

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF061918\
 Data File : BF106584.D
 Acq On : 19 Jun 2018 16:13
 Operator : JU/SJ
 Sample : PB110378BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB110378BL

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	5.213	483	489	496	rBV	181518	243856	7.35%	1.002%
2	5.607	552	556	561	rBV	2529334	2418210	72.87%	9.941%
3	6.601	721	725	729	rBV	2346326	2448572	73.79%	10.066%
4	6.737	744	748	752	rBV	2476264	2502357	75.41%	10.287%
5	6.960	782	786	789	rBV	746479	599679	18.07%	2.465%
6	7.119	808	813	816	rBV	2118139	2056920	61.99%	8.456%
7	7.525	876	882	885	rBV	1620568	1726140	52.02%	7.096%
8	8.242	999	1004	1007	rBV	848420	764276	23.03%	3.142%
9	9.319	1181	1187	1190	rBV	3031501	3082495	92.89%	12.672%
10	10.001	1298	1303	1307	rBV	1005698	877845	26.45%	3.609%
11	10.795	1433	1438	1441	rBV	1964005	1969628	59.36%	8.097%
12	11.495	1553	1557	1560	rBV	1079006	943205	28.42%	3.878%
13	13.083	1822	1827	1830	rBV	3133059	3318352	100.00%	13.642%
14	14.148	2004	2008	2012	rBV	787572	747747	22.53%	3.074%
15	15.701	2266	2272	2281	rBV	465164	625716	18.86%	2.572%

