

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG102319\
 Data File : BG043310.D
 Acq On : 23 Oct 2019 16:06
 Operator : JU
 Sample : PB124203BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB124203BL

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	5.608	422	429	445	rBV	474830	851599	24.43%	6.342%
2	7.200	691	700	711	rBV	499614	927360	26.60%	6.907%
3	7.564	754	762	776	rBV	510525	949582	27.24%	7.072%
4	8.023	834	840	850	rBV	115276	202977	5.82%	1.512%
5	8.346	887	895	904	rBV	405464	750984	21.54%	5.593%
6	9.186	1030	1038	1055	rBV	246094	542842	15.57%	4.043%
7	10.837	1309	1319	1330	rBV	99716	230132	6.60%	1.714%
8	13.275	1727	1734	1750	rBV	835228	1498828	42.99%	11.163%
9	14.650	1962	1968	1983	rBV2	209011	394759	11.32%	2.940%
10	16.142	2215	2222	2242	rBV	757190	1418677	40.69%	10.566%
11	17.400	2429	2436	2452	rBV	280024	545406	15.64%	4.062%
12	20.009	2873	2880	2895	rBV	2435953	3486355	100.00%	25.965%
13	21.665	3156	3162	3177	rBV	393931	804537	23.08%	5.992%
14	24.868	3696	3707	3721	rBV2	256295	823260	23.61%	6.131%

