

Data Path : Z:\HPCHEM1\BNA G\DATA\BG113017\
 Data File : BG031297.D
 Acq On : 30 Nov 2017 21:00
 Operator : SJ/JU
 Sample : I6634-05
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 112717-TIDO-E-D-B

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:\HPCHEM1\BNA_G\METHODS\8270-BG111017.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	5.210	320	327	340	rBV2	71448	117664	1.82%	0.501%
2	5.691	403	409	421	rBV	320499	571598	8.82%	2.435%
3	7.295	675	682	701	rBV	189607	401093	6.19%	1.709%
4	7.666	737	745	755	rBV	632599	1135816	17.53%	4.839%
5	8.130	817	824	833	rBV	243757	422578	6.52%	1.800%
6	8.453	872	879	892	rBV	965410	1740759	26.87%	7.417%
7	9.299	1015	1023	1037	rBV	677533	1326120	20.47%	5.650%
8	10.950	1296	1304	1320	rBV	282660	626924	9.68%	2.671%
9	13.377	1708	1717	1739	rBV	2366100	3799559	58.65%	16.189%
10	14.751	1941	1951	1960	rBV2	531390	952325	14.70%	4.058%
11	16.238	2198	2204	2224	rBV	1132399	1838843	28.38%	7.835%
12	17.495	2410	2418	2424	rBV	768311	1210776	18.69%	5.159%
13	20.086	2852	2859	2871	rBV	4806377	6478612	100.00%	27.604%
14	21.767	3138	3145	3158	rBV	841571	1457382	22.50%	6.210%
15	25.069	3694	3707	3720	rBV	483206	1390019	21.46%	5.923%

