

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG122818\
 Data File : BG038890.D
 Acq On : 29 Dec 2018 1:42
 Operator : JU/SJ
 Sample : PB116060TB
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB116060TB

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-BG121218.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	5.138	343	349	355	rBV	22367	35016	1.61%	0.384%
2	5.608	421	429	443	rBV	403071	680465	31.29%	7.458%
3	7.218	695	703	713	rBV	438380	769295	35.37%	8.431%
4	7.588	758	766	776	rBV	407993	731057	33.61%	8.012%
5	8.052	838	845	853	rVB	72241	126119	5.80%	1.382%
6	8.376	892	900	908	rBV	283911	503911	23.17%	5.523%
7	9.233	1037	1046	1058	rBV	231495	427916	19.68%	4.690%
8	10.873	1318	1325	1336	rBV	96816	184451	8.48%	2.022%
9	13.311	1733	1740	1751	rBV	736971	1186362	54.55%	13.002%
10	14.692	1968	1975	1985	rBV2	184266	309370	14.22%	3.391%
11	16.184	2223	2229	2255	rBV3	274311	609162	28.01%	6.676%
12	17.441	2436	2443	2454	rBV	275965	422382	19.42%	4.629%
13	20.044	2879	2886	2894	rBV	1628498	2174847	100.00%	23.836%
14	21.719	3164	3171	3181	rBV	293884	498047	22.90%	5.459%
15	25.021	3723	3733	3748	rVB	156315	465799	21.42%	5.105%

