

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF062018\
 Data File : BF106625.D
 Acq On : 20 Jun 2018 17:50
 Operator : JU/SJ
 Sample : PB110384BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB110384BL

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-BF061918.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	5.201	482	487	494	rBV	138535	165695	4.53%	0.760%
2	5.601	551	555	565	rBV	1885592	1650458	45.15%	7.573%
3	6.595	720	724	727	rBV	1805372	1652552	45.21%	7.583%
4	6.731	743	747	751	rBV	1849872	1709850	46.77%	7.846%
5	6.954	782	785	789	rBB	652905	542452	14.84%	2.489%
6	7.113	808	812	815	rBV	2086387	1973877	53.99%	9.057%
7	7.519	876	881	884	rBV	1484093	1642683	44.94%	7.538%
8	8.237	999	1003	1007	rBV	710545	698192	19.10%	3.204%
9	9.313	1181	1186	1190	rBV	2701833	2952108	80.75%	13.546%
10	9.995	1298	1302	1309	rVB	852456	836602	22.88%	3.839%
11	10.789	1432	1437	1441	rBV	1594158	1490274	40.77%	6.838%
12	11.489	1552	1556	1560	rBV	1132415	957930	26.20%	4.396%
13	13.083	1821	1827	1830	rBV	3144382	3655682	100.00%	16.774%
14	14.142	2003	2007	2011	rBV	1078705	1056538	28.90%	4.848%
15	15.683	2263	2269	2276	rBV	608834	808449	22.11%	3.710%