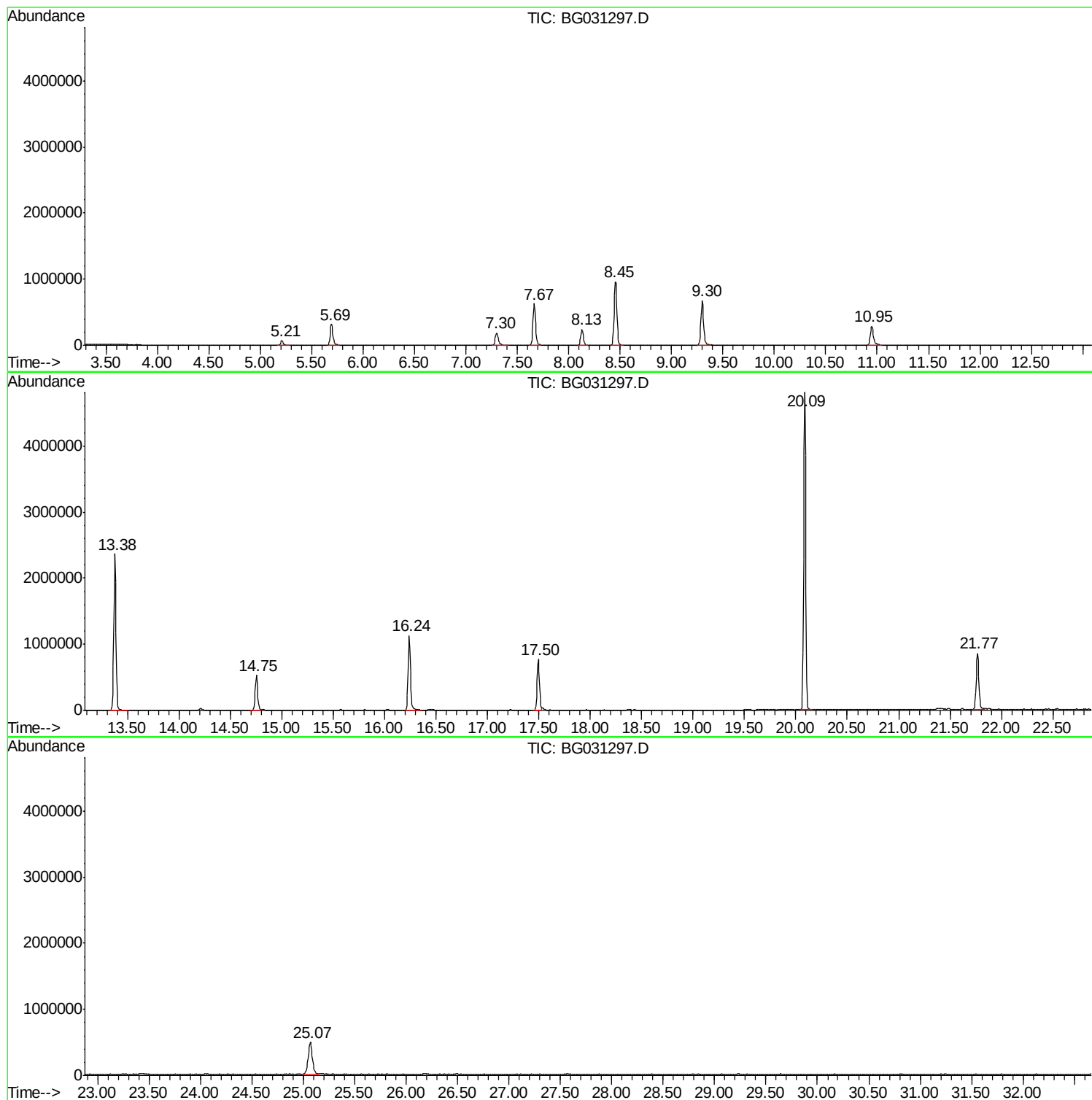
Sum of corrected areas: 23470068

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG111017.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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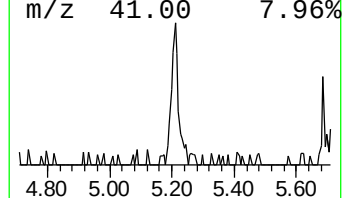
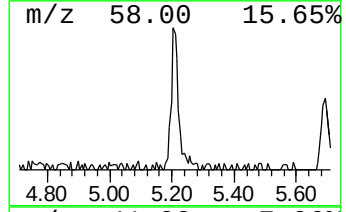
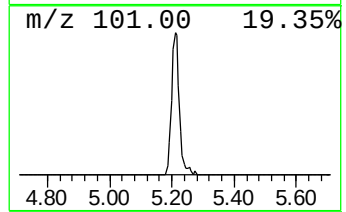
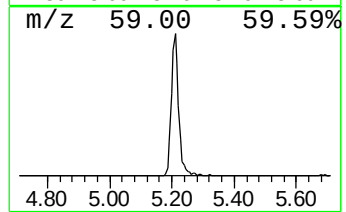
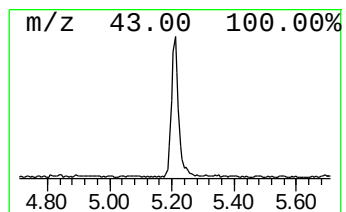
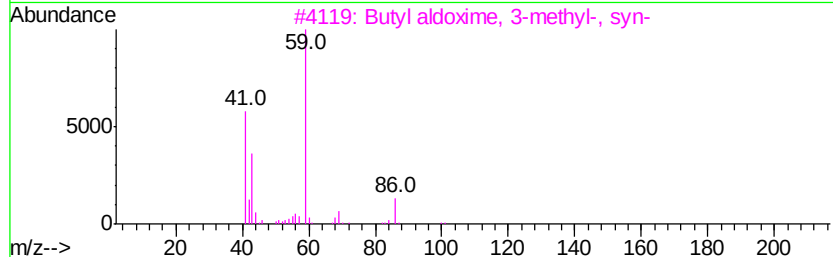
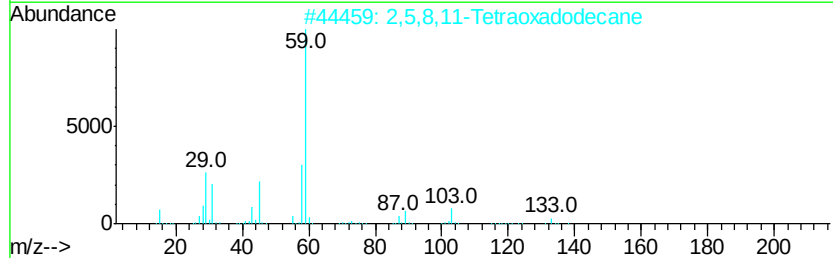
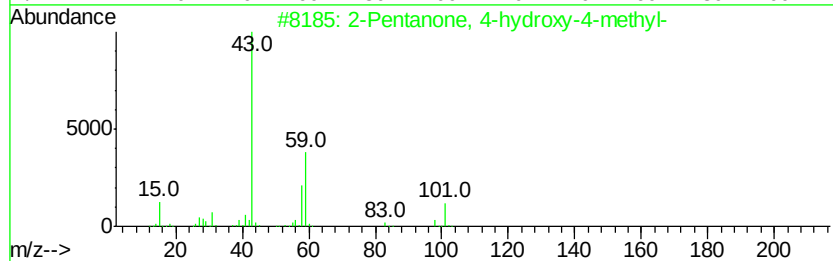
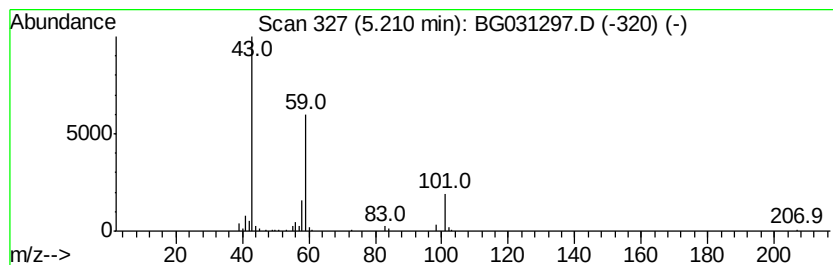
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.21	5.57 ng	117664	1,4-Dichlorobenzene-d4	8.13