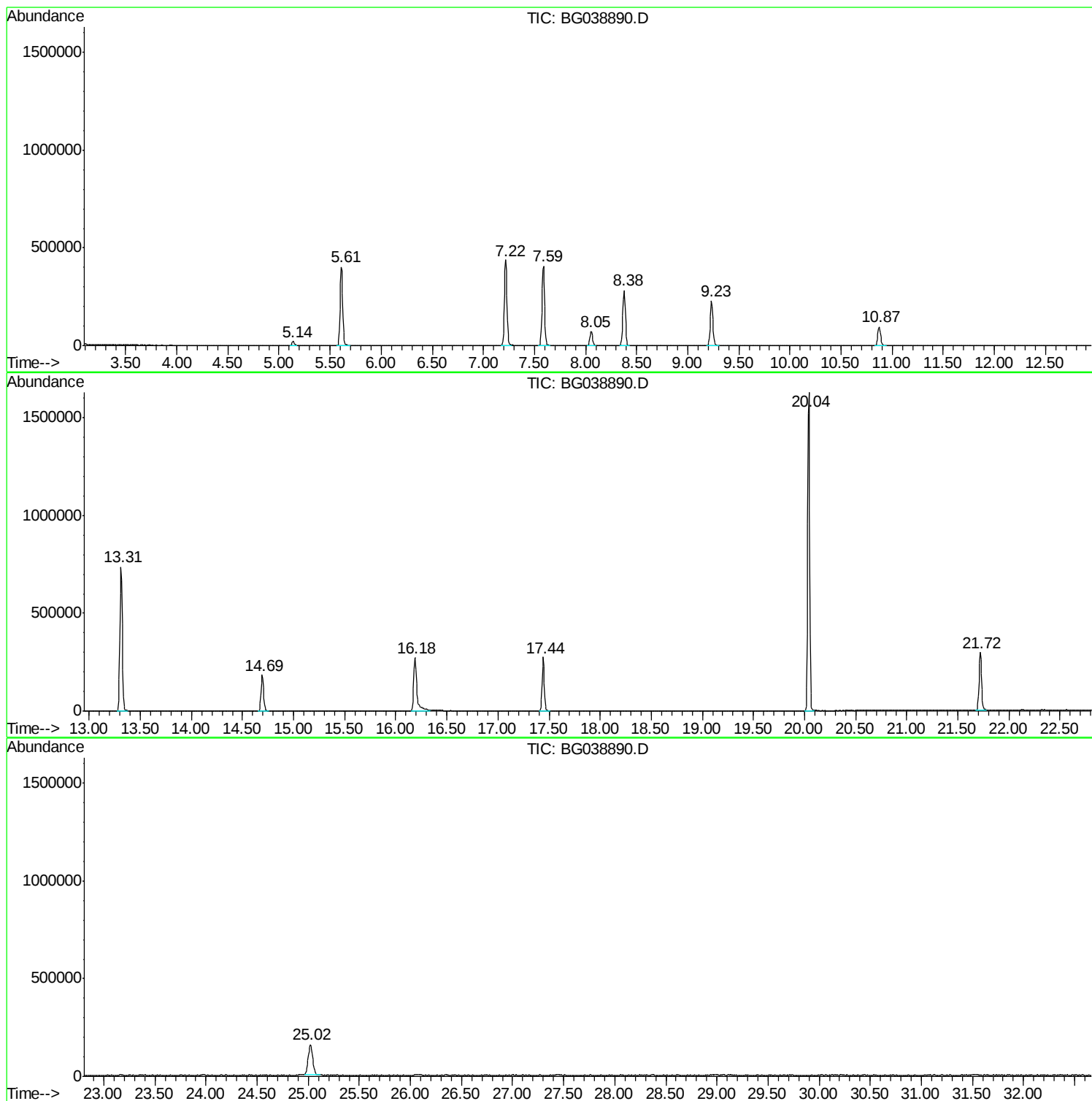
Sum of corrected areas: 9124199

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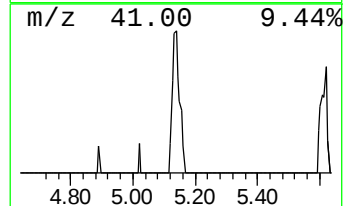
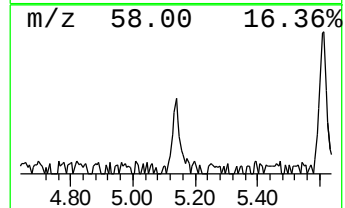
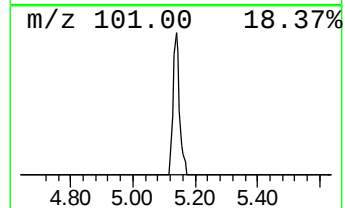
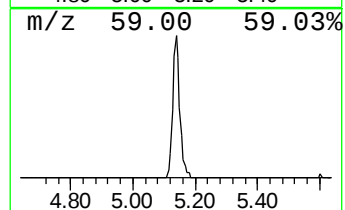
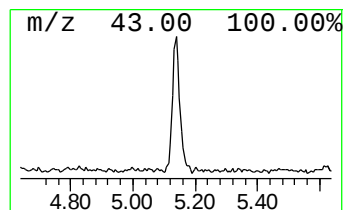
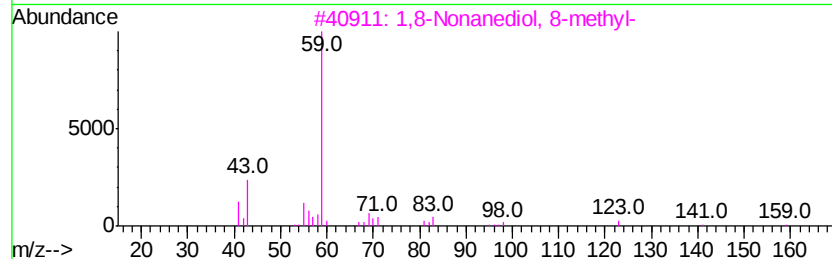
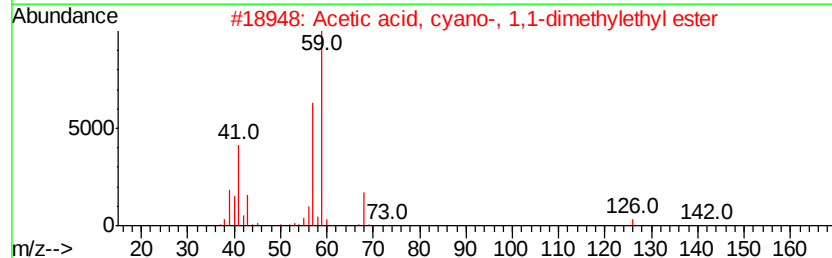
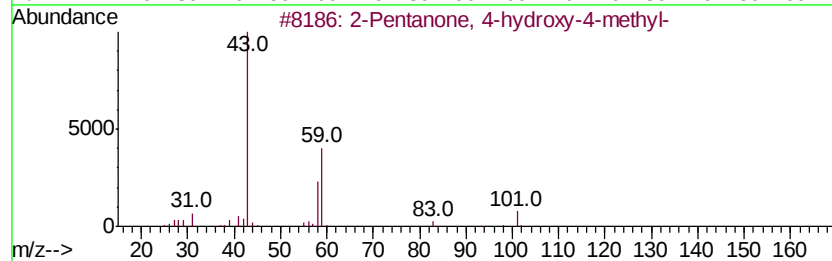
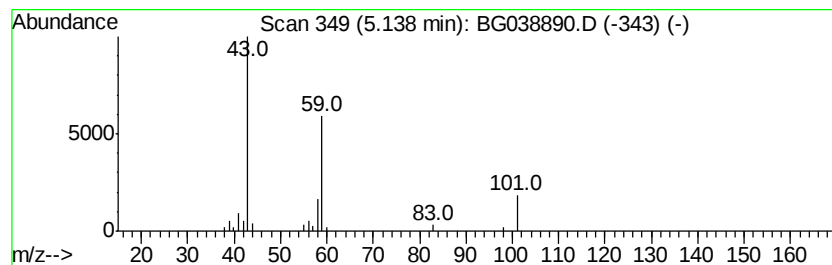
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	5.55 ng	35016	1,4-Dichlorobenzene-d4	8.05