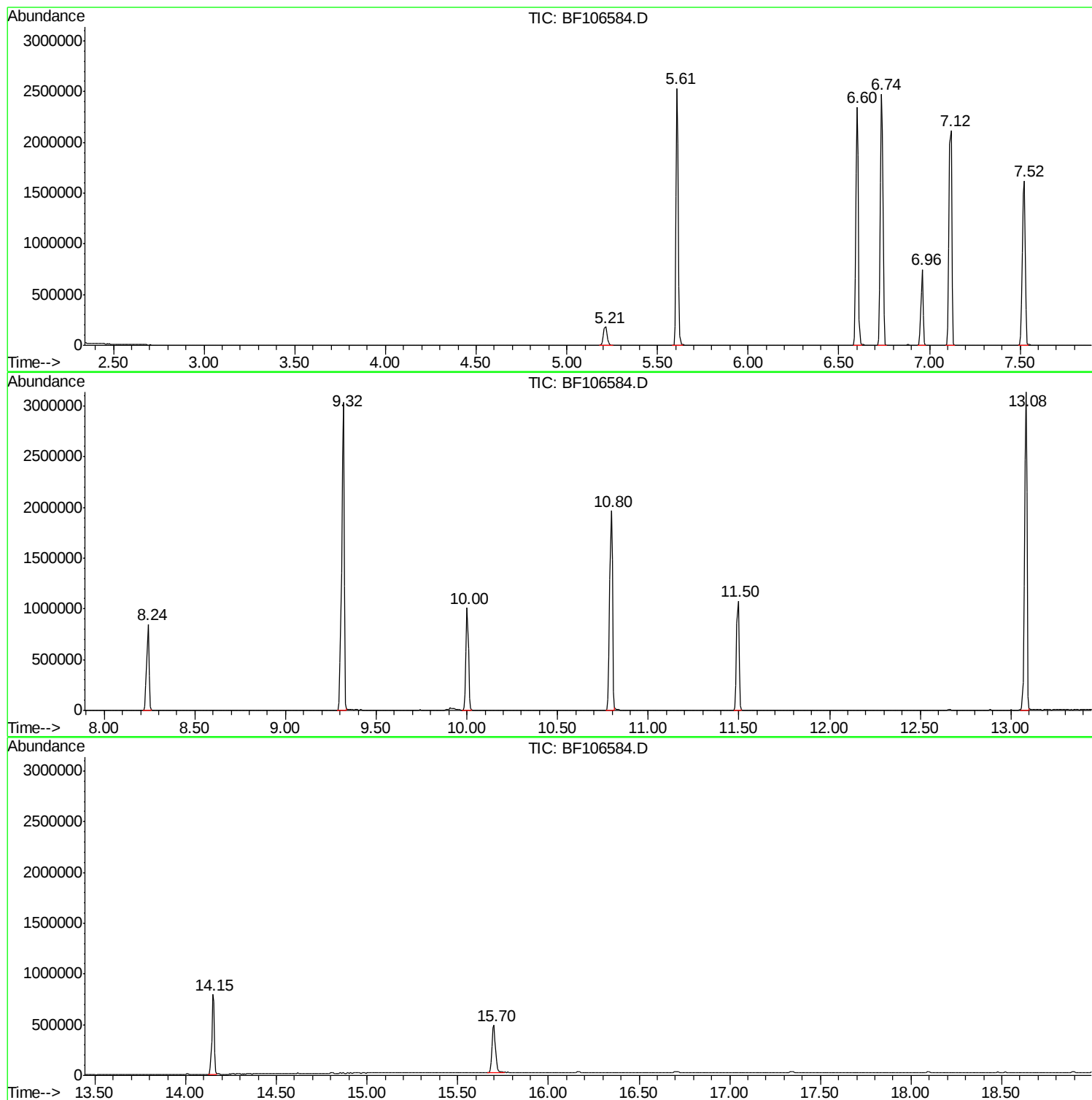
Sum of corrected areas: 24324998

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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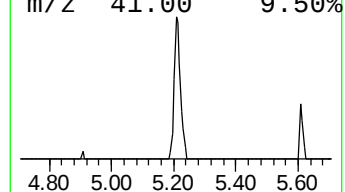
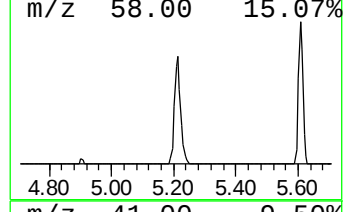
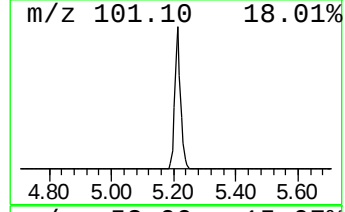
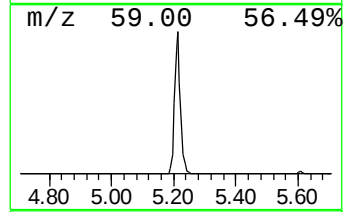
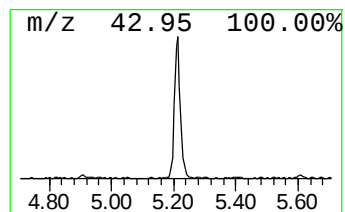
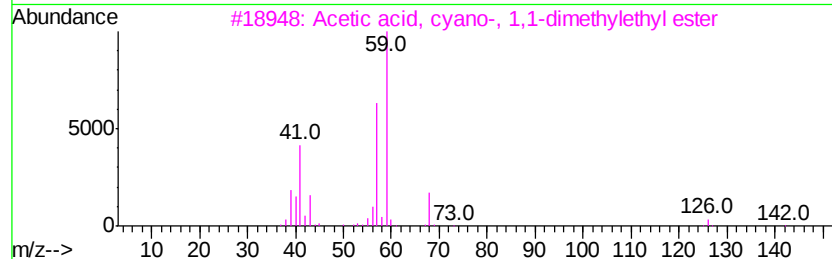
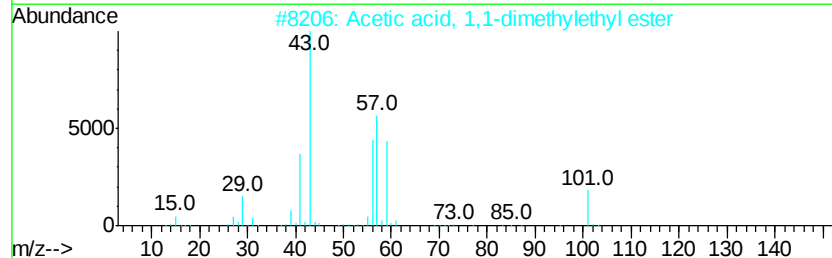
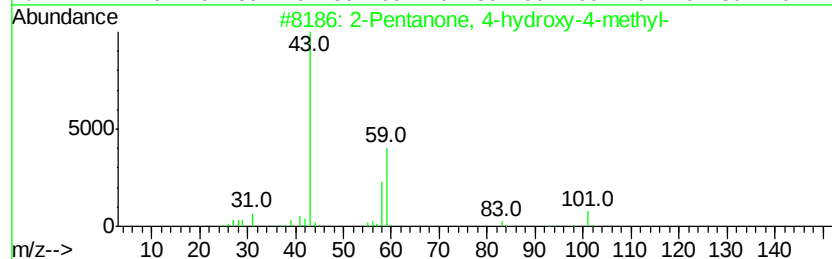
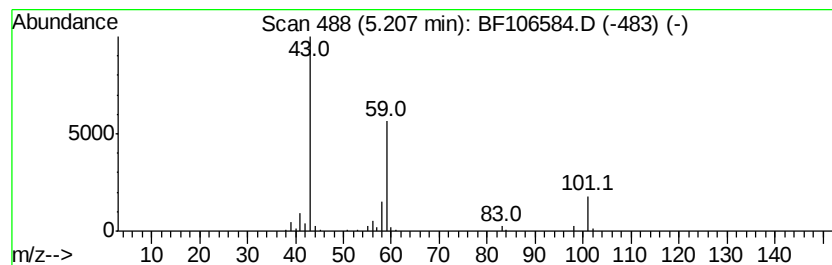
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.21	8.13 ng	243856	1,4-Dichlorobenzene-d4	6.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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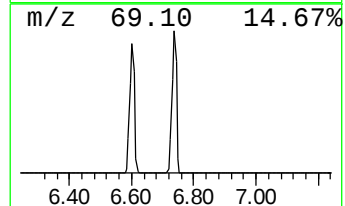
