

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG090619\
Data File : BG042749.D
Acq On : 6 Sep 2019 15:39
Operator : HP/JU
Sample : PB122825BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PB122825BL

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-BG082219.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	5.568	415	422	439	rBV	378597	623094	20.02%	5.592%
2	7.178	688	696	709	rBV	385385	684738	22.00%	6.146%
3	7.542	750	758	770	rBV	415594	723040	23.23%	6.489%
4	8.006	831	837	845	rBV	86393	147726	4.75%	1.326%
5	8.335	884	893	903	rBV	328488	576420	18.52%	5.173%
6	9.175	1028	1036	1051	rBV	206621	404728	13.00%	3.632%
7	10.832	1310	1318	1334	rBV	97725	204222	6.56%	1.833%
8	13.282	1728	1735	1761	rBV	825479	1373827	44.13%	12.330%
9	14.669	1964	1971	1985	rBV	191346	358110	11.50%	3.214%
10	16.161	2217	2225	2248	rBV	590023	1173393	37.69%	10.531%
11	17.419	2432	2439	2456	rBV2	282298	532137	17.09%	4.776%
12	20.027	2870	2883	2901	rBV	2264809	3112904	100.00%	27.938%
13	21.696	3159	3167	3184	rBV2	313494	616807	19.81%	5.536%
14	24.951	3709	3721	3744	rBV2	184960	610896	19.62%	5.483%

