

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111119\
 Data File : BG043605.D
 Acq On : 12 Nov 2019 6:06
 Operator : JU
 Sample : PB124694BL
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB124694BL

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-BG102119.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	5.596	420	427	441	rBV	336417	602152	23.04%	5.628%
2	7.195	691	699	712	rBV	341296	662781	25.36%	6.195%
3	7.547	751	759	772	rBV	372000	675224	25.83%	6.311%
4	7.999	829	836	849	rBV	103311	185645	7.10%	1.735%
5	8.323	883	891	903	rBV	415891	751416	28.75%	7.024%
6	9.163	1025	1034	1049	rBV	268339	571662	21.87%	5.343%
7	10.808	1306	1314	1328	rBV	104729	230966	8.84%	2.159%
8	13.252	1723	1730	1744	rBV	892461	1518306	58.09%	14.192%
9	14.633	1958	1965	1976	rBV2	196906	363731	13.92%	3.400%
10	16.125	2212	2219	2232	rBV	443593	857435	32.81%	8.015%
11	17.377	2425	2432	2452	rBV	240158	461036	17.64%	4.309%
12	19.985	2870	2876	2887	rBV	1772091	2613648	100.00%	24.430%
13	21.642	3152	3158	3175	rBV	279475	555011	21.24%	5.188%
14	23.893	3535	3541	3549	rBV7	12762	33932	1.30%	0.317%
15	24.827	3689	3700	3718	rBV	187646	615543	23.55%	5.754%