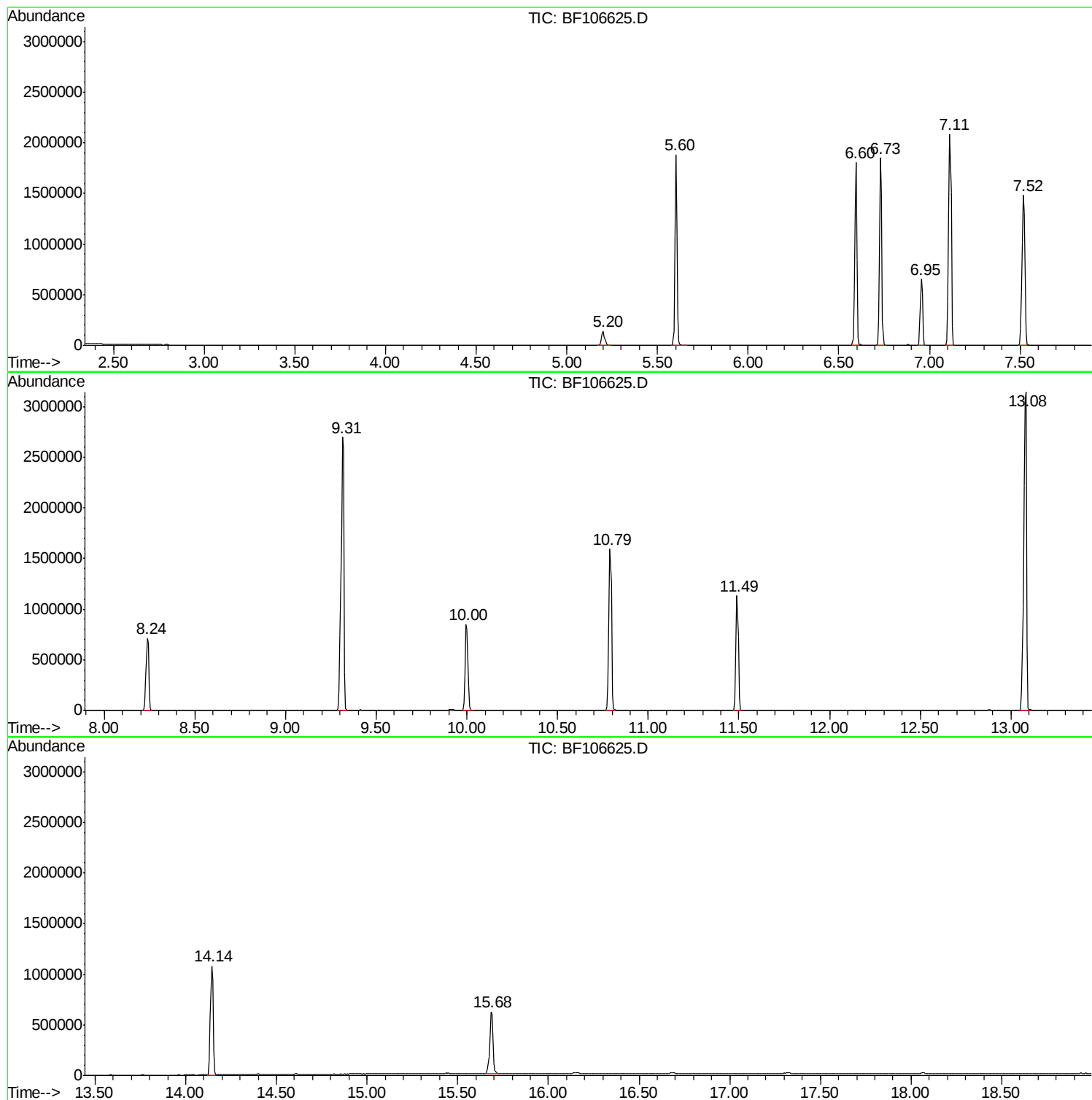
Sum of corrected areas: 21793342

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



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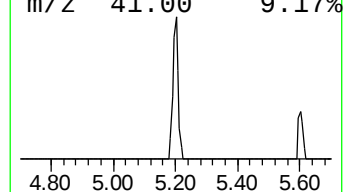
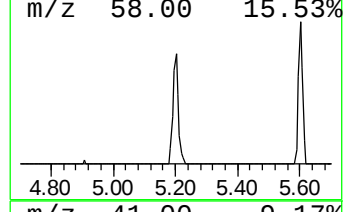
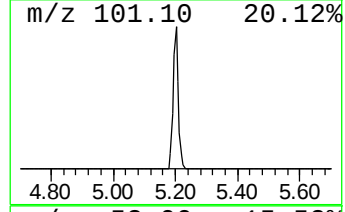
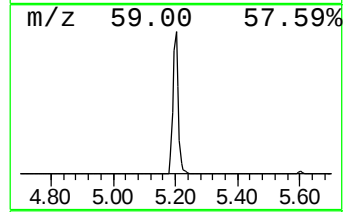
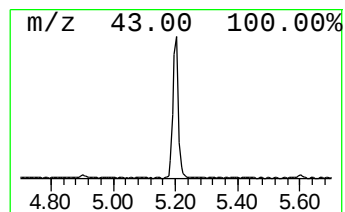
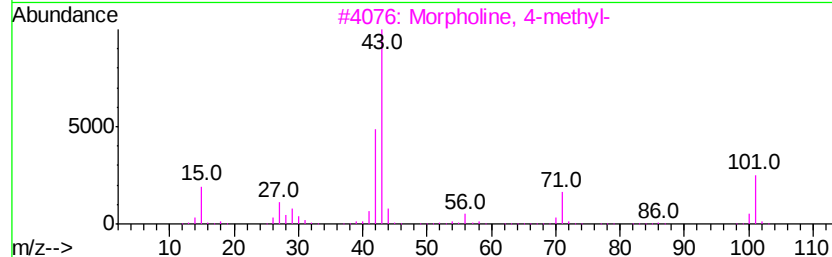
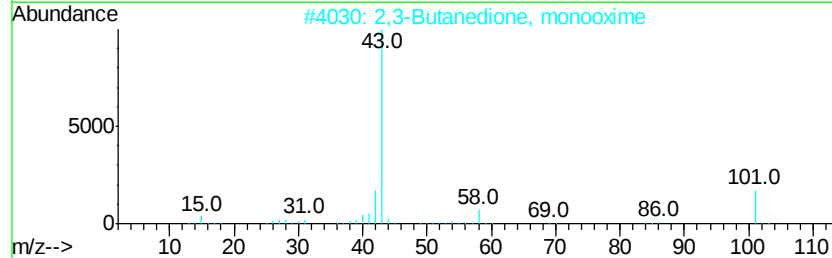
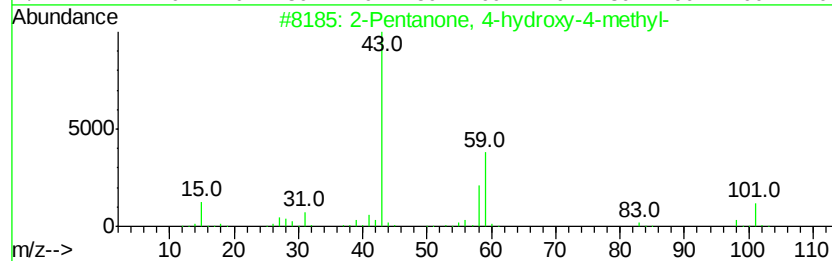
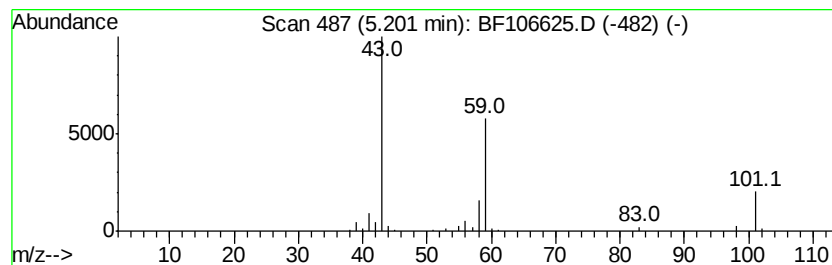
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	6.11 ng	165695	1,4-Dichlorobenzene-d4	6.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
3			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4			3-Octanol	130	C8H18O	000589-98-0	9
5			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9