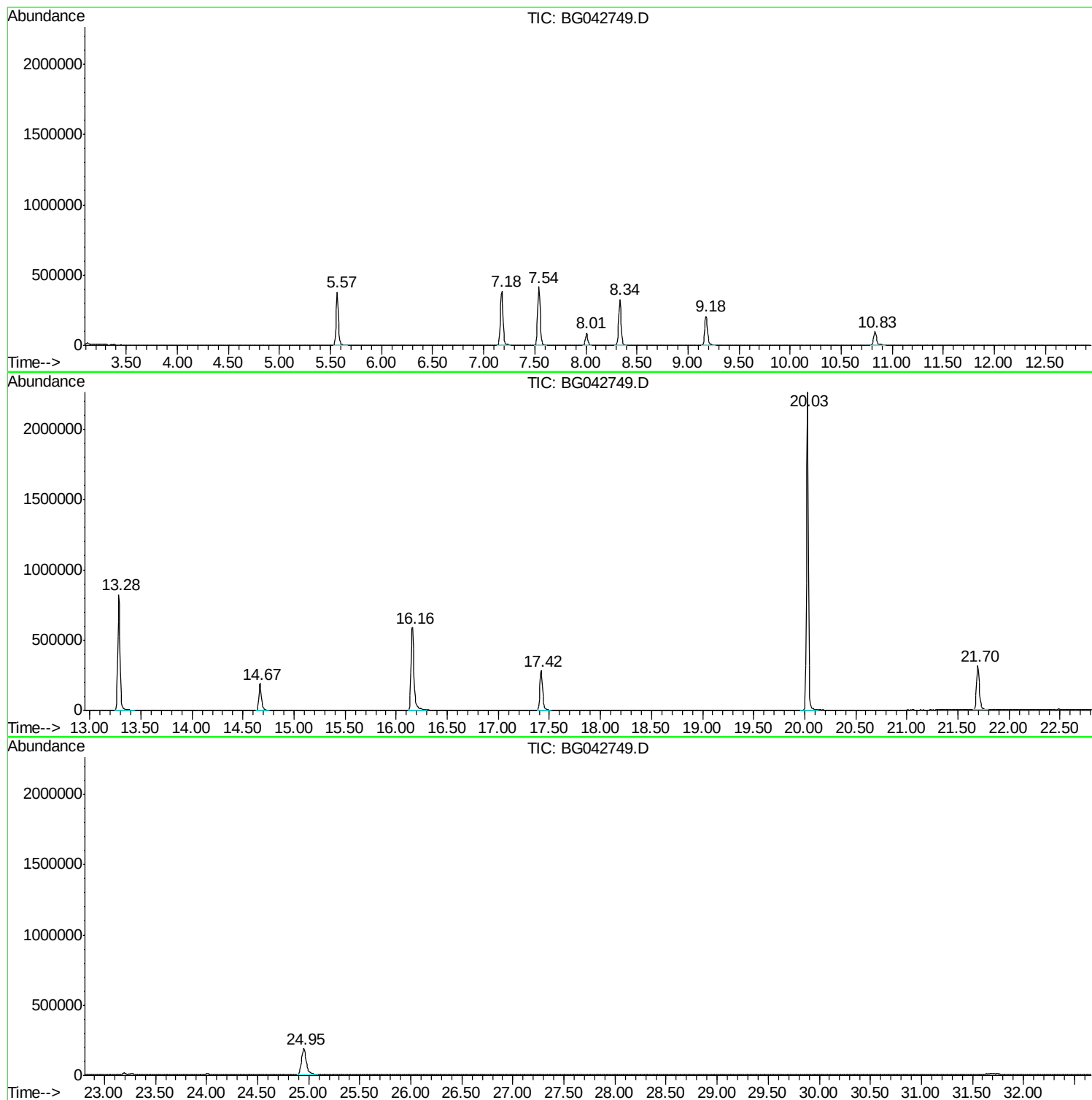
Sum of corrected areas: 11142042

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082219.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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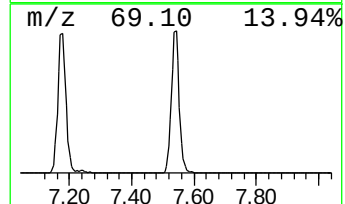
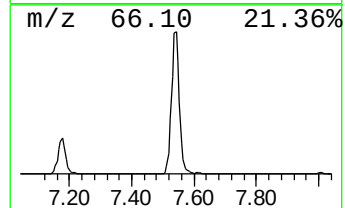
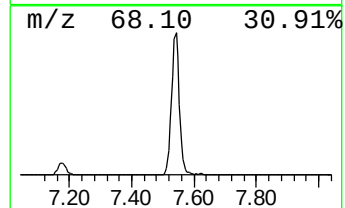
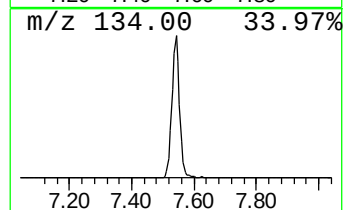
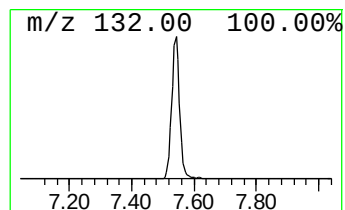
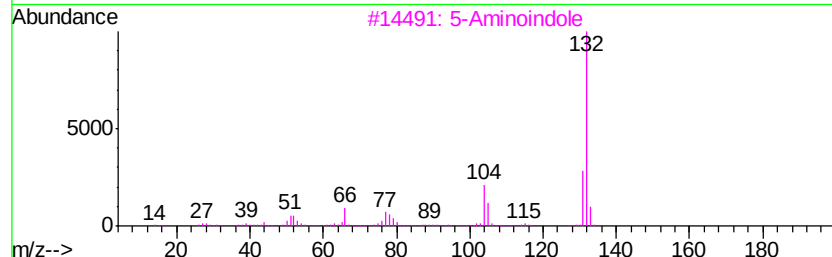
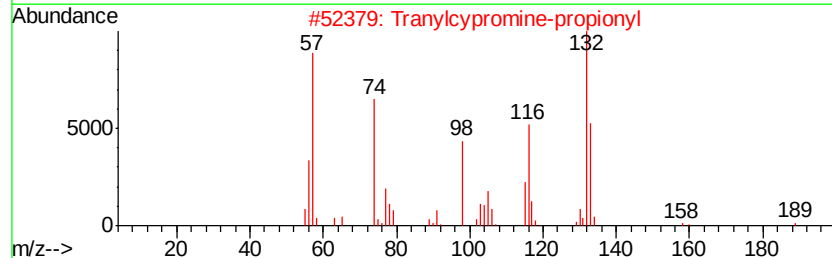
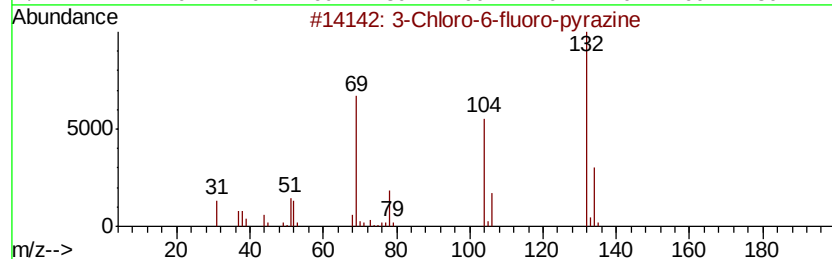
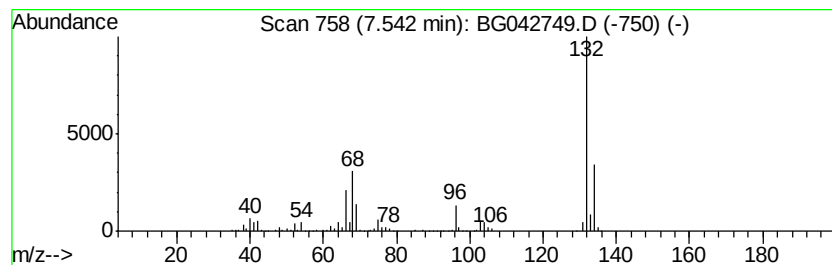
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Peak Number 1 unknown7.54 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.54	97.89 ng	723040	1,4-Dichlorobenzene-d4	8.01

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	16
3			5-Aminoindole	132	C8H8N2	005192-03-0	16
4			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
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
unknown7.54	7.54	97.9	ng	723040	1	8.01	147726	20.0