

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG090618\
 Data File : BG036572.D
 Acq On : 6 Sep 2018 16:51
 Operator : JU/SJ
 Sample : J4801-04
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 CHRT-26957

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_G\METHODS\8270-BG082318.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	5.165	310	319	326	rBV	140867	219550	4.92%	1.027%
2	5.624	390	397	406	rBV	917311	1468381	32.92%	6.867%
3	7.210	660	667	677	rBV	893149	1505140	33.74%	7.039%
4	7.586	724	731	740	rBV	964725	1648944	36.97%	7.712%
5	8.056	805	811	819	rVB	202402	347080	7.78%	1.623%
6	8.379	858	866	875	rVB	738390	1294922	29.03%	6.056%
7	9.219	1000	1009	1018	rVB	624827	1128918	25.31%	5.280%
8	10.864	1279	1289	1295	rBV	285116	511480	11.47%	2.392%
9	13.309	1697	1705	1714	rBV	1748892	2779579	62.31%	13.000%
10	14.684	1932	1939	1946	rBV2	450813	754742	16.92%	3.530%
11	16.164	2184	2191	2214	rBV	1341453	2085607	46.75%	9.754%
12	17.421	2399	2405	2417	rBV2	663143	1002321	22.47%	4.688%
13	20.030	2843	2849	2859	rBV	3280137	4460774	100.00%	20.862%
14	21.687	3125	3131	3143	rBV	691414	1151159	25.81%	5.384%
15	24.901	3666	3678	3701	rBV2	291341	1023583	22.95%	4.787%

