

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF090718\
 Data File : BF108875.D
 Acq On : 7 Sep 2018 21:48
 Operator : JU/SJ
 Sample : PB112683BL
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB112683BL

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 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-BF090518.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	4.942	438	443	452	rVB	448975	611373	9.65%	1.449%
2	5.354	508	513	516	rBV	3910122	3919799	61.86%	9.292%
3	6.372	680	686	689	rBV	4100862	4013484	63.34%	9.514%
4	6.507	704	709	712	rBV	4209193	4049510	63.91%	9.599%
5	6.736	744	748	751	rBV	1568660	1265708	19.98%	3.000%
6	6.895	770	775	778	rBV	5146155	4542289	71.69%	10.767%
7	7.295	837	843	846	rBV	3589040	3752373	59.22%	8.895%
8	8.019	961	966	969	rBV	1676449	1564329	24.69%	3.708%
9	9.095	1143	1149	1152	rBV	6648572	6336154	100.00%	15.020%
10	9.766	1259	1263	1267	rBV	1997939	1636205	25.82%	3.879%
11	10.548	1392	1396	1408	rBV	2400594	2119172	33.45%	5.023%
12	11.242	1510	1514	1517	rBV	1746438	1369584	21.62%	3.247%
13	12.824	1779	1783	1787	rBV	4626510	4317224	68.14%	10.234%
14	13.865	1956	1960	1964	rBV	1487956	1217291	19.21%	2.886%
15	15.271	2194	2199	2204	rBV	1105111	1326658	20.94%	3.145%
16	15.742	2274	2279	2283	rBV2	47379	72000	1.14%	0.171%
17	16.206	2355	2358	2364	rVB2	46301	72207	1.14%	0.171%