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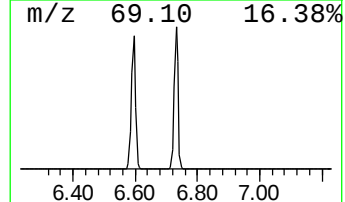
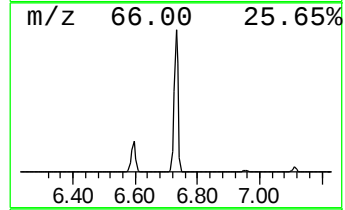
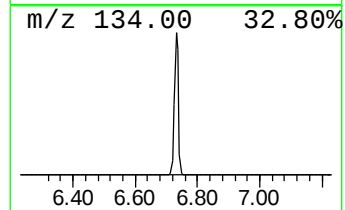
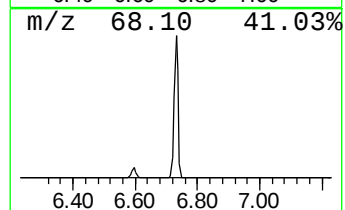
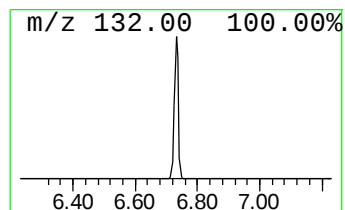
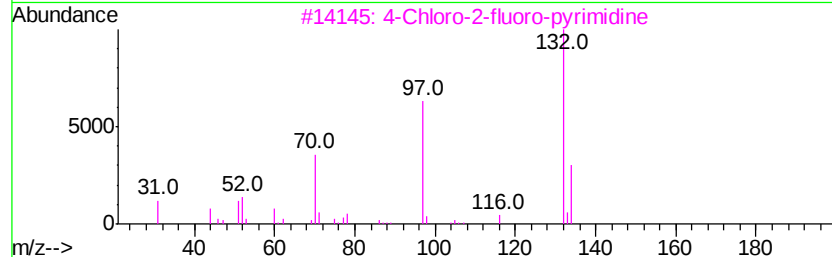
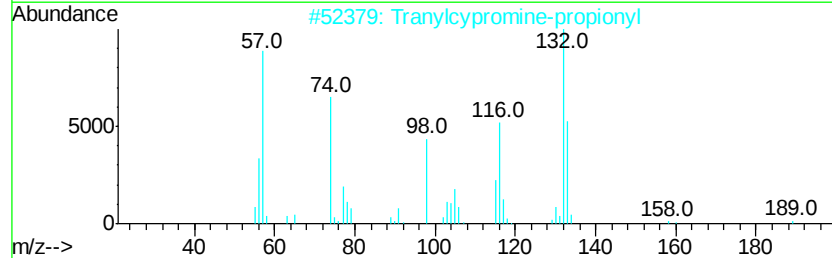
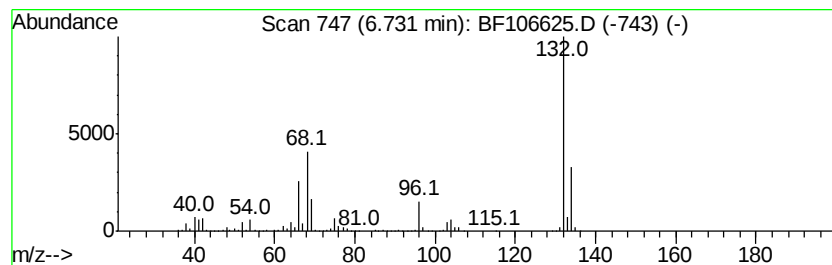
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Peak Number 2 unknown6.73 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.73	63.04 ng	1709850	1,4-Dichlorobenzene-d4	6.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
4		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.20	6.1	ng	165695	1	6.95	542452	20.0
unknown6.73	6.73	63.0	ng	1709850	1	6.95	542452	20.0