

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF122618\
 Data File : BF111711.D
 Acq On : 26 Dec 2018 13:02
 Operator : JU/SJ
 Sample : PB115960BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB115960BL

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-BF121918.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	5.128	511	517	522	rBV	142095	179898	4.45%	0.564%
2	5.540	582	587	590	rBV	3127071	3002342	74.30%	9.420%
3	6.540	752	757	760	rBV	3366252	3120528	77.23%	9.791%
4	6.675	776	780	783	rBV	3809122	3178187	78.65%	9.972%
5	6.898	815	818	822	rBV	1162042	989793	24.49%	3.105%
6	7.057	841	845	849	rBV	3013137	2540534	62.87%	7.971%
7	7.457	908	913	916	rBV	2335110	1987067	49.18%	6.234%
8	8.181	1032	1036	1040	rBV	1444938	1235878	30.58%	3.878%
9	9.257	1214	1219	1222	rBV	4580170	3865520	95.66%	12.128%
10	9.939	1331	1335	1338	rBV	1550049	1424411	35.25%	4.469%
11	10.722	1464	1468	1472	rBV	2525439	2278933	56.40%	7.150%
12	11.422	1583	1587	1591	rBV	1750847	1440074	35.64%	4.518%
13	13.010	1852	1857	1860	rBV	4086225	4040803	100.00%	12.678%
14	14.063	2032	2036	2039	rBV	1497498	1392784	34.47%	4.370%
15	15.545	2283	2288	2292	rBV	938969	1195809	29.59%	3.752%

