

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF061318\
 Data File : BF106417.D
 Acq On : 14 Jun 2018 3:29
 Operator : JU/SJ
 Sample : PB110157BL
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB110157BL

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-BF061118.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	5.196	482	486	494	rVB	167164	185527	5.64%	0.759%
2	5.613	553	557	570	rVB	1907513	1998403	60.78%	8.177%
3	6.607	722	726	739	rBV	1928503	2017766	61.37%	8.256%
4	6.742	746	749	753	rBV	2265509	1952876	59.40%	7.990%
5	6.966	783	787	790	rBV	1012250	844926	25.70%	3.457%
6	7.119	809	813	817	rBV	2400223	2255554	68.60%	9.229%
7	7.525	877	882	885	rBV	1896930	1881725	57.23%	7.699%
8	8.248	1001	1005	1008	rBV	1272334	1094297	33.28%	4.477%
9	9.319	1183	1187	1191	rBV	3308477	3287757	100.00%	13.452%
10	10.007	1300	1304	1308	rBV	1462091	1247961	37.96%	5.106%
11	10.795	1434	1438	1442	rBV	1571965	1401179	42.62%	5.733%
12	11.495	1553	1557	1561	rBV	1503930	1273705	38.74%	5.211%
13	13.083	1822	1827	1830	rBV	3316258	3093615	94.10%	12.658%
14	14.142	2003	2007	2011	rBV	1135999	952453	28.97%	3.897%
15	15.671	2261	2267	2275	rBV	771072	952847	28.98%	3.899%