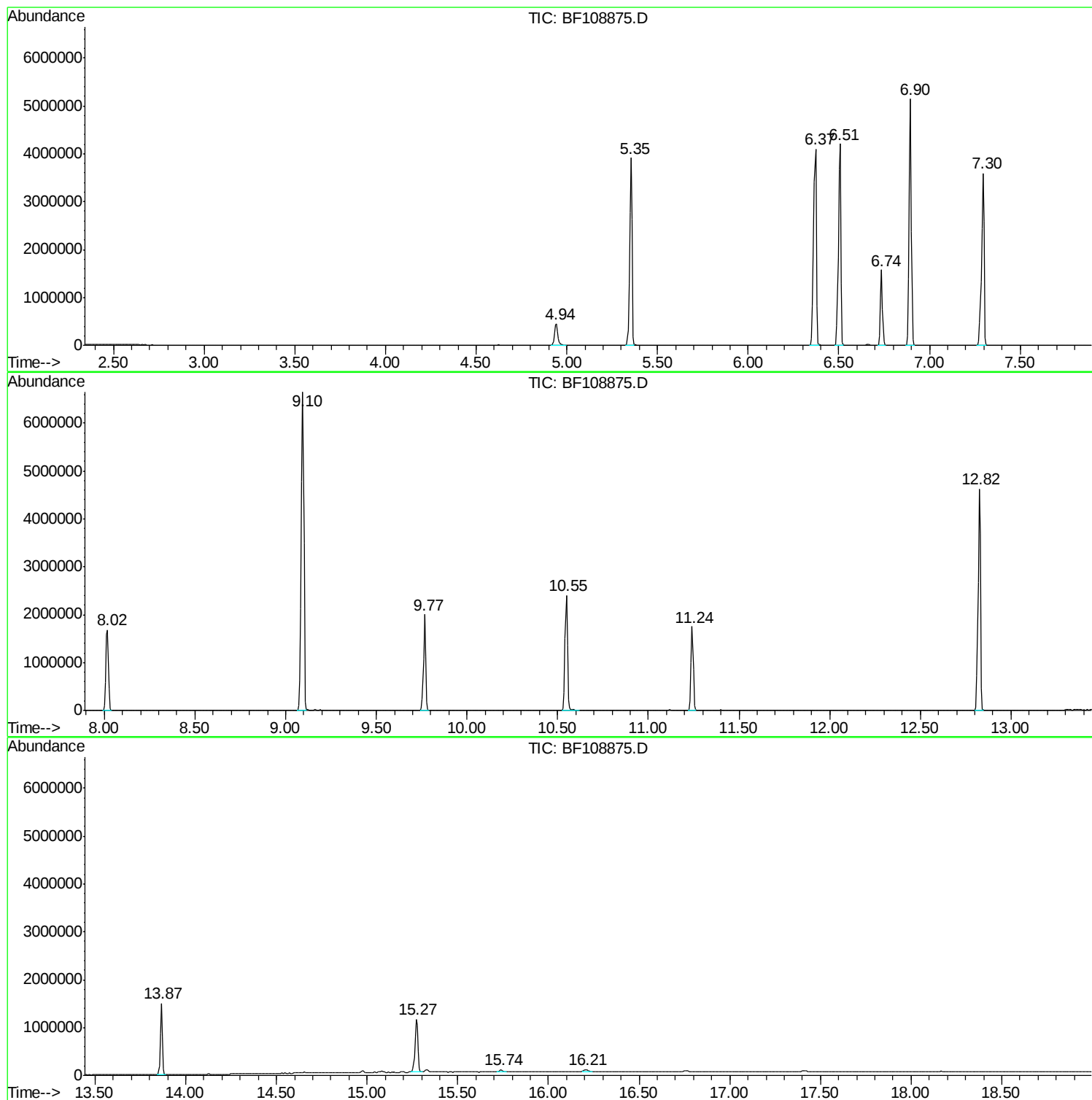
Sum of corrected areas: 42185360

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF090518.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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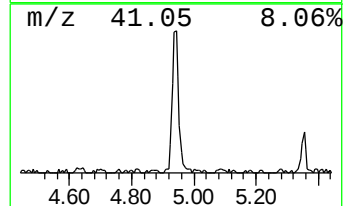
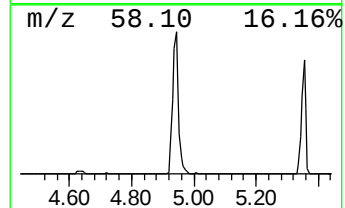
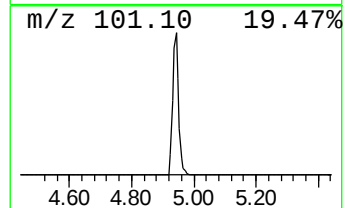
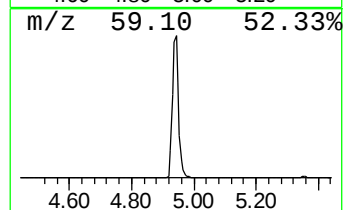
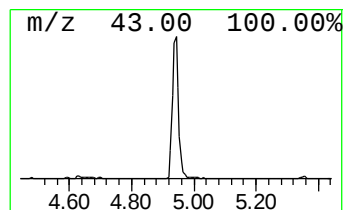
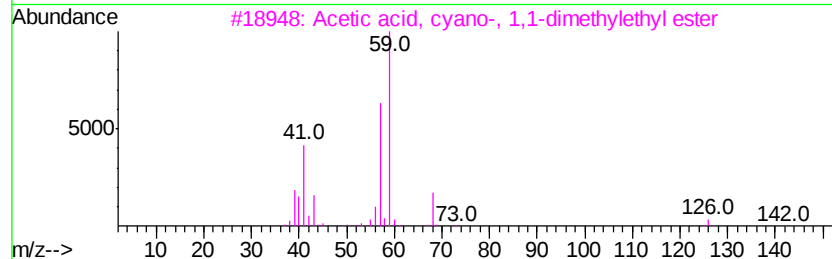
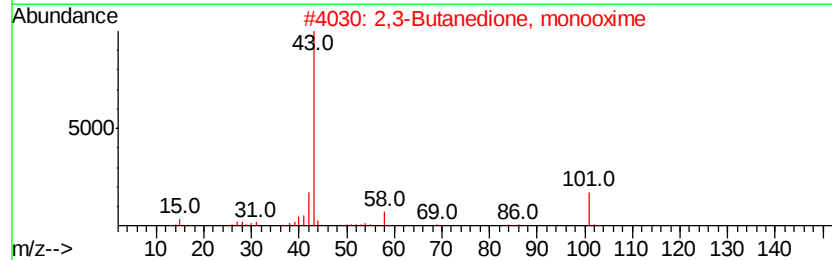
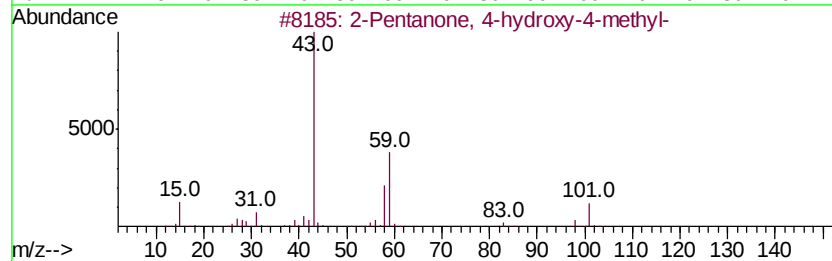
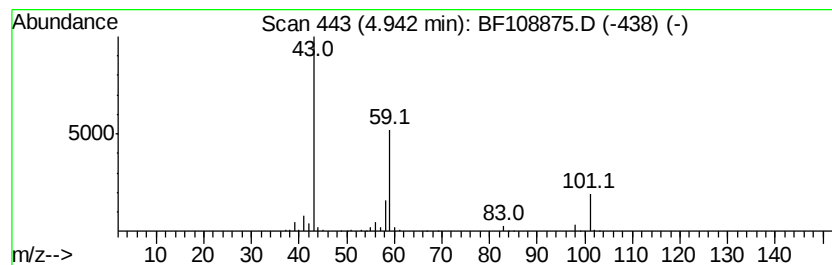
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.94	9.66 ng	611373	1,4-Dichlorobenzene-d4	6.74