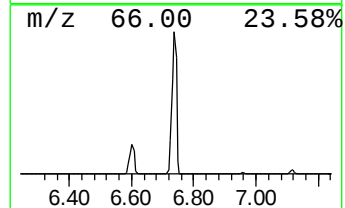
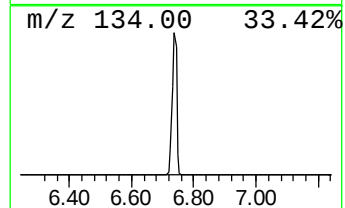
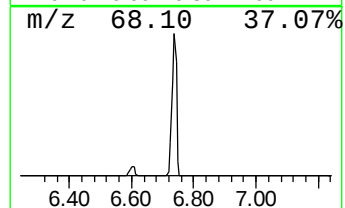
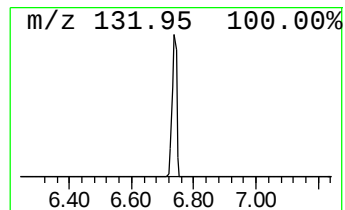
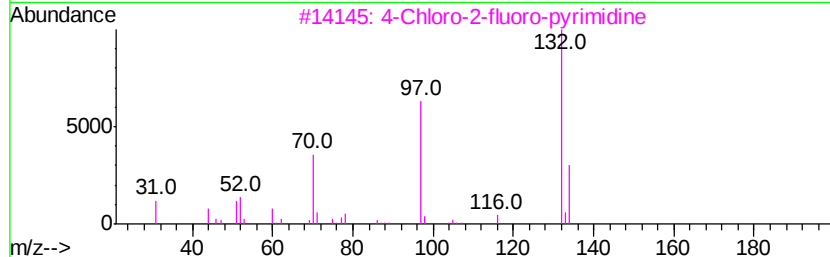
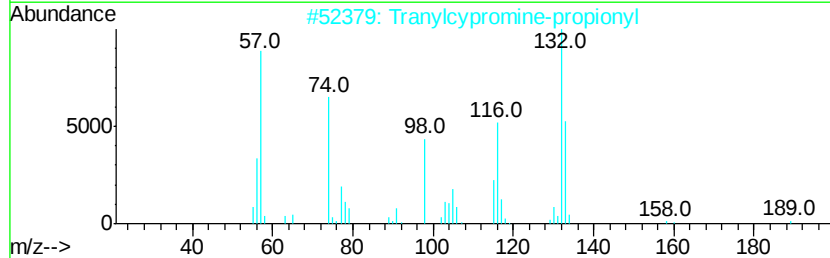
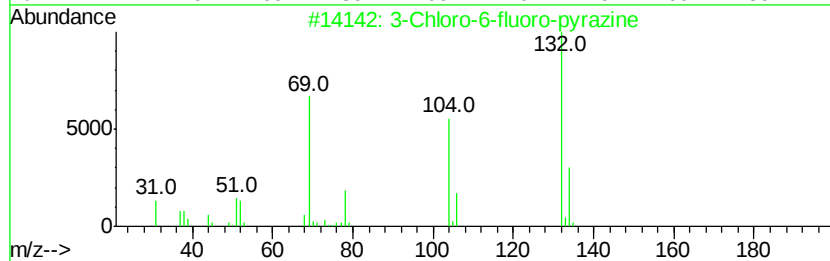
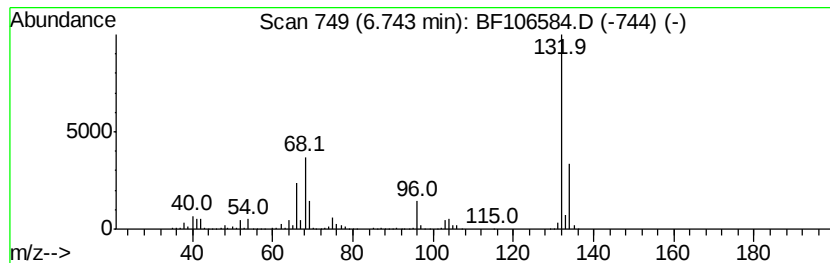
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Peak Number 2 unknown6.74 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	83.46 ng	2502360	1,4-Dichlorobenzene-d4	6.96

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		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	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...	5.21	8.1	ng	243856	1	6.96	599679	20.0
unknown6.74	6.74	83.5	ng	2502360	1	6.96	599679	20.0