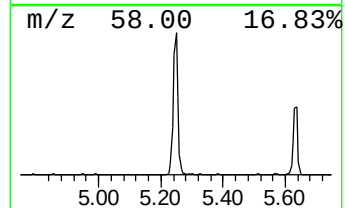
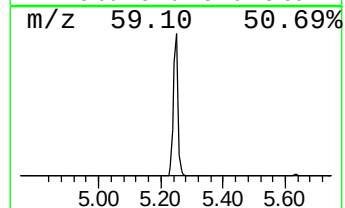
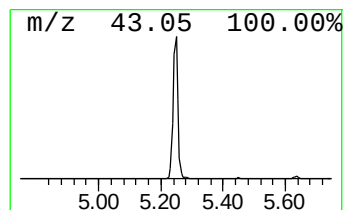
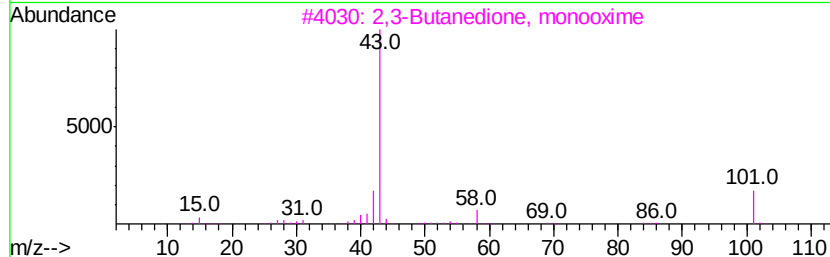
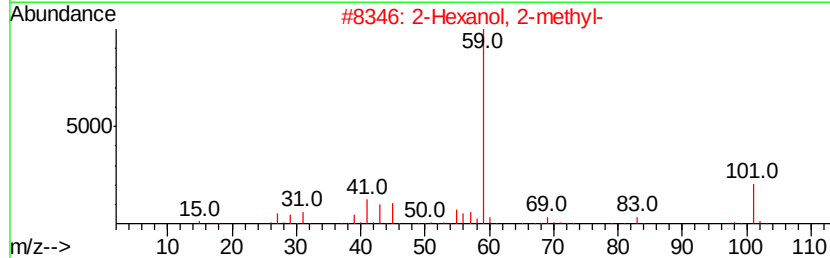
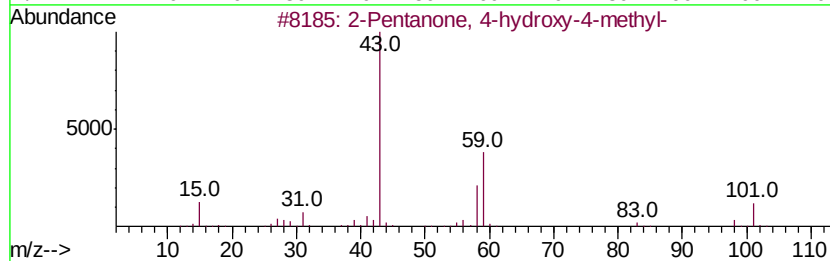
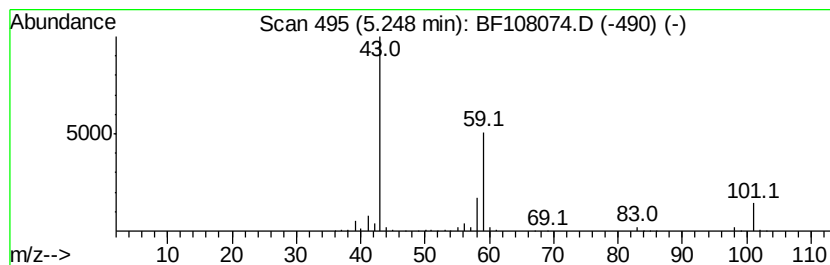
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.25	8.52 ng	763741	1,4-Dichlorobenzene-d4	7.01

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	27
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5			Acetone	58	C3H6O	000067-64-1	9



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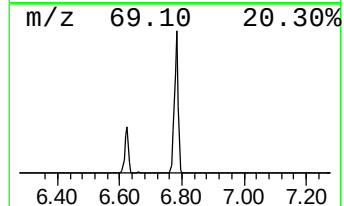
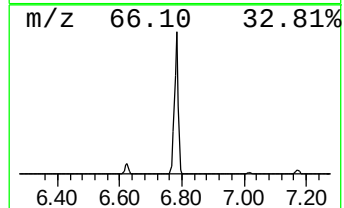
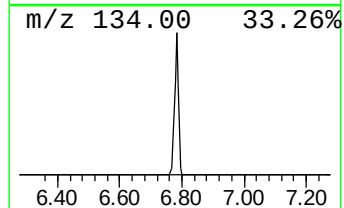