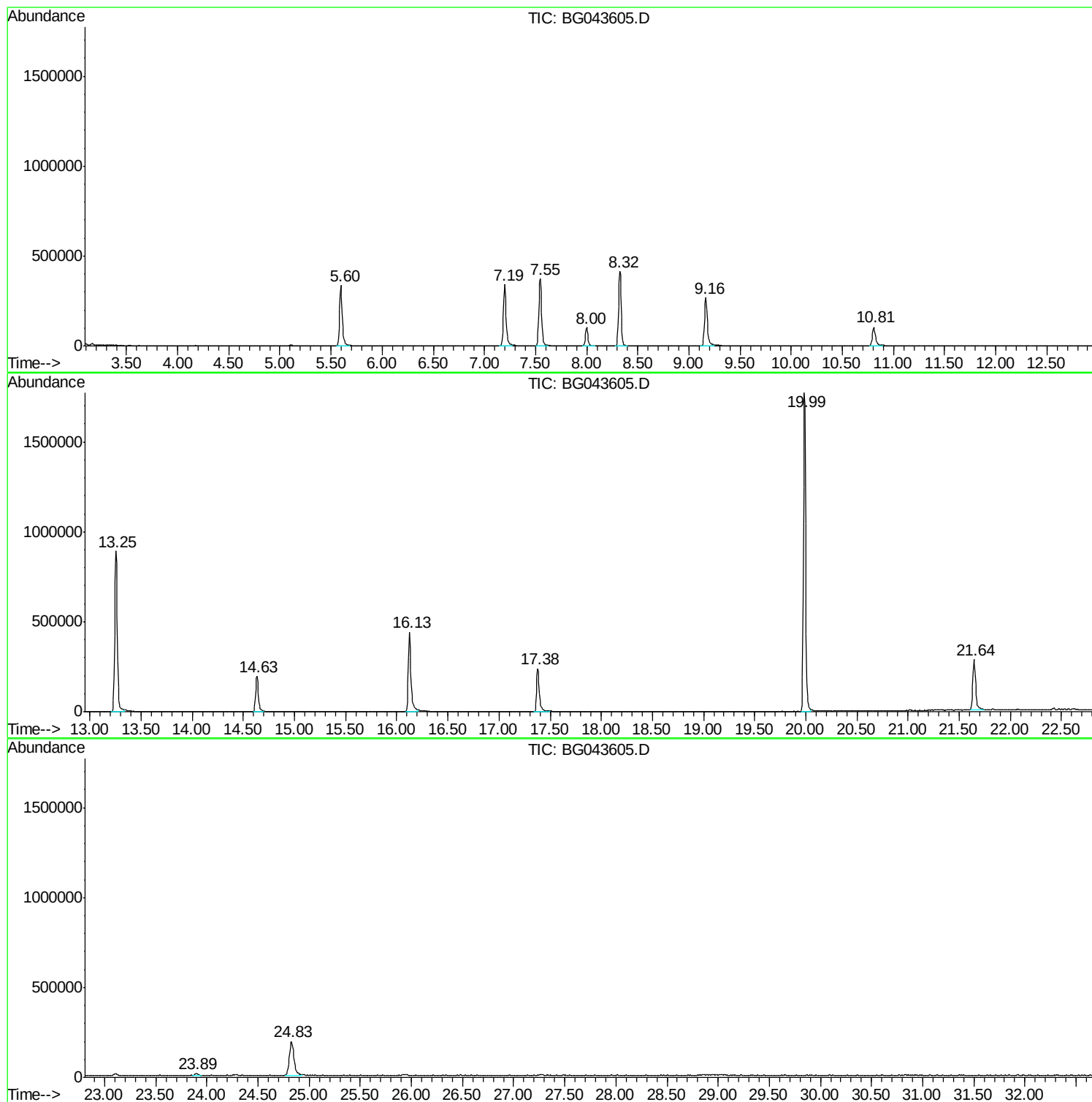
Sum of corrected areas: 10698488

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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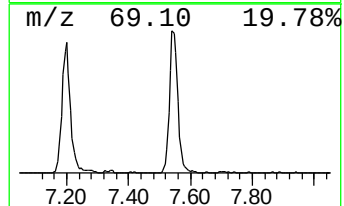
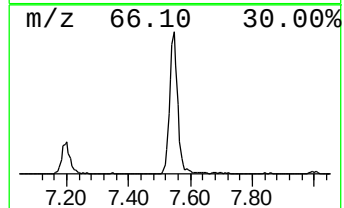
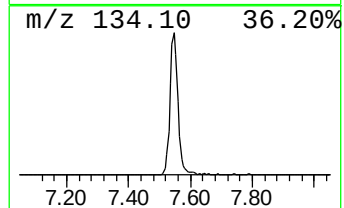
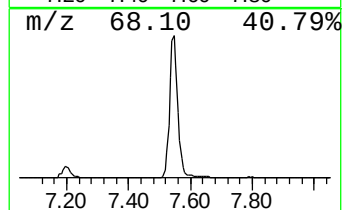
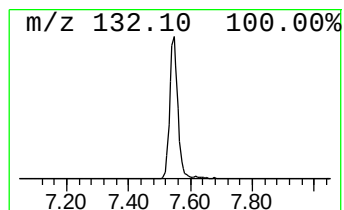
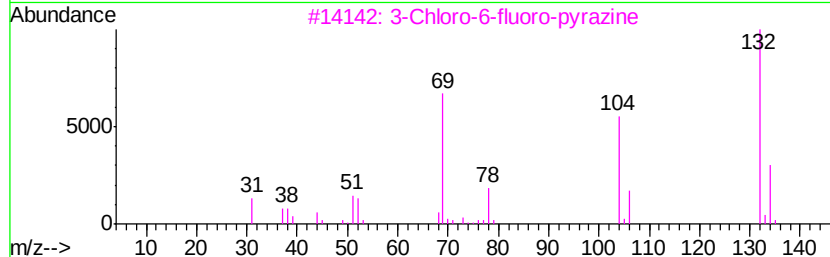
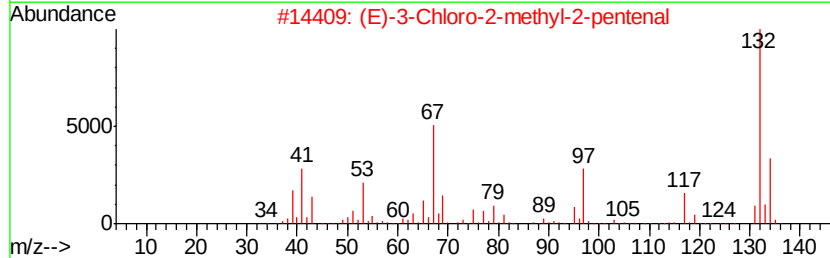
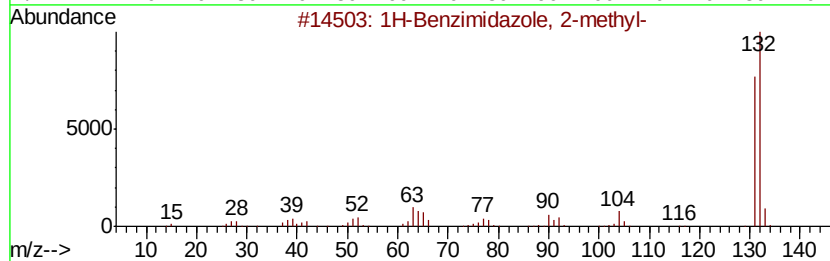
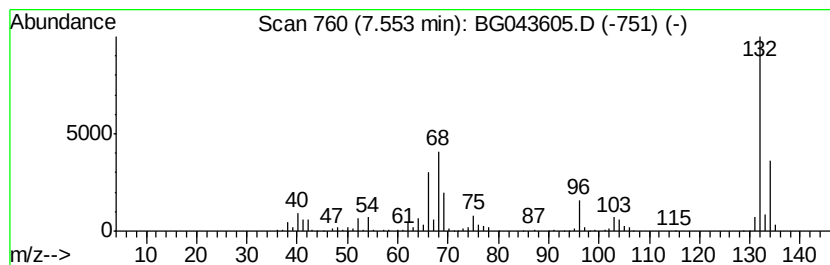
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Peak Number 1 unknown7.55 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.55	72.74 ng	675224	1,4-Dichlorobenzene-d4	8.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
4		5-Aminoindole	132	C8H8N2	005192-03-0	16
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown7.55	7.55	72.7	ng	675224	1	8.00	185645	20.0