

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN032719\
 Data File : BN004999.D
 Acq On : 27 Mar 2019 17:30
 Operator : JU/SJ
 Sample : PB118322BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB118322BL

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_N\METHODS\8270-BN032619.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	5.269	382	388	395	rVB	147331	199240	45.21%	8.141%
2	6.852	651	657	666	rBV	133195	208785	47.38%	8.531%
3	7.211	711	718	724	rBV	153943	228636	51.88%	9.343%
4	7.675	791	797	802	rVB	32323	46098	10.46%	1.884%
5	7.993	845	851	857	rBV	114674	175190	39.76%	7.159%
6	8.840	989	995	1003	rVB	82620	126927	28.80%	5.186%
7	10.457	1264	1270	1278	rVB	41351	65796	14.93%	2.689%
8	12.934	1685	1691	1700	rVB	232338	332791	75.52%	13.598%
9	14.322	1921	1927	1933	rBV2	58964	87189	19.79%	3.563%
10	15.816	2176	2181	2190	rBV	179951	241527	54.81%	9.869%
11	17.069	2389	2394	2402	rVB2	74620	103936	23.59%	4.247%
12	19.704	2836	2842	2847	rBV	376838	440672	100.00%	18.007%
13	21.269	3103	3108	3115	rBV	79218	101498	23.03%	4.147%
14	23.510	3482	3489	3500	rVB	48674	88982	20.19%	3.636%

