

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031319\
Data File : BM019149.D
Acq On : 13 Mar 2019 17:03
Operator : JU/SJ
Sample : PB117864BL
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
PB117864BL

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_M\METHODS\8270-BM031119.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	5.257	380	386	405	rBV	3411510	4932640	36.14%	7.402%
2	6.840	648	655	675	rBV	3443355	5428427	39.77%	8.146%
3	7.193	708	715	732	rBV	3533040	5536152	40.56%	8.307%
4	7.657	787	794	804	rBV	802643	1249210	9.15%	1.874%
5	7.975	840	848	861	rBV	2431633	3881204	28.43%	5.824%
6	8.828	986	993	1013	rBV	1766456	3090992	22.64%	4.638%
7	10.445	1260	1268	1290	rBV	888637	1708428	12.52%	2.564%
8	12.922	1683	1689	1711	rBV	5090956	7897383	57.85%	11.850%
9	14.310	1919	1925	1943	rBV2	1545569	2579626	18.90%	3.871%
10	15.810	2174	2180	2208	rBV	4542609	7050201	51.65%	10.579%
11	17.068	2387	2394	2417	rBV	2224112	3398731	24.90%	5.100%
12	19.704	2836	2842	2855	rBV	10911536	13650464	100.00%	20.483%
13	21.268	3102	3108	3122	rBV	2504139	3577831	26.21%	5.369%
14	23.503	3482	3488	3499	rVB	1323564	2661259	19.50%	3.993%

