

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF102618\
 Data File : BF110248.D
 Acq On : 27 Oct 2018 8:43
 Operator : JU/SJ
 Sample : PB114167BL
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB114167BL

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-BF102318.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	2.369	44	48	57	rVB	26306	43465	1.44%	0.232%
2	5.210	526	531	542	rBV	266244	338854	11.26%	1.809%
3	5.593	592	596	600	rBV	1817489	1766223	58.69%	9.427%
4	6.587	760	765	768	rBV	1896832	1714966	56.98%	9.153%
5	6.610	768	769	783	rVB	68249	67297	2.24%	0.359%
6	6.734	786	790	793	rBV	2182807	1841639	61.19%	9.829%
7	6.963	826	829	833	rBV	661669	565606	18.79%	3.019%
8	7.122	852	856	860	rBV	2455406	2182058	72.50%	11.646%
9	7.528	920	925	928	rBV	1622216	1638320	54.44%	8.744%
10	8.251	1043	1048	1064	rBV	699194	741930	24.65%	3.960%
11	9.322	1225	1230	1245	rBV	3155135	3009596	100.00%	16.063%
12	10.010	1343	1347	1359	rVB	760683	709474	23.57%	3.787%
13	10.798	1477	1481	1500	rBV	787318	786007	26.12%	4.195%
14	11.498	1597	1600	1608	rBV	592733	583086	19.37%	3.112%
15	13.086	1866	1870	1873	rBV	2336657	1963903	65.25%	10.482%
16	14.151	2047	2051	2059	rBV	464409	458167	15.22%	2.445%
17	15.674	2305	2310	2321	rBV2	230378	325751	10.82%	1.739%

