

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF070318\
 Data File : BF107114.D
 Acq On : 4 Jul 2018 6:39
 Operator : JU/SJ
 Sample : J3804-05
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-03

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	5.184	480	484	497	rBV	51302	59065	3.58%	0.587%
2	5.596	551	554	571	rBV	552798	597565	36.19%	5.940%
3	6.601	722	725	743	rBV	361594	383661	23.24%	3.814%
4	6.731	743	747	751	rBV	1108869	1085451	65.74%	10.789%
5	6.954	781	785	788	rBV	388396	305948	18.53%	3.041%
6	7.113	807	812	815	rBV	1188085	1091526	66.11%	10.850%
7	7.525	877	882	885	rBV	827410	830074	50.27%	8.251%
8	8.237	999	1003	1021	rBV	350271	338843	20.52%	3.368%
9	9.307	1181	1185	1200	rBV	1380573	1425975	86.36%	14.174%
10	9.701	1249	1252	1258	rBV	93658	90265	5.47%	0.897%
11	9.995	1298	1302	1310	rVB	370993	342058	20.72%	3.400%
12	10.795	1434	1438	1450	rBV	921792	895259	54.22%	8.899%
13	11.489	1553	1556	1562	rBV	412600	380081	23.02%	3.778%
14	13.077	1821	1826	1829	rBV	1709229	1651106	100.00%	16.412%
15	14.142	2004	2007	2013	rBV	365773	343604	20.81%	3.415%
16	15.683	2265	2269	2275	rVB2	174634	240002	14.54%	2.386%