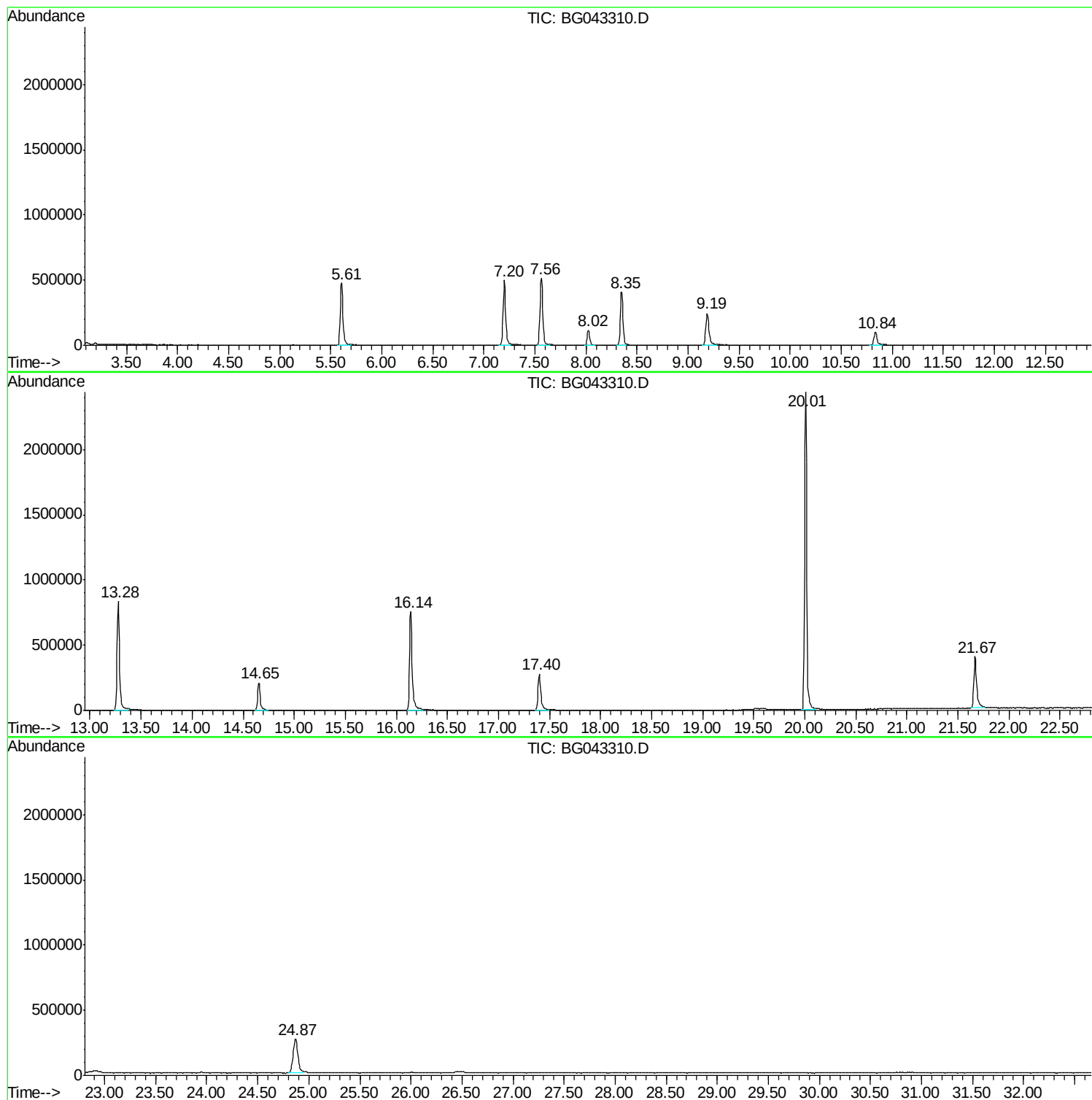
Sum of corrected areas: 13427298

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.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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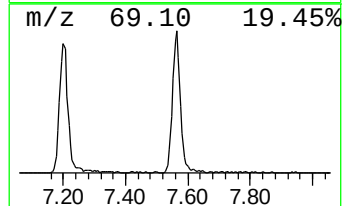
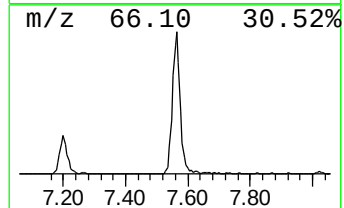
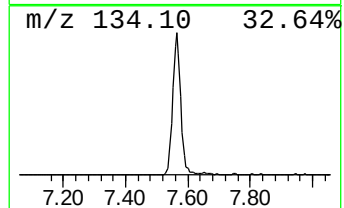
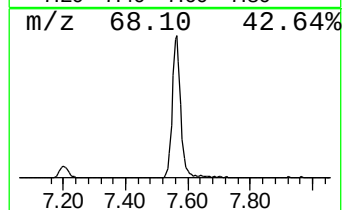
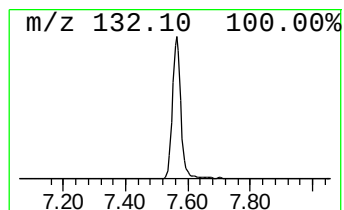
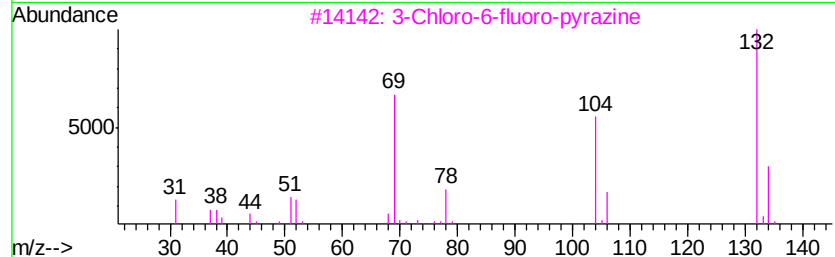
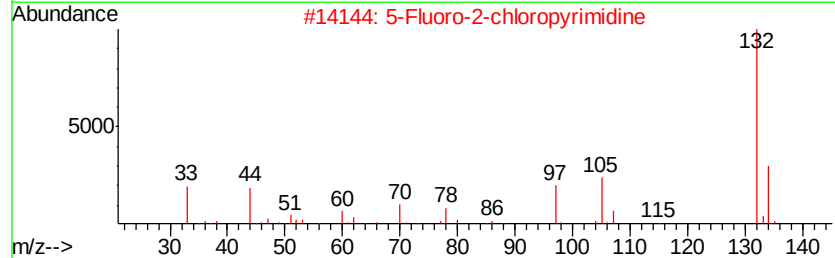
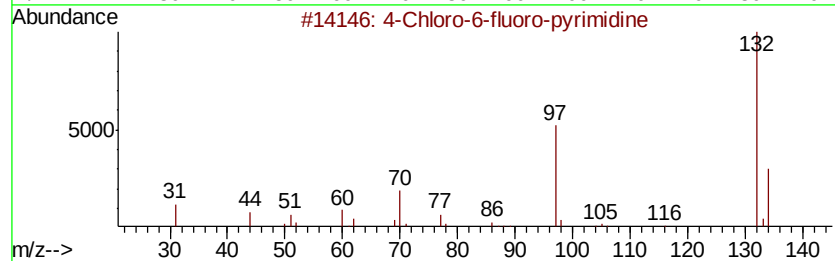
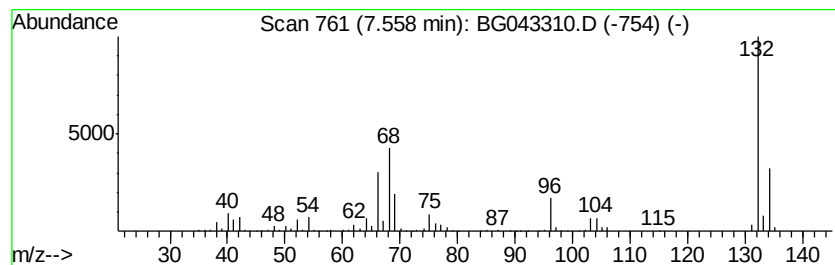
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Peak Number 1 unknown7.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	93.57 ng	949582	1,4-Dichlorobenzene-d4	8.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	27
3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
4		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	17
5		Tranlylcypromine-propionyl	189	C12H15NO	1000123-86-3	12



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
unknown7.56	7.56	93.6	ng	949582	1	8.02	202977	20.0