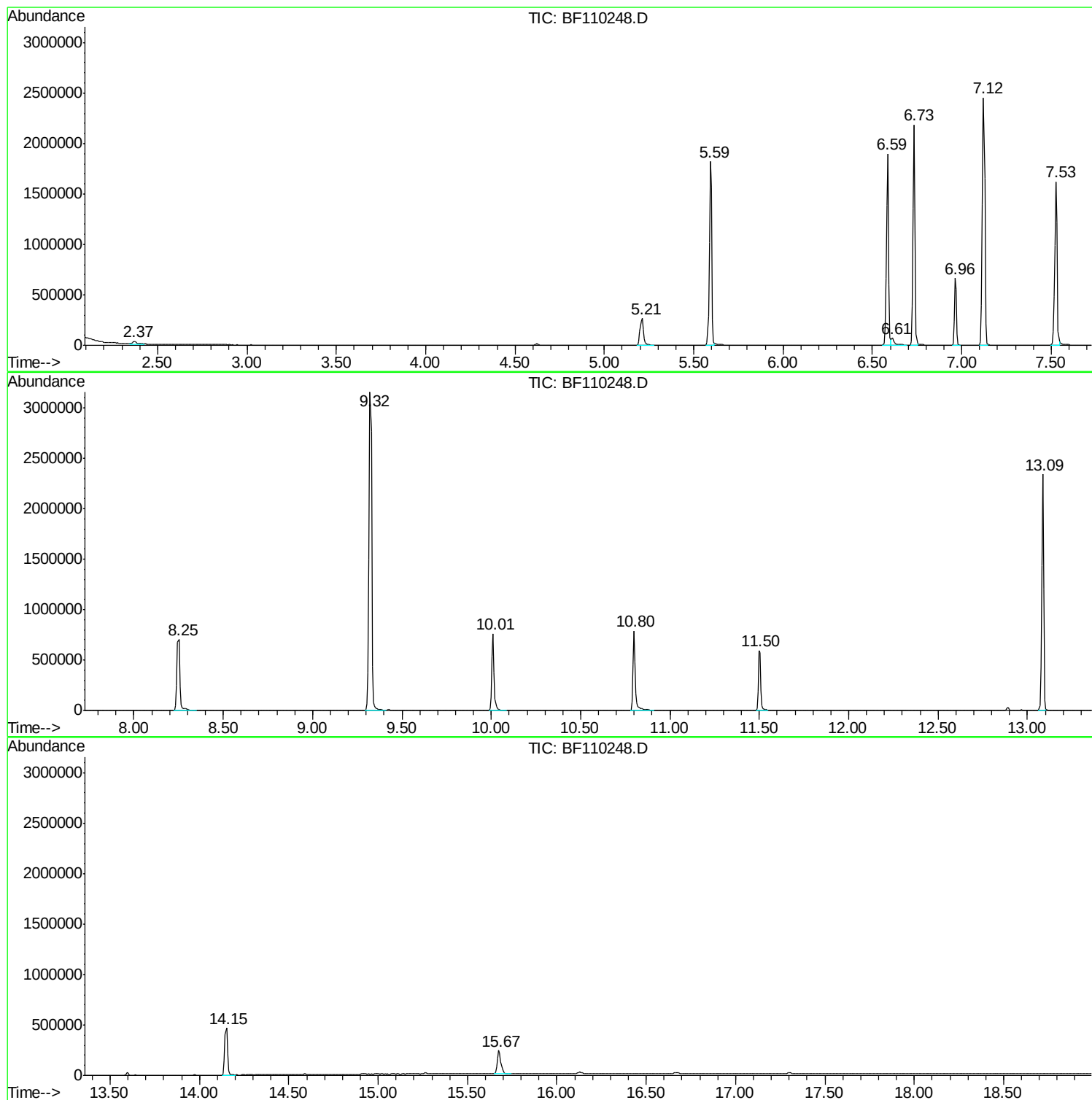
Sum of corrected areas: 18736342

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



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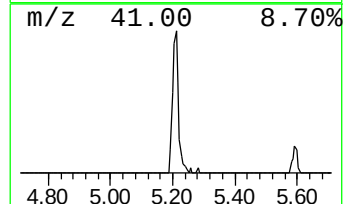
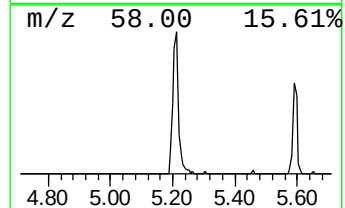
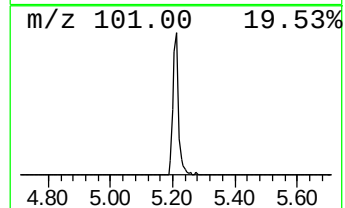
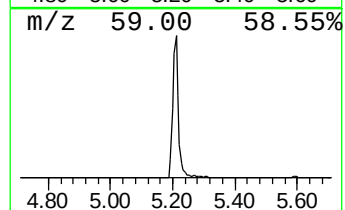
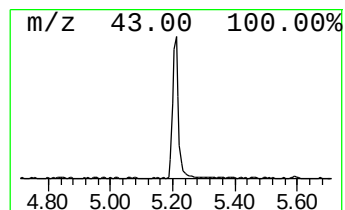
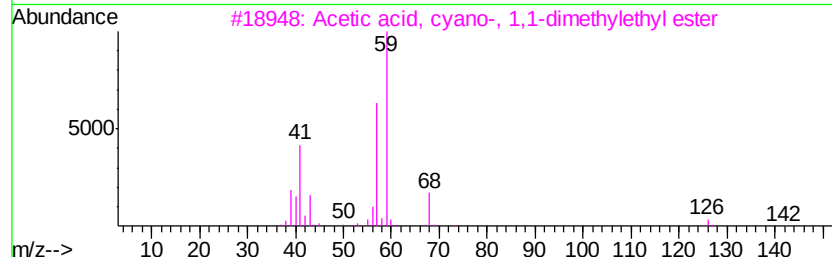
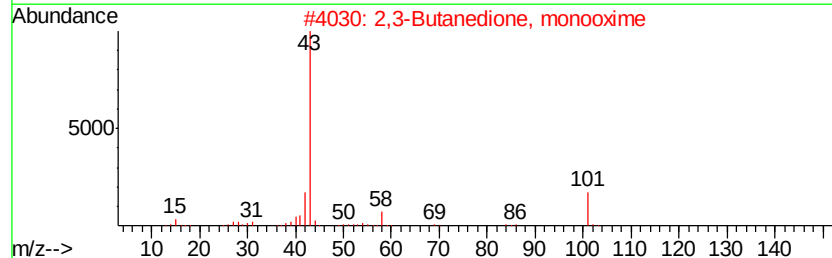
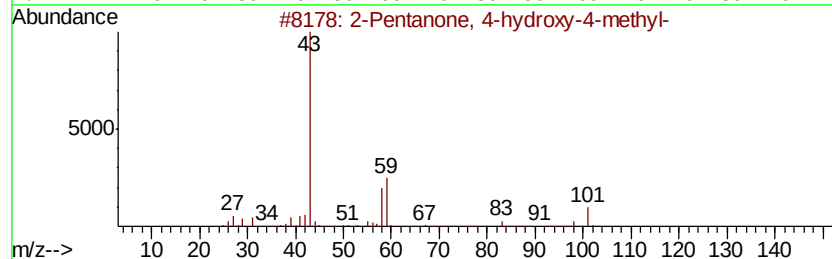
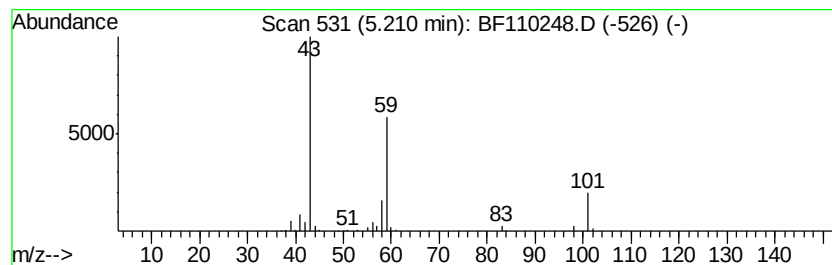
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.21	11.98 ng	338854	1,4-Dichlorobenzene-d4	6.96