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
4			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	9
5			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	9



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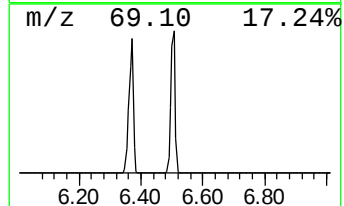
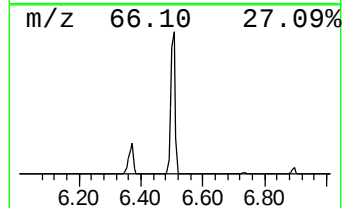
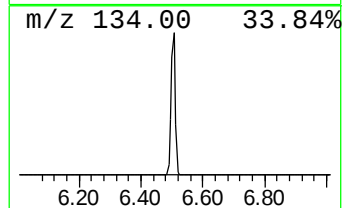
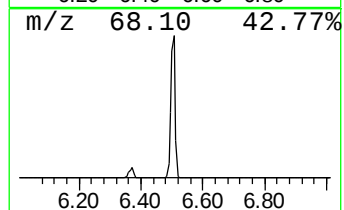
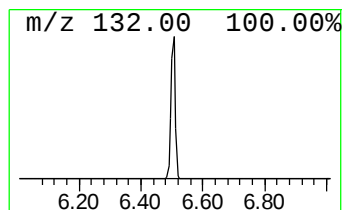
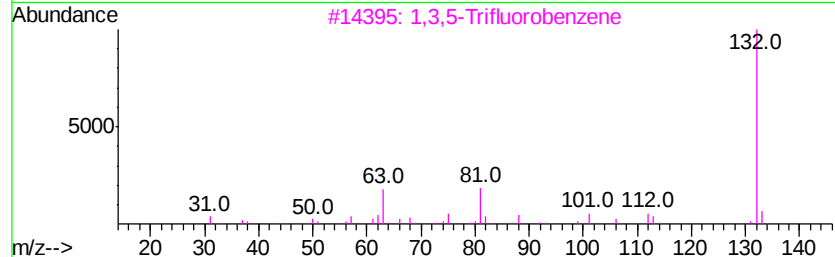
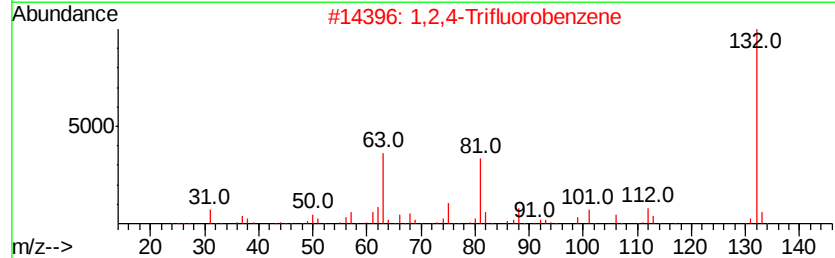
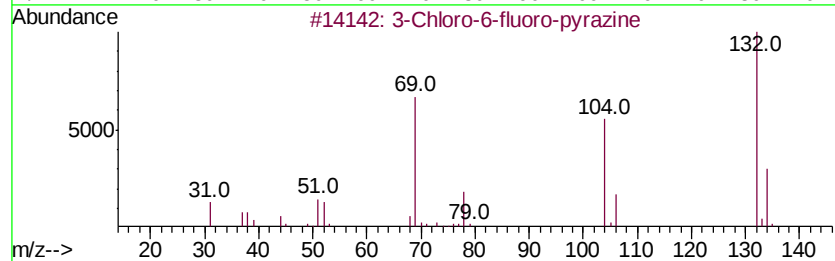
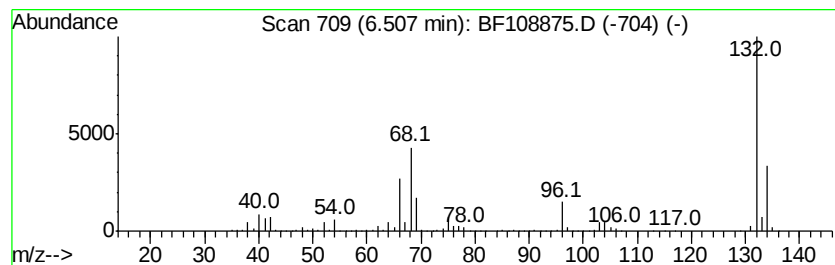
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Peak Number 2 unknown6.51 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.51	63.99 ng	4049510	1,4-Dichlorobenzene-d4	6.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2		1,2,4-Trifluorobenzene	132	C6H3F3	000367-23-7	9
3		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
4		Xenon	132	Xe	007440-63-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.94	9.7	ng	611373	1	6.74	1265710	20.0
unknown6.51	6.51	64.0	ng	4049510	1	6.74	1265710	20.0