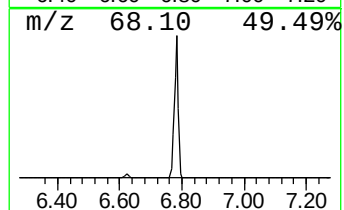
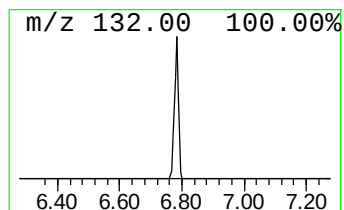
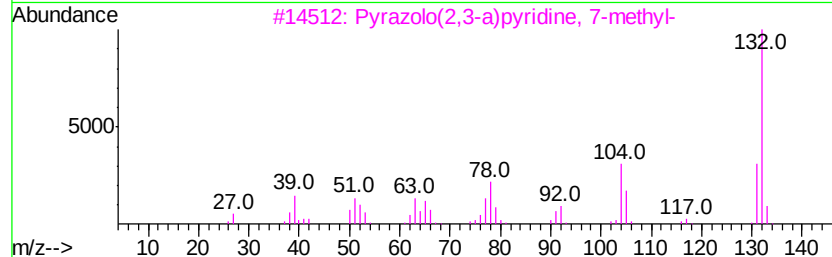
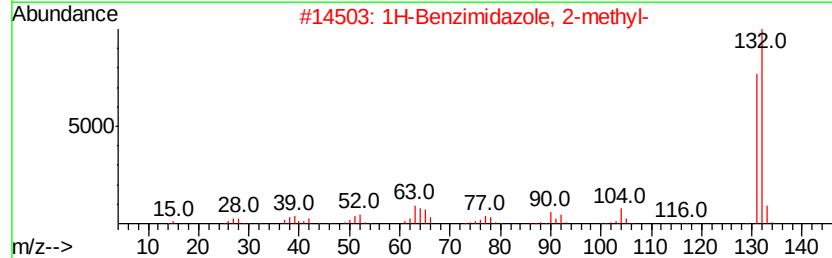
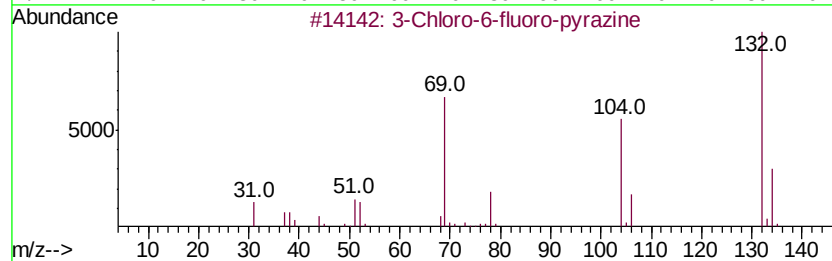
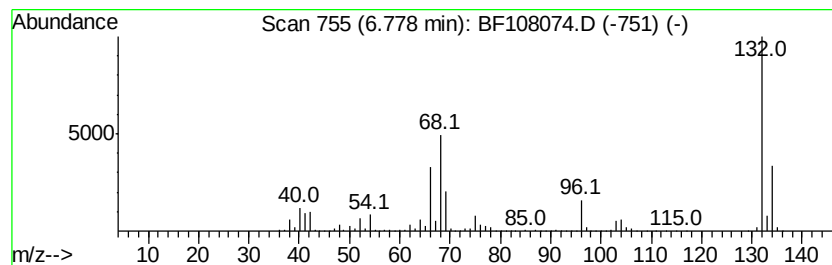
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Peak Number 2 unknown6.78 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.78	75.93 ng	6803690	1,4-Dichlorobenzene-d4	7.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		Pyrazolo(2,3-a)pyridine, 7-methyl-	132	C8H8N2	034760-58-2	12
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.25	8.5	ng	763741	1	7.01	1792210	20.0
unknown6.78	6.78	75.9	ng	6803690	1	7.01	1792210	20.0