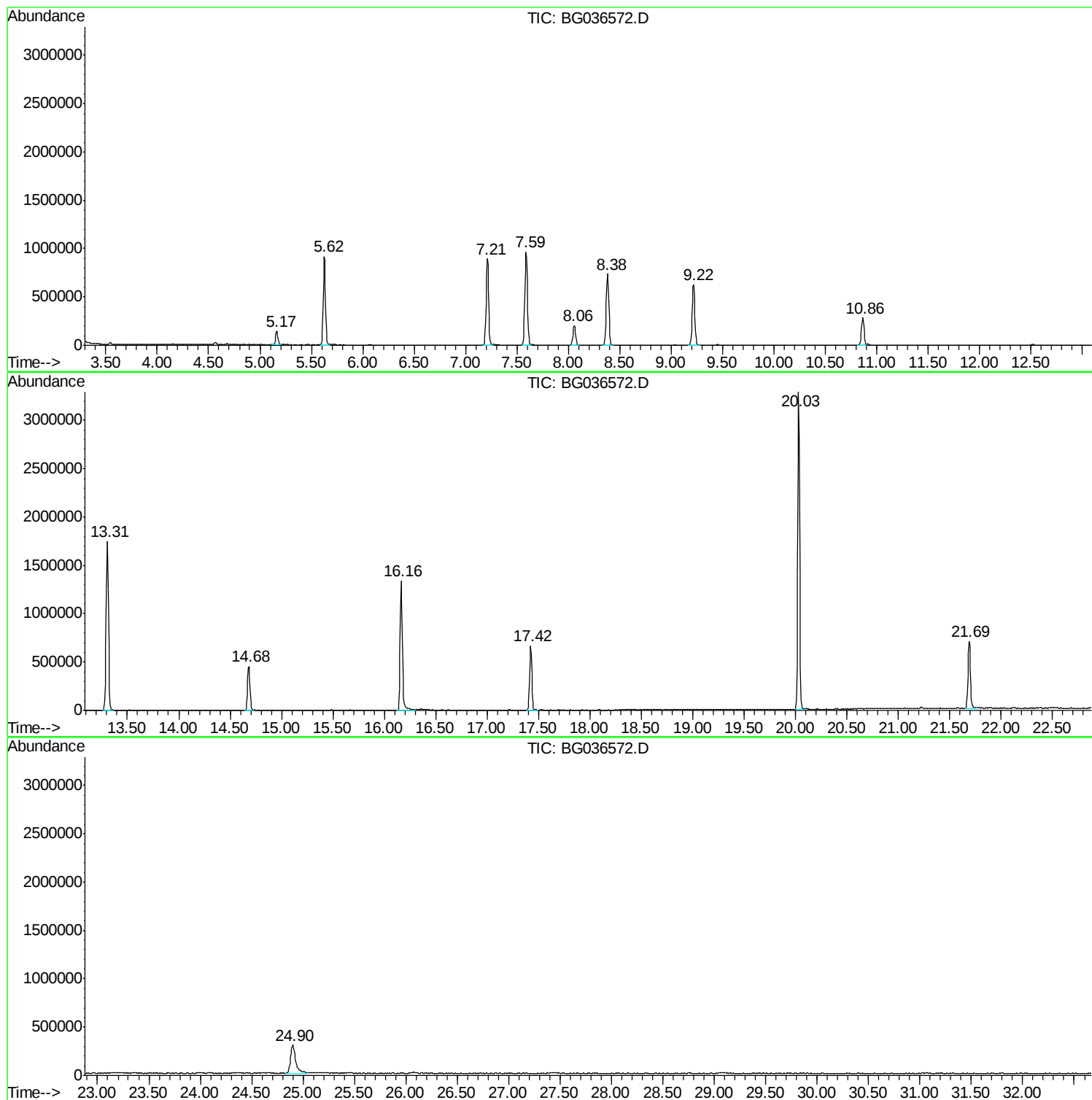
Sum of corrected areas: 21382180

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082318.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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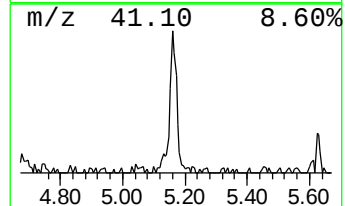
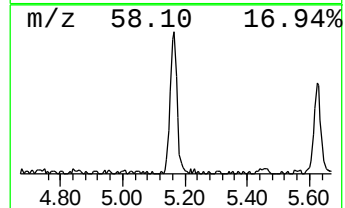
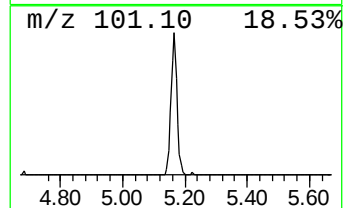
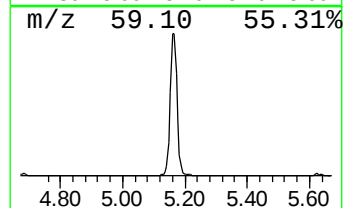
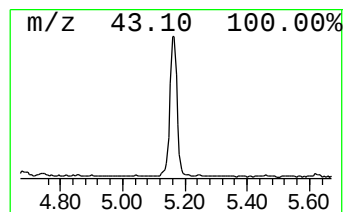
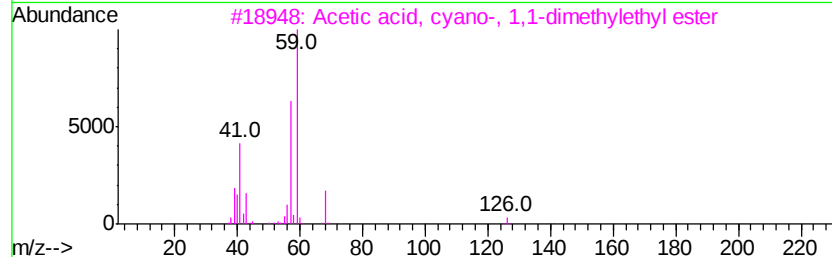
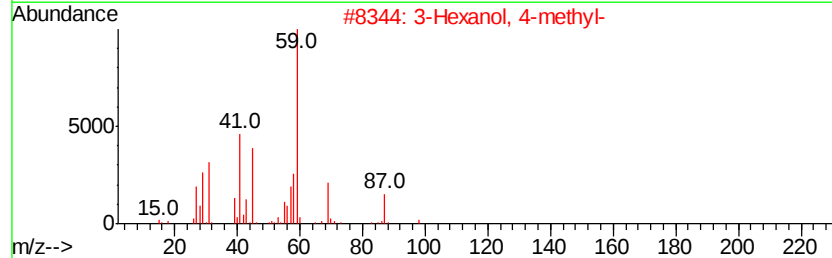
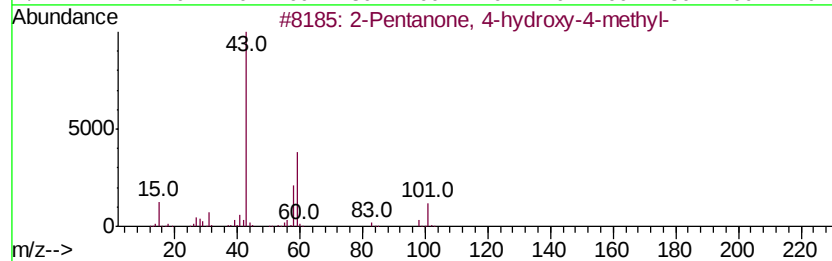
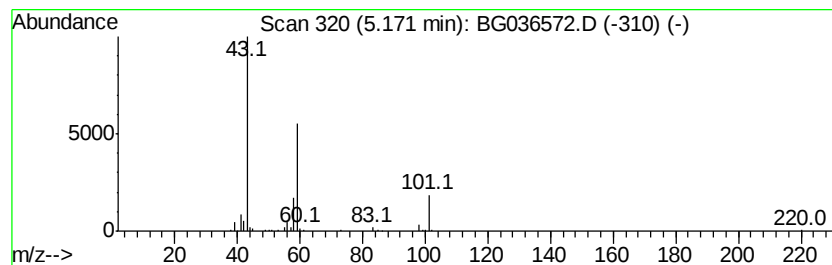
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	12.65 ng	219550	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
3		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
4		2,5,8,11-Tetraoxadodecane	178	C8H18O4	000112-49-2	23
5		3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	17