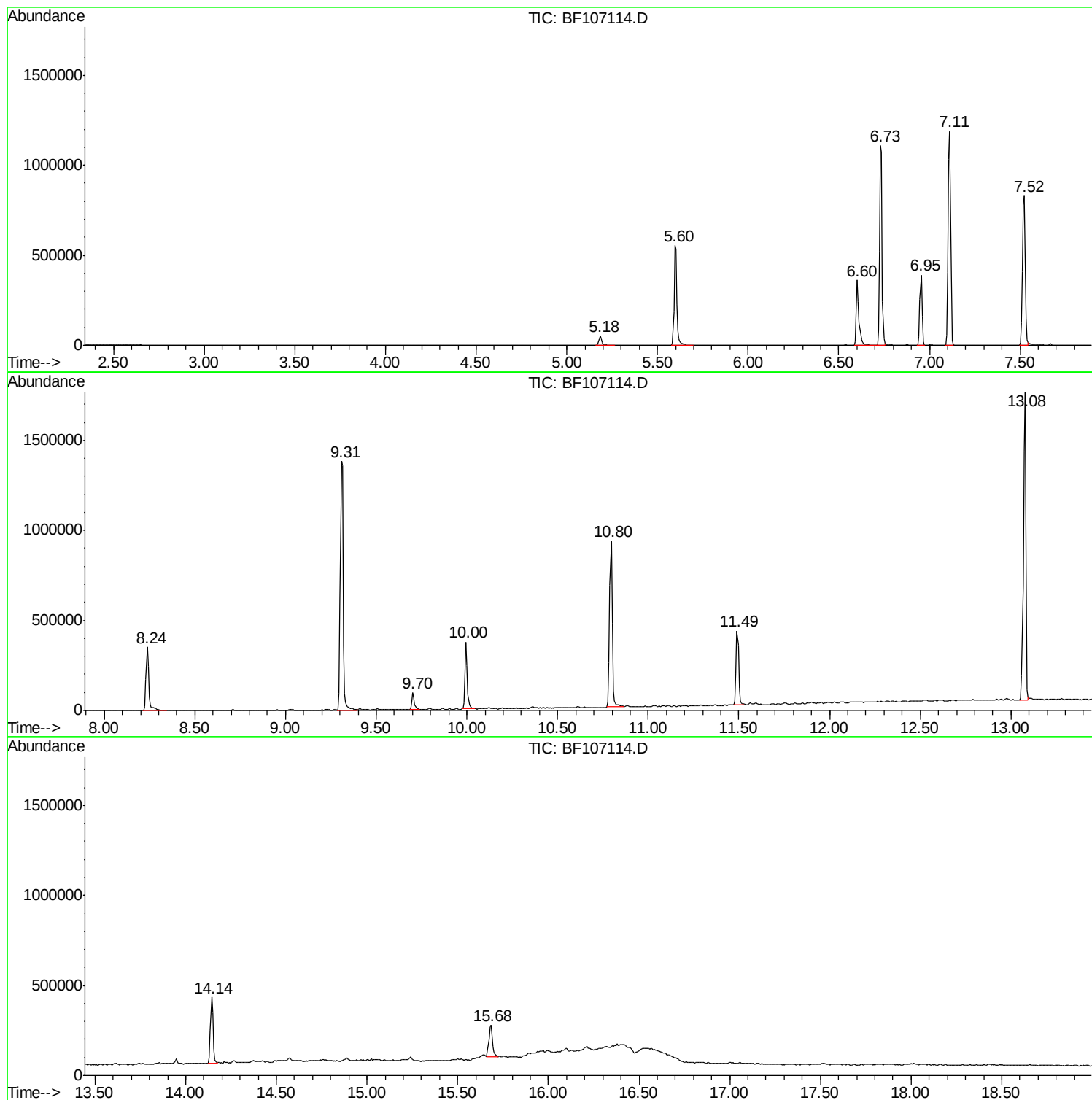
Sum of corrected areas: 10060483

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.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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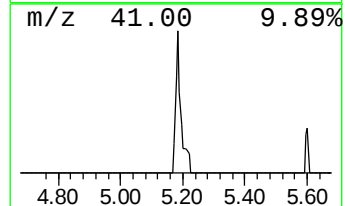
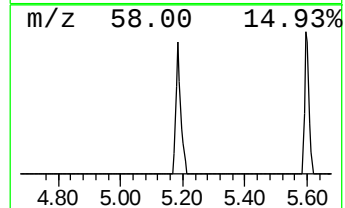
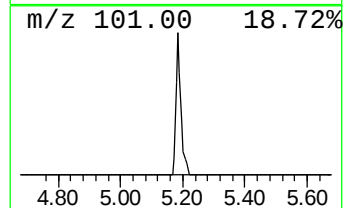
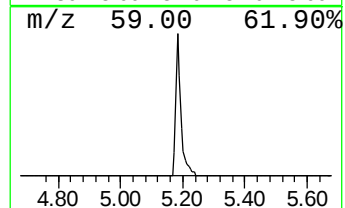
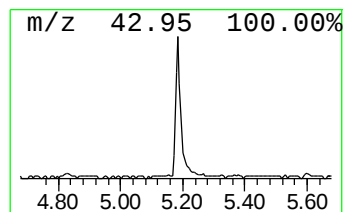
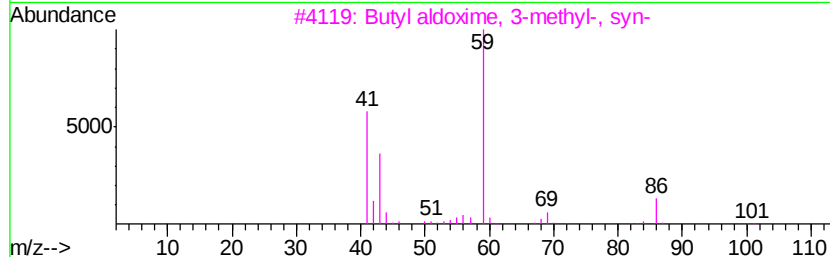
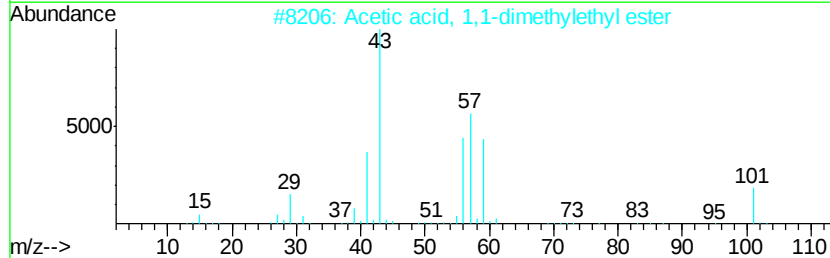
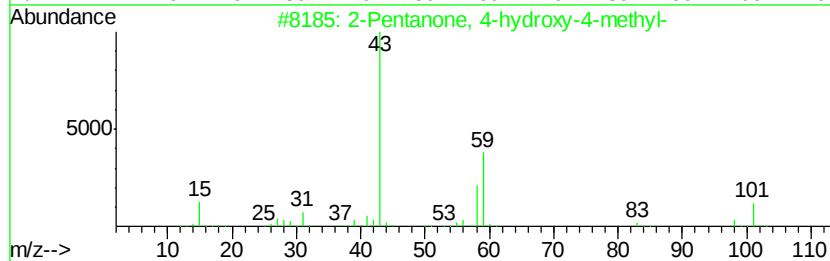
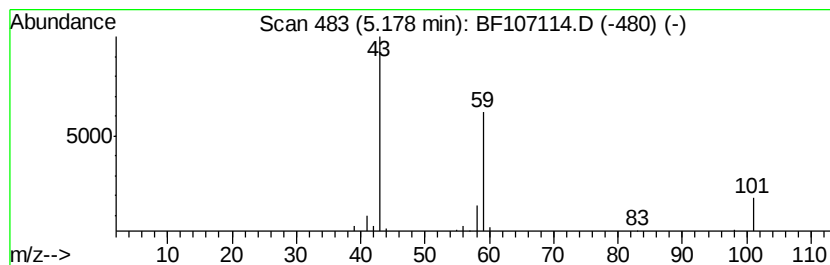
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	3.86 ng	59065	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	42
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	32
4			3-Octanol	130	C8H18O	000589-98-0	28
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