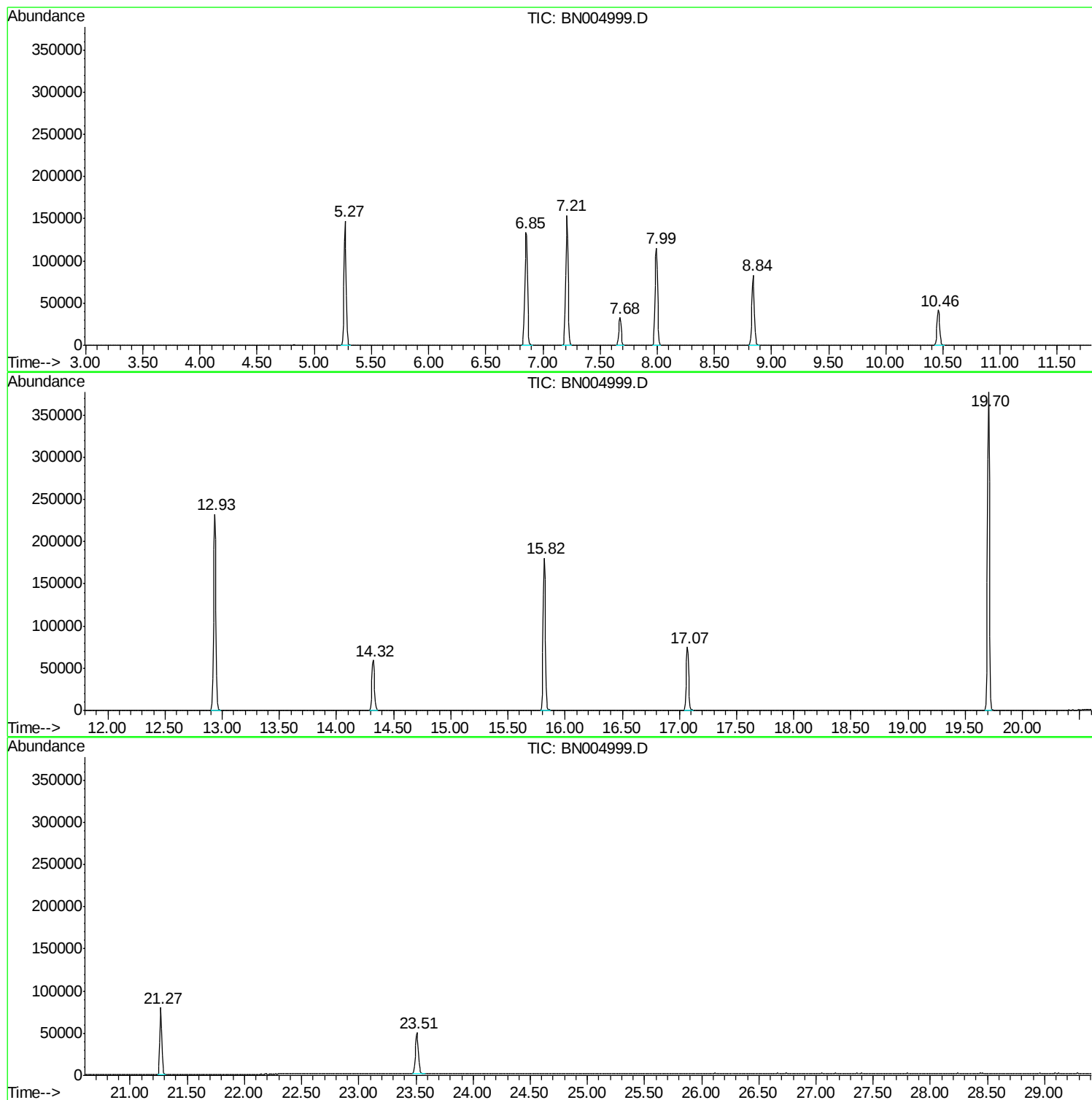
Sum of corrected areas: 2447267

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-BN032619.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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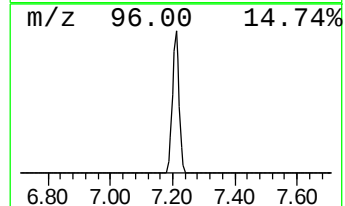
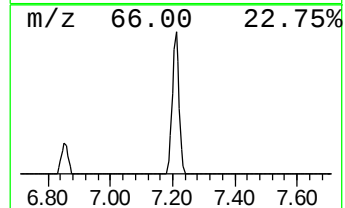
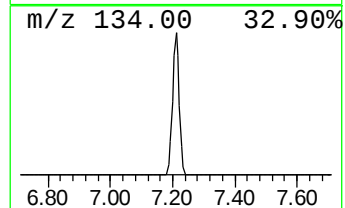
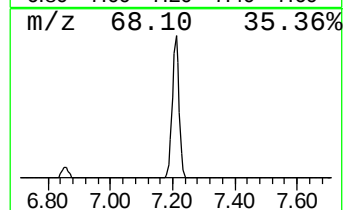
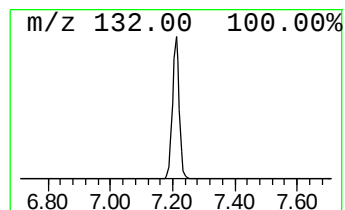
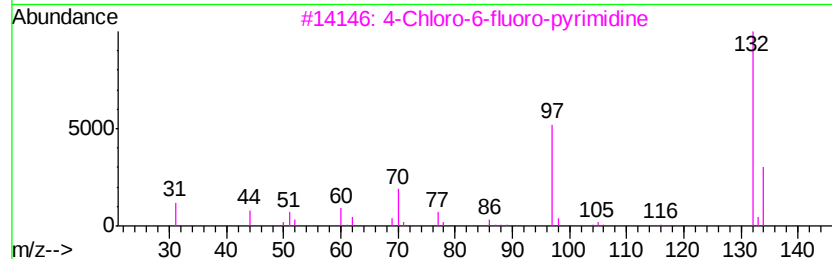
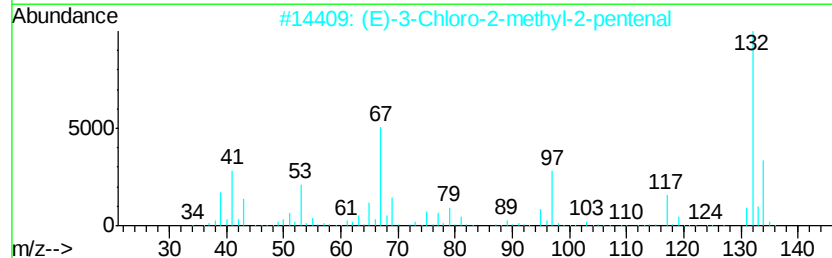
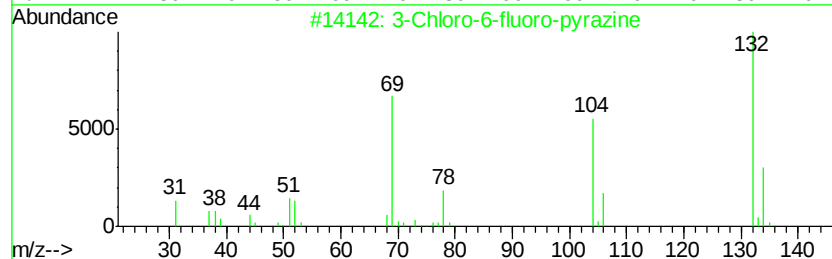
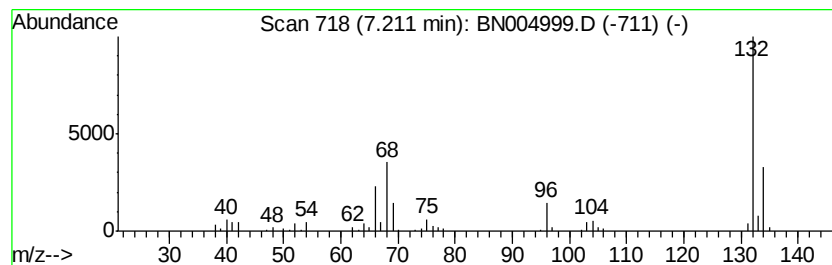
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Peak Number 1 unknown7.21 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.21	99.20 ng	228636	1,4-Dichlorobenzene-d4	7.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		Xenon	132	Xe	007440-63-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown7.21	7.21	99.2	ng	228636	1	7.68	46098	20.0