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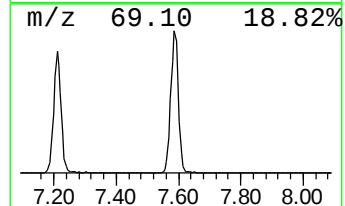
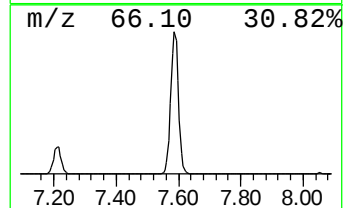
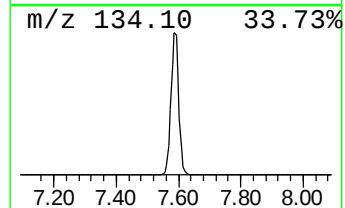
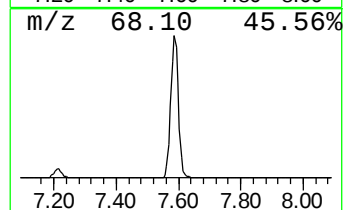
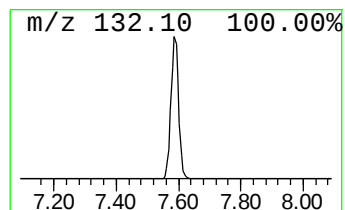
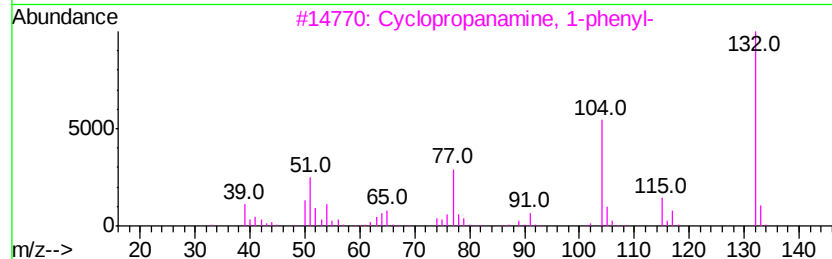
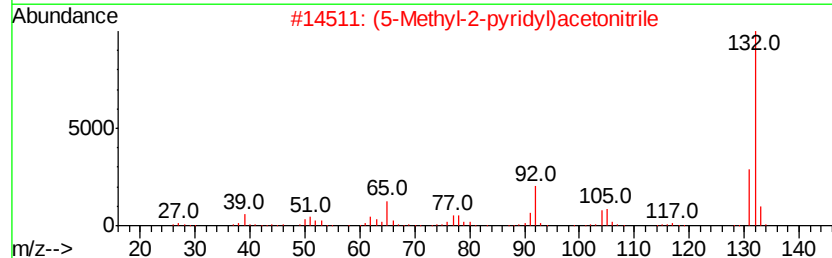
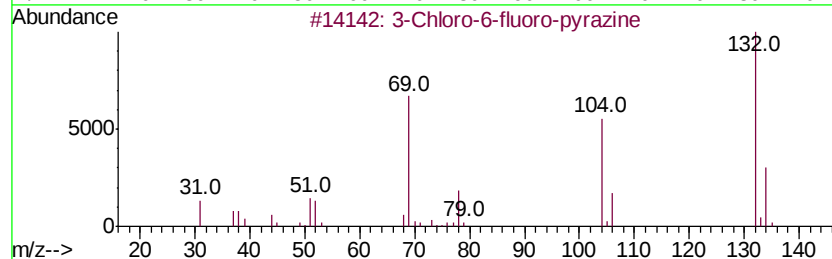
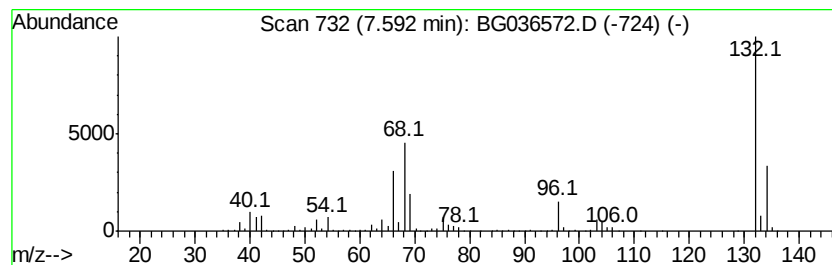
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Peak Number 2 unknown7.59 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	95.02 ng	1648940	1,4-Dichlorobenzene-d4	8.06

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
3			Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
4			5-Aminoindole	132	C8H8N2	005192-03-0	10
5			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.17	12.7	ng	219550	1	8.06	347080	20.0
unknown7.59	7.59	95.0	ng	1648940	1	8.06	347080	20.0