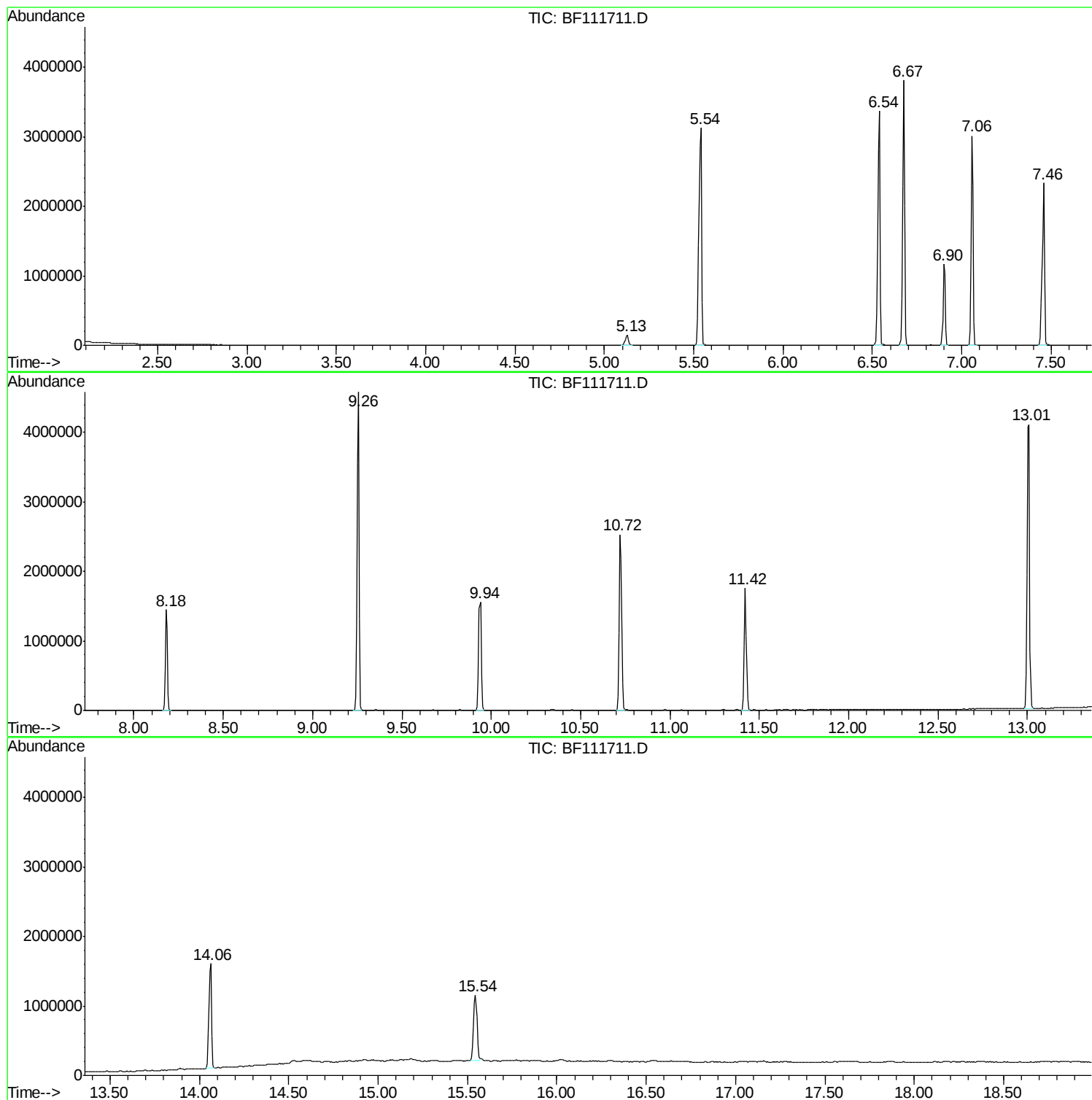
Sum of corrected areas: 31872561

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



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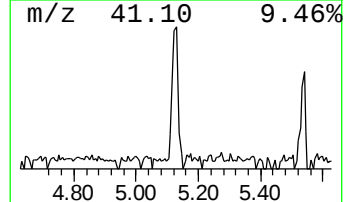
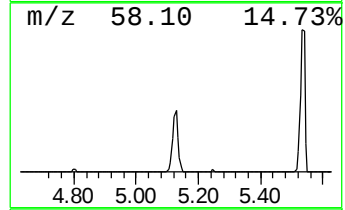
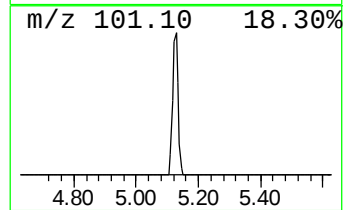
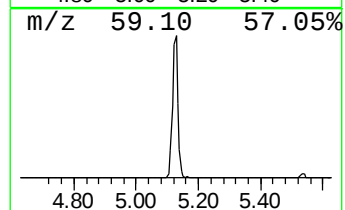
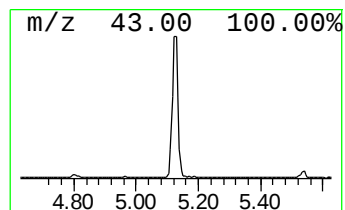
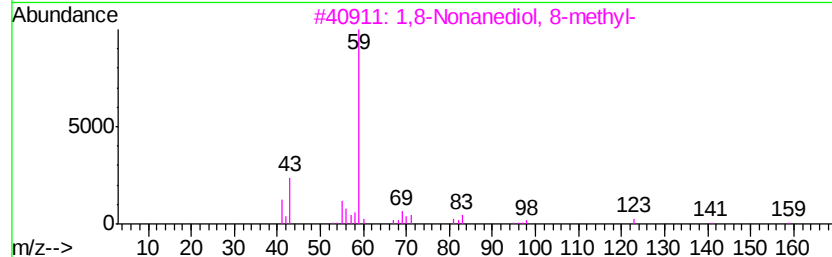
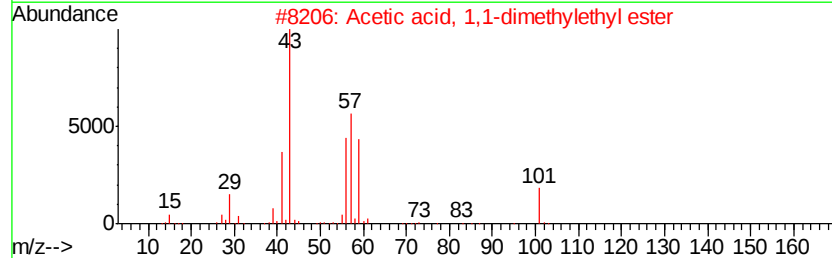
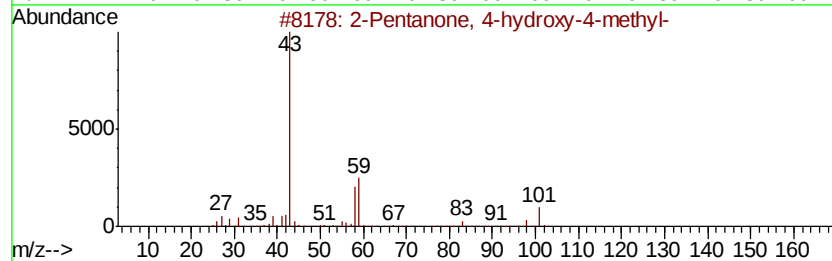
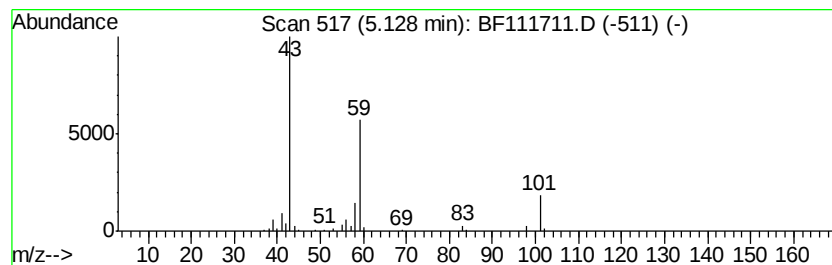
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.13	3.64 ng	179898	1,4-Dichlorobenzene-d4	6.90