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	25
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9



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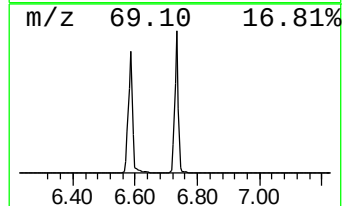
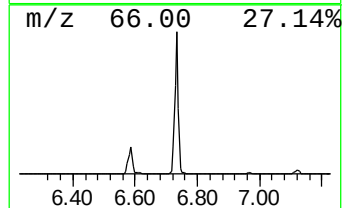
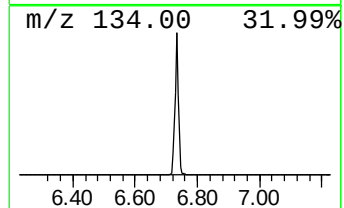
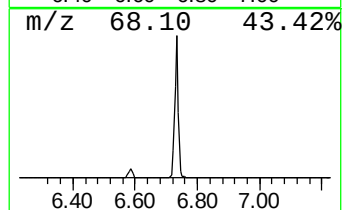
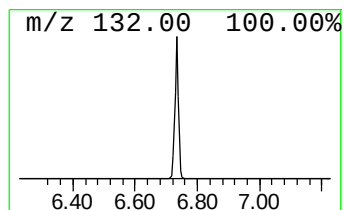
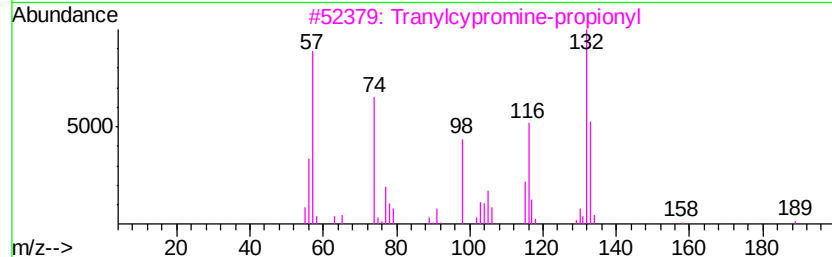
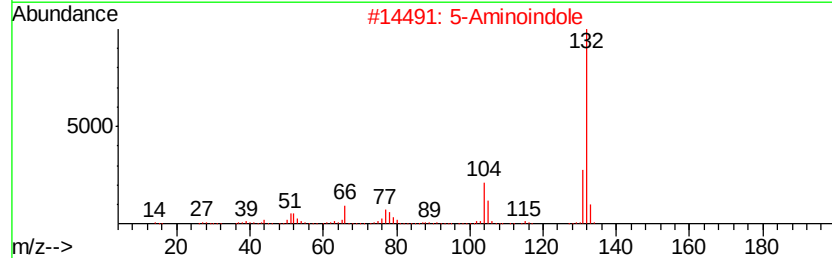
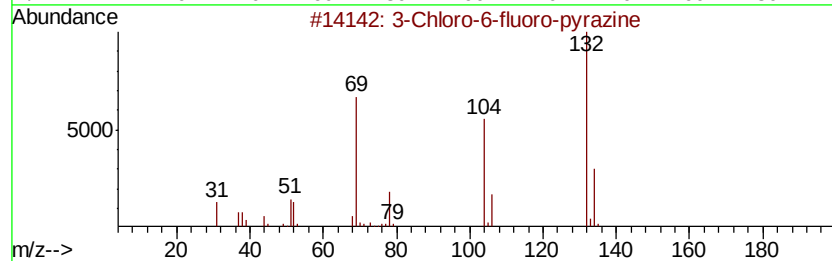
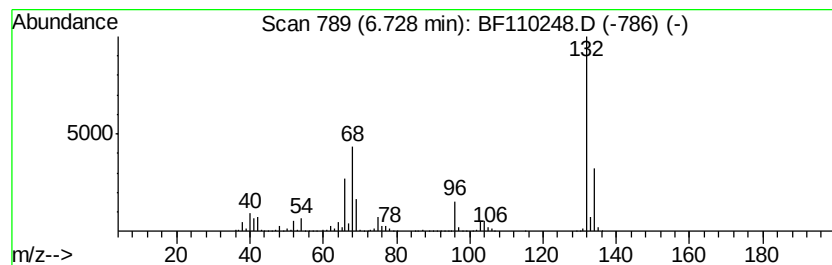
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Peak Number 2 unknown6.73 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.73	65.12 ng	1841640	1,4-Dichlorobenzene-d4	6.96

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10
4			2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5			4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.21	12.0	ng	338854	1	6.96	565606	20.0
unknown6.73	6.73	65.1	ng	1841640	1	6.96	565606	20.0