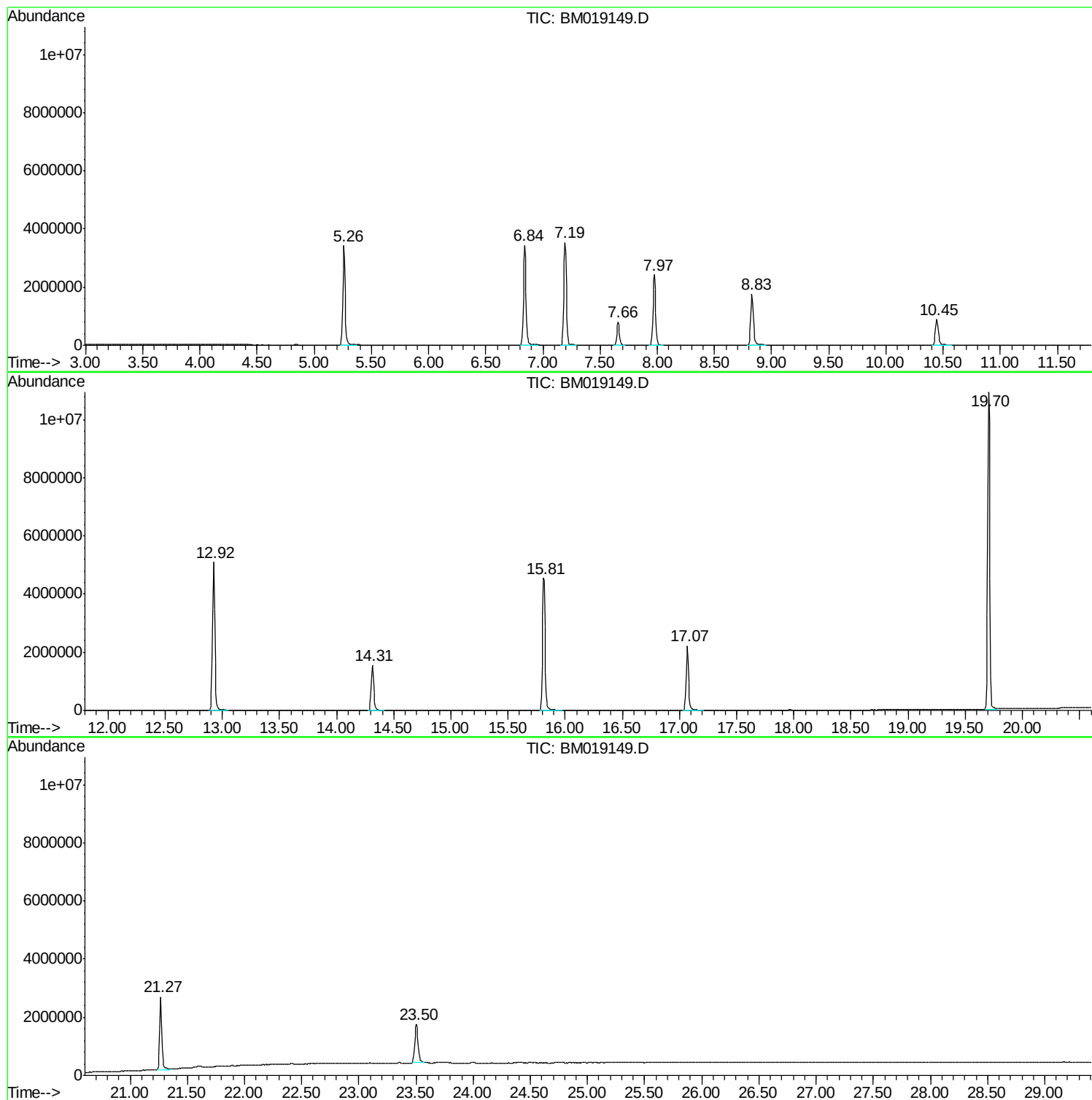
Sum of corrected areas: 66642548

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM031119.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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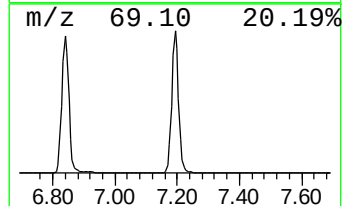
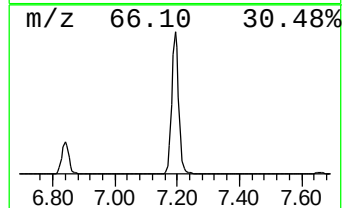
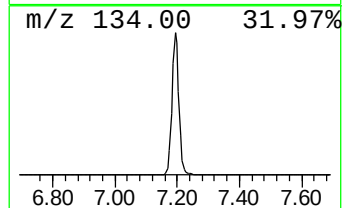
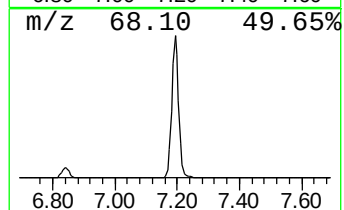
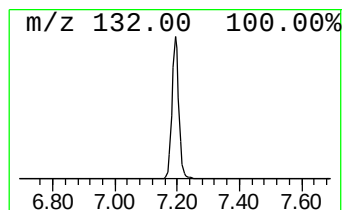
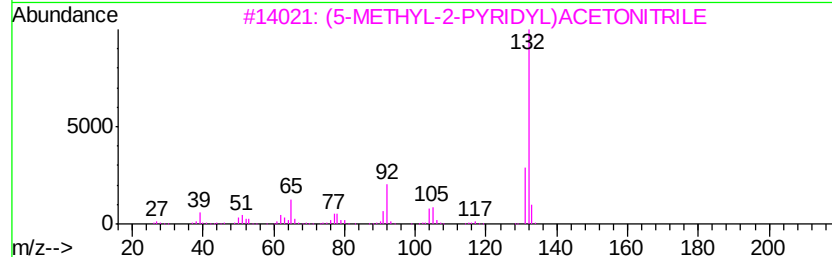
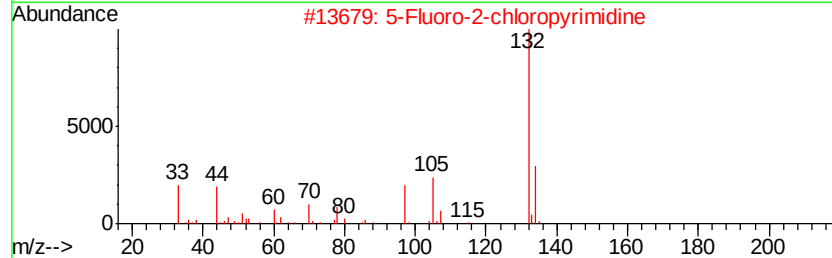
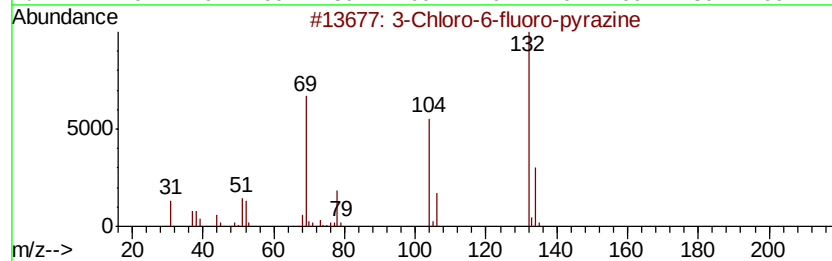
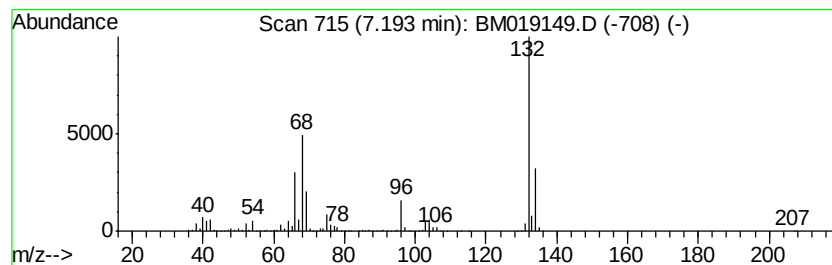
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TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown7.19 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.19	88.63 ng	5536150	1,4-Dichlorobenzene-d4	7.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
4		Benzene, 1,2,4-trifluoro-	132	C6H3F3	000367-23-7	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
unknown7.19	7.19	88.6	ng	5536150	1	7.66	1249210	20.0