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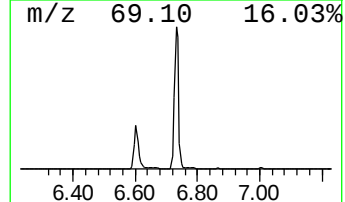
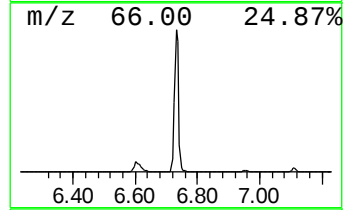
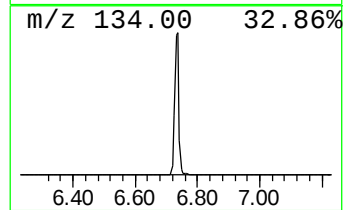
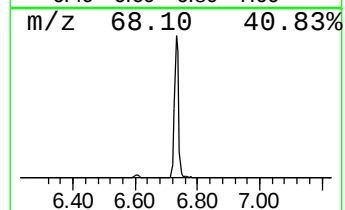
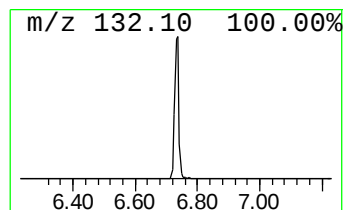
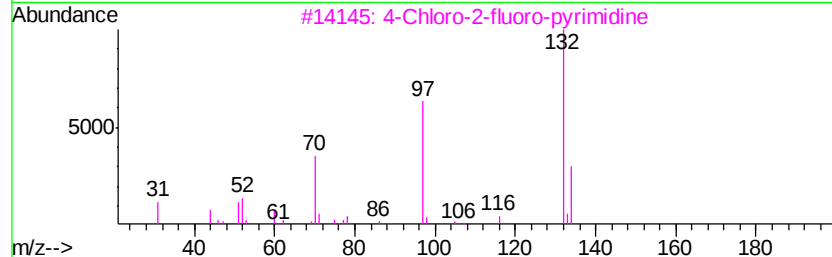
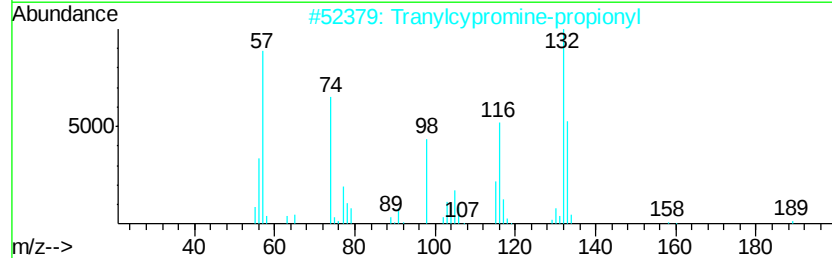
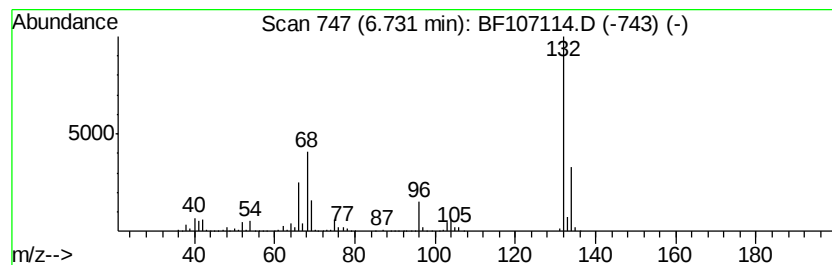
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Peak Number 2 unknown6.73 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.73	70.96 ng	1085450	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		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.18	3.9	ng	59065	1	6.95	305948	20.0
unknown6.73	6.73	71.0	ng	1085450	1	6.95	305948	20.0