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3			1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	23
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	10



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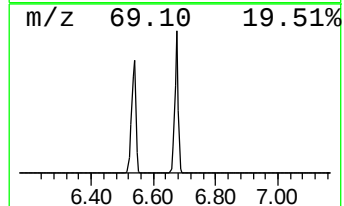
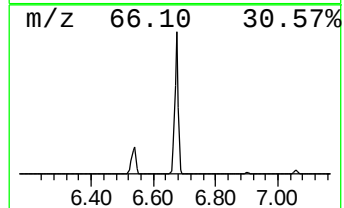
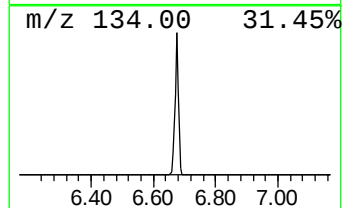
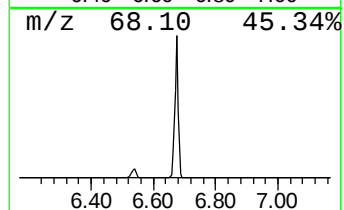
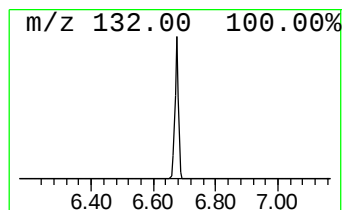
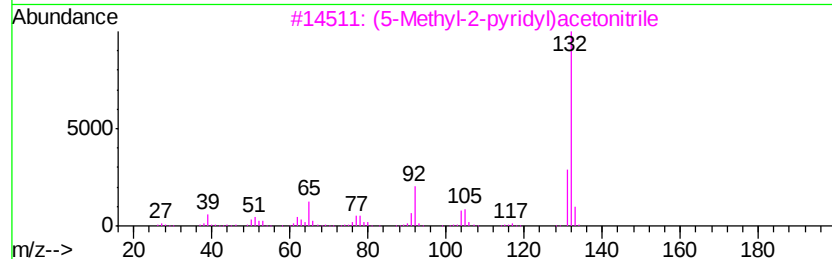
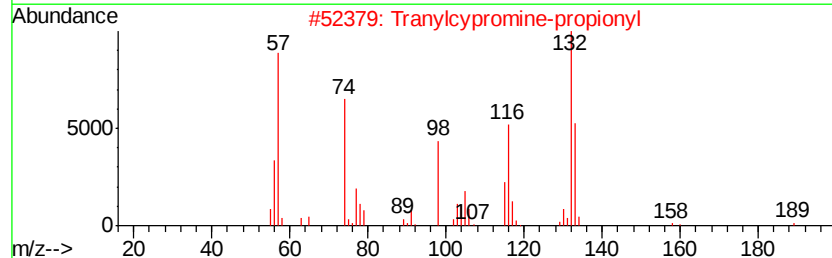
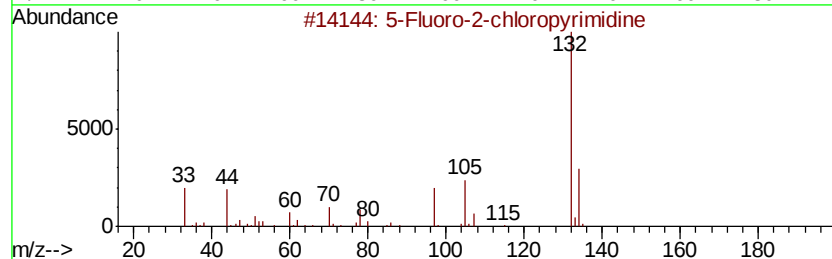
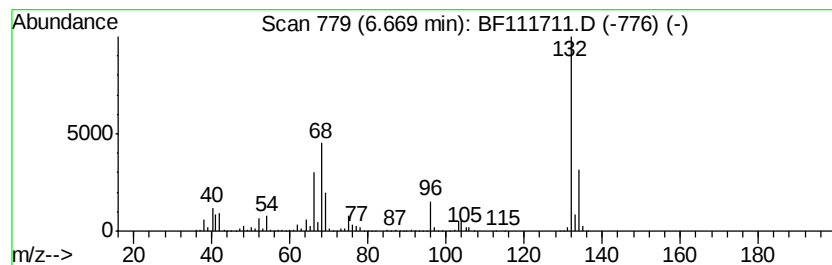
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Peak Number 2 unknown6.67 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.67	64.22 ng	3178190	1,4-Dichlorobenzene-d4	6.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
4		1,2,4-Trifluorobenzene	132	C6H3F3	000367-23-7	9
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.13	3.6	ng	179898	1	6.90	989793	20.0
unknown6.67	6.67	64.2	ng	3178190	1	6.90	989793	20.0