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56	
2	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25	
3	1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	23	
4	Propylamine	59	C3H9N	000107-10-8	9	
5	4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9	



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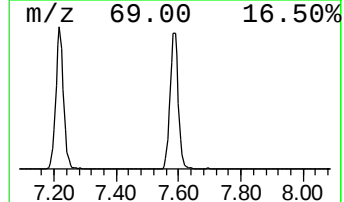
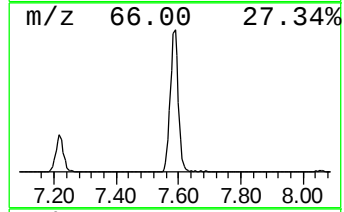
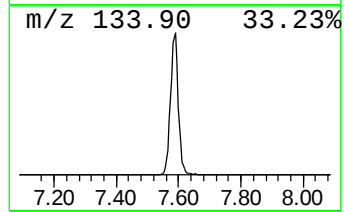
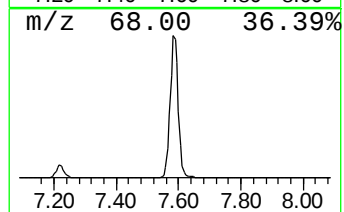
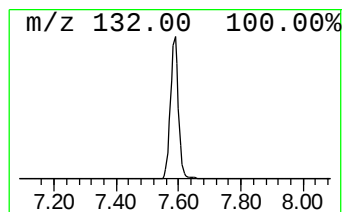
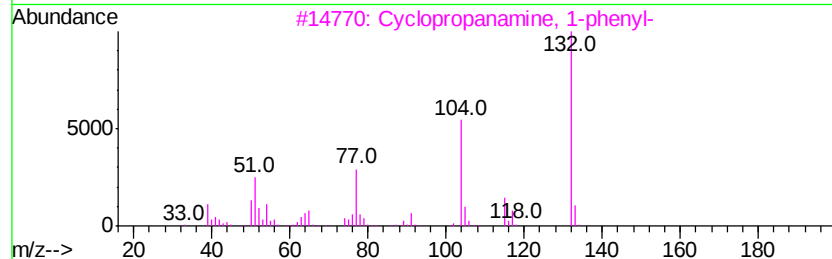
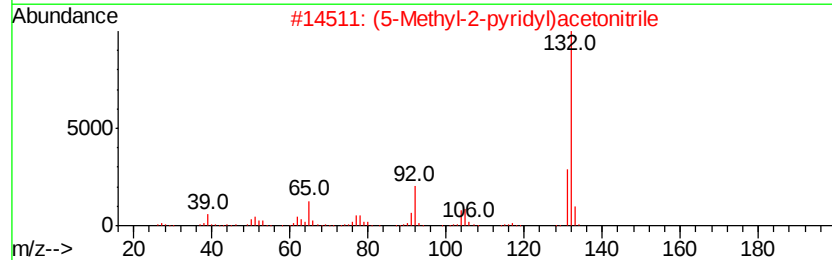
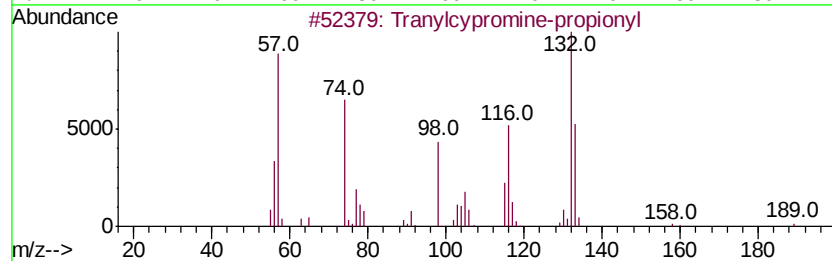
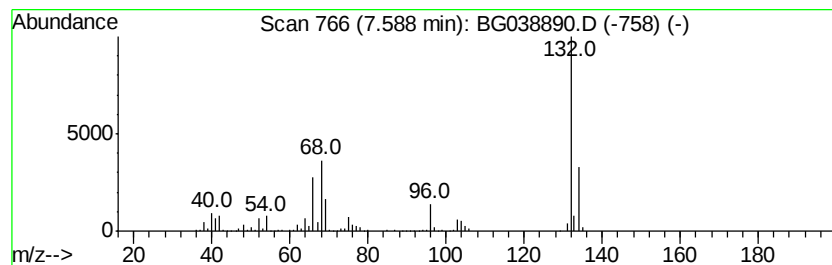
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Peak Number 2 unknown7.59 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	115.93 ng	731057	1,4-Dichlorobenzene-d4	8.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	16
2		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
3		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.14	5.5	ng	35016	1	8.05	126119	20.0
unknown7.59	7.59	115.9	ng	731057	1	8.05	126119	20.0