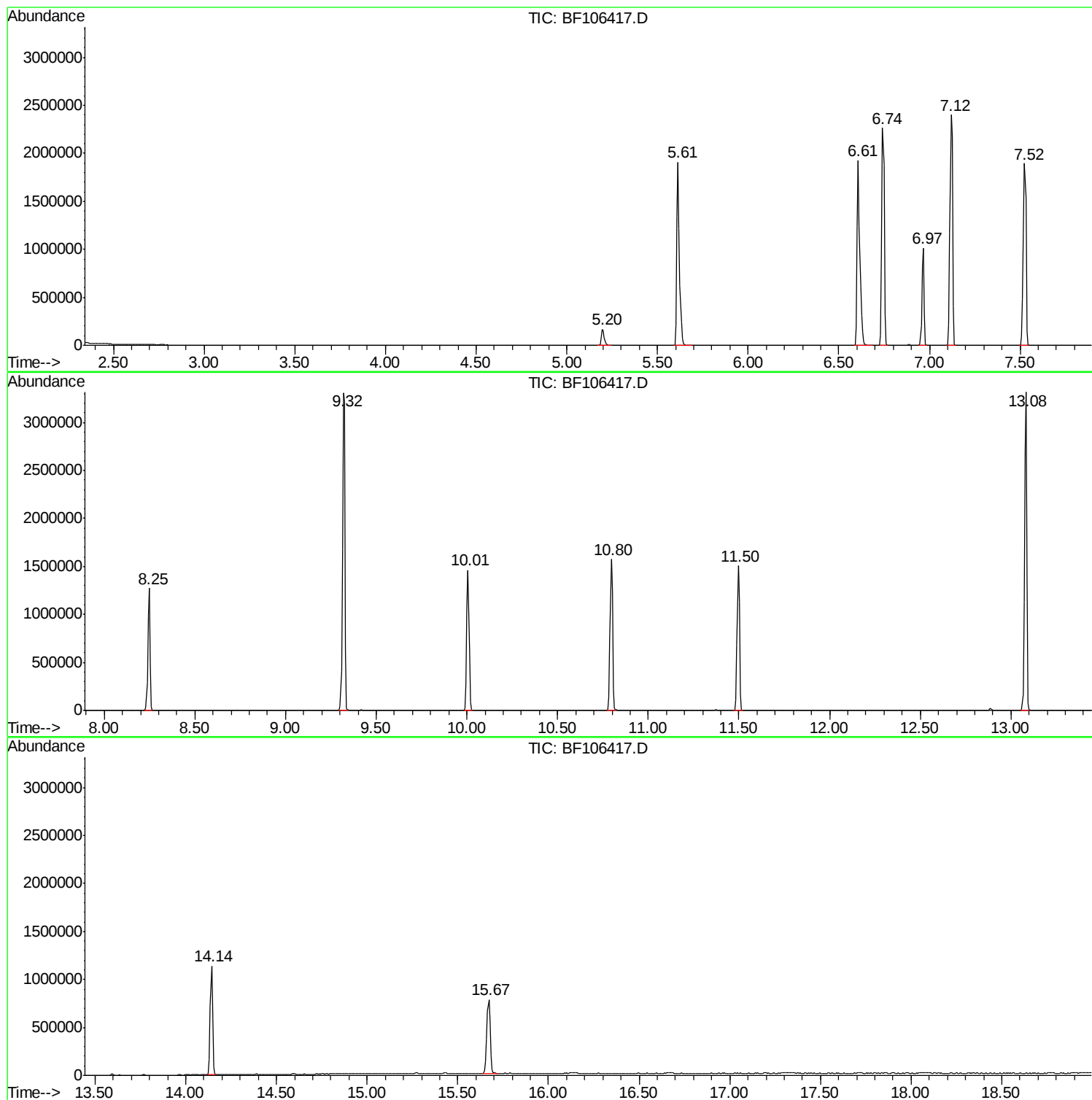
Sum of corrected areas: 24440591

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



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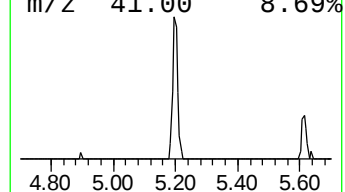
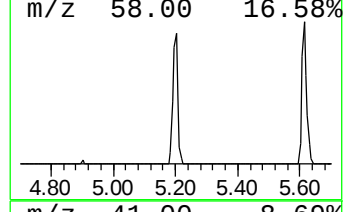
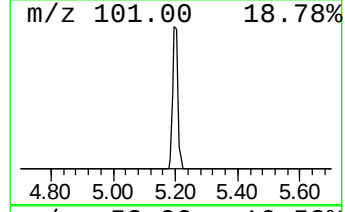
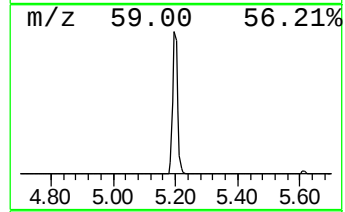
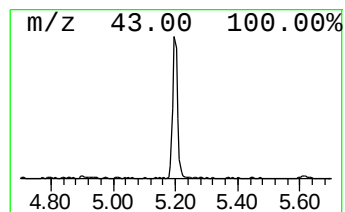
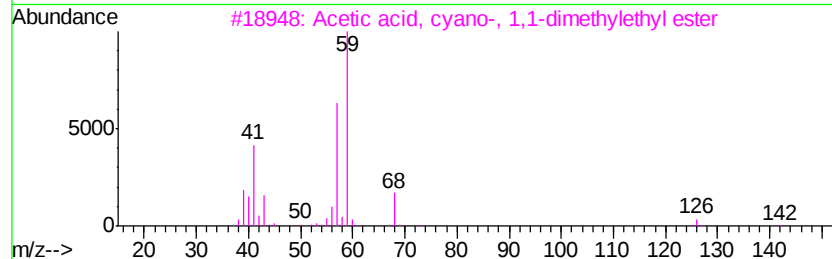
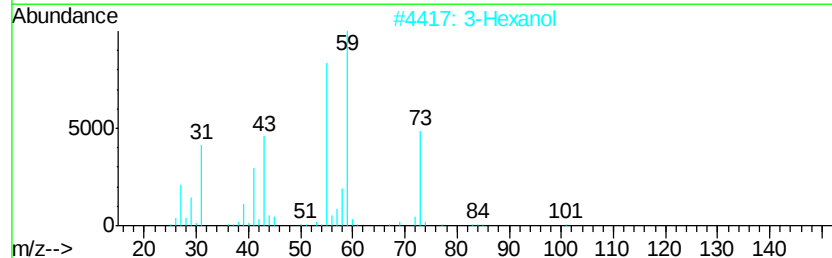
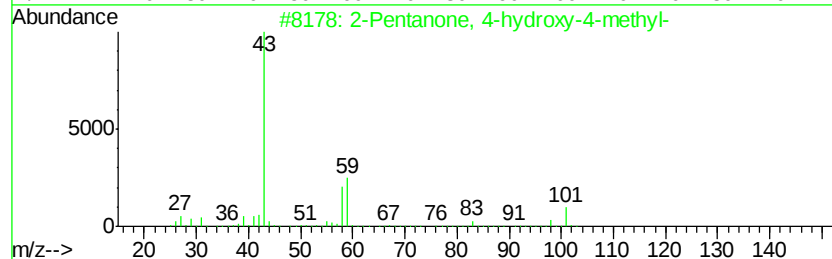
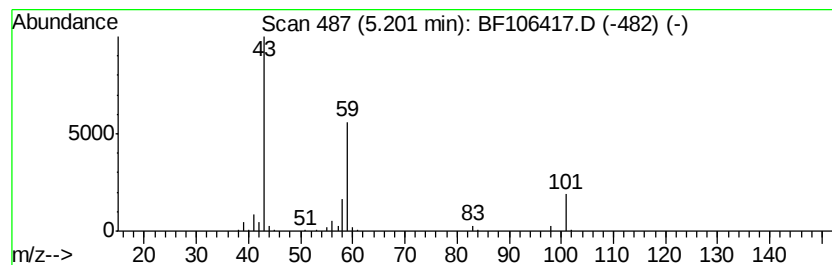
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	4.39 ng	185527	1,4-Dichlorobenzene-d4	6.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			3-Hexanol	102	C6H14O	000623-37-0	33
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
5			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	9