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			2,5,8,11-Tetraoxadodecane	178	C8H18O4	000112-49-2	23
3			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	17
4			Silane, dimethyl-	60	C2H8Si	001111-74-6	9
5			Formamide, N,N-diethyl-	101	C5H11NO	000617-84-5	9



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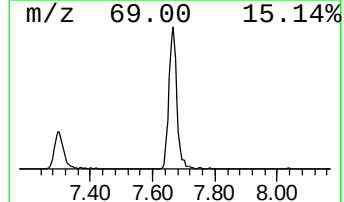
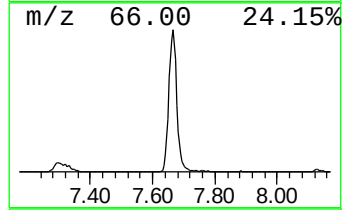
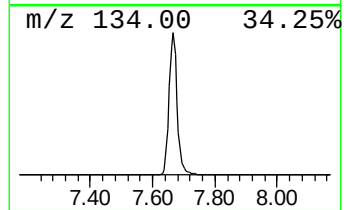
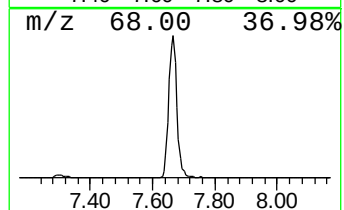
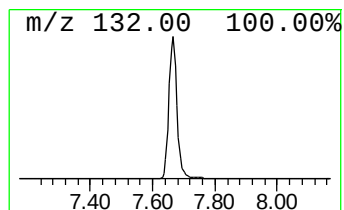
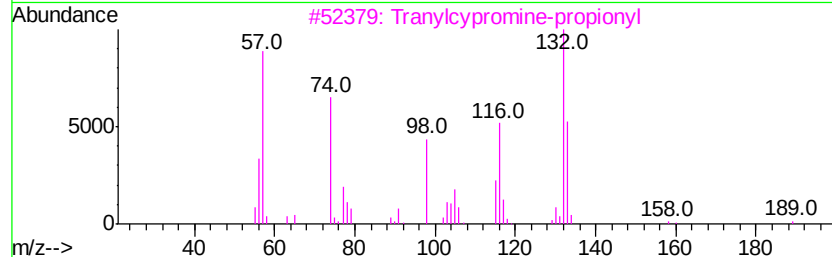
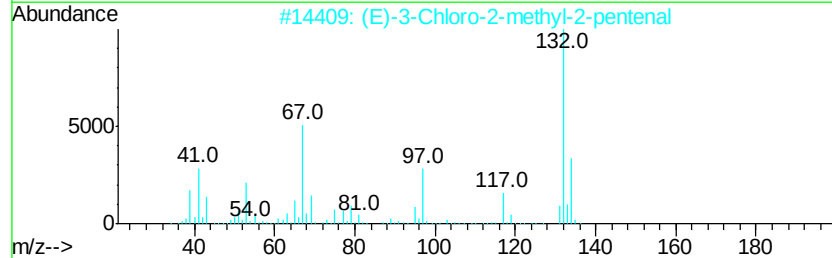
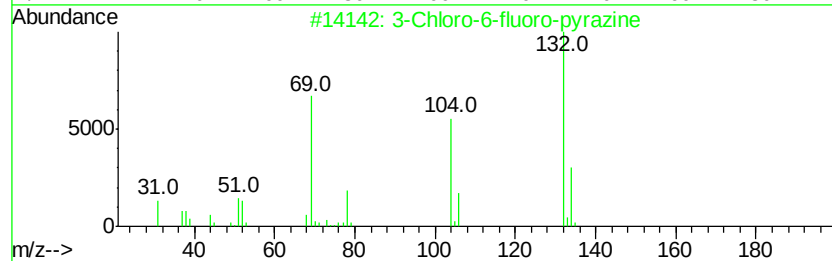
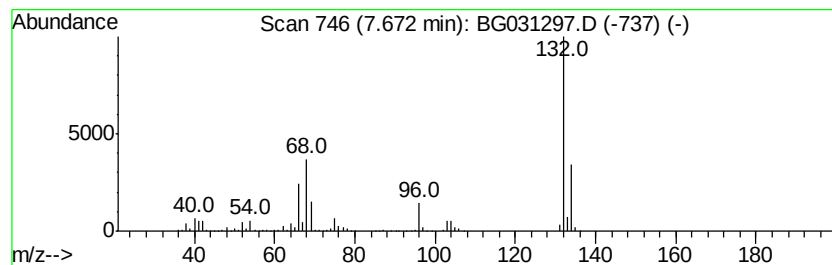
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Peak Number 2 unknown7.67 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.67	53.76 ng	1135820	1,4-Dichlorobenzene-d4	8.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	10
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.21	5.6	ng	117664	1	8.13	422578	20.0
unknown7.67	7.67	53.8	ng	1135820	1	8.13	422578	20.0