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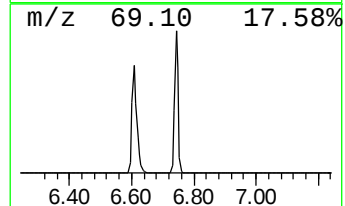
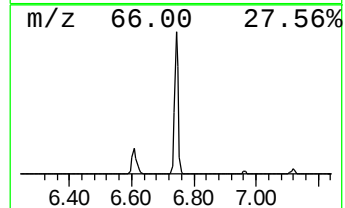
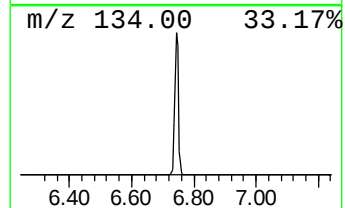
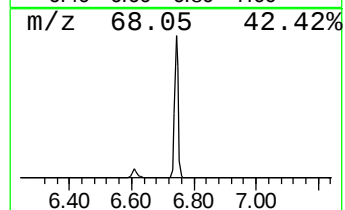
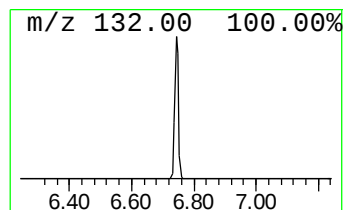
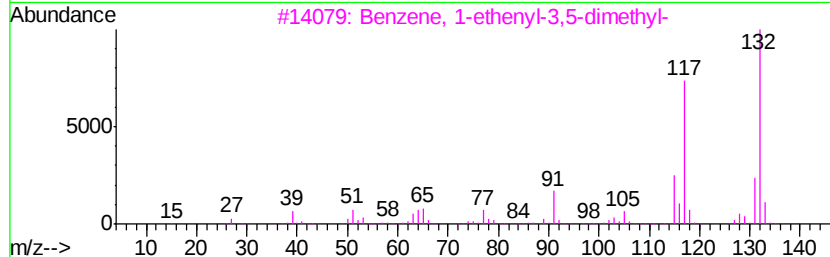
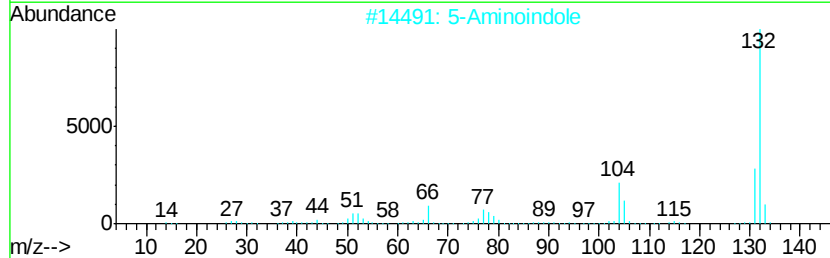
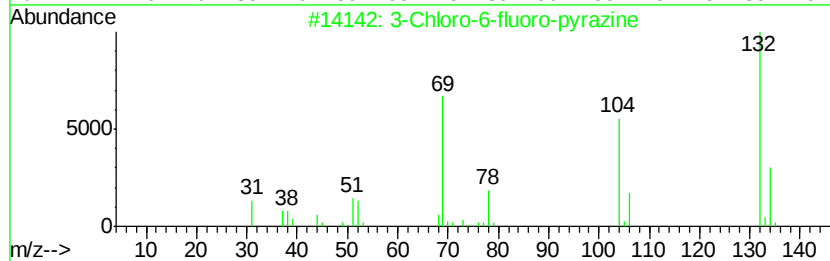
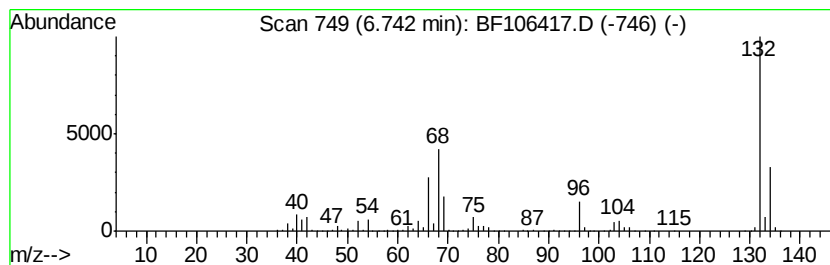
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Peak Number 2 unknown6.74 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	46.23 ng	1952880	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	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...	5.20	4.4	ng	185527	1	6.97	844926	20.0
unknown6.74	6.74	46.2	ng	1952880	1	6.97	844926	20.0