

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN110918\
Data File : BN003471.D
Acq On : 09 Nov 2018 14:40
Operator : MJ/SJ
Sample : PB114663BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
PB114663BL

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-BN102418.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	4.963	331	336	344	rBV	60382	82331	1.67%	0.256%
2	5.416	407	413	422	rBV	1702647	2420401	49.02%	7.529%
3	7.010	676	684	695	rBV	1612781	2584455	52.34%	8.039%
4	7.381	739	747	760	rBV	1739671	2758404	55.86%	8.581%
5	7.851	820	827	835	rVB	480437	755127	15.29%	2.349%
6	8.169	874	881	890	rBV	1384336	2167566	43.90%	6.743%
7	9.016	1018	1025	1040	rVB	995444	1619491	32.80%	5.038%
8	10.645	1295	1302	1314	rBV	664322	1093841	22.15%	3.403%
9	13.104	1713	1720	1727	rBV	2974853	4358751	88.27%	13.559%
10	14.486	1948	1955	1963	rBV2	1014089	1510385	30.59%	4.698%
11	15.974	2201	2208	2225	rVB	2100941	2933232	59.40%	9.124%
12	17.227	2415	2421	2433	rVB2	1221661	1659104	33.60%	5.161%
13	19.845	2860	2866	2871	rBV	4033472	4938060	100.00%	15.361%
14	21.404	3126	3131	3139	rBV	1271703	1607723	32.56%	5.001%
15	23.721	3518	3525	3537	rVB	910974	1658461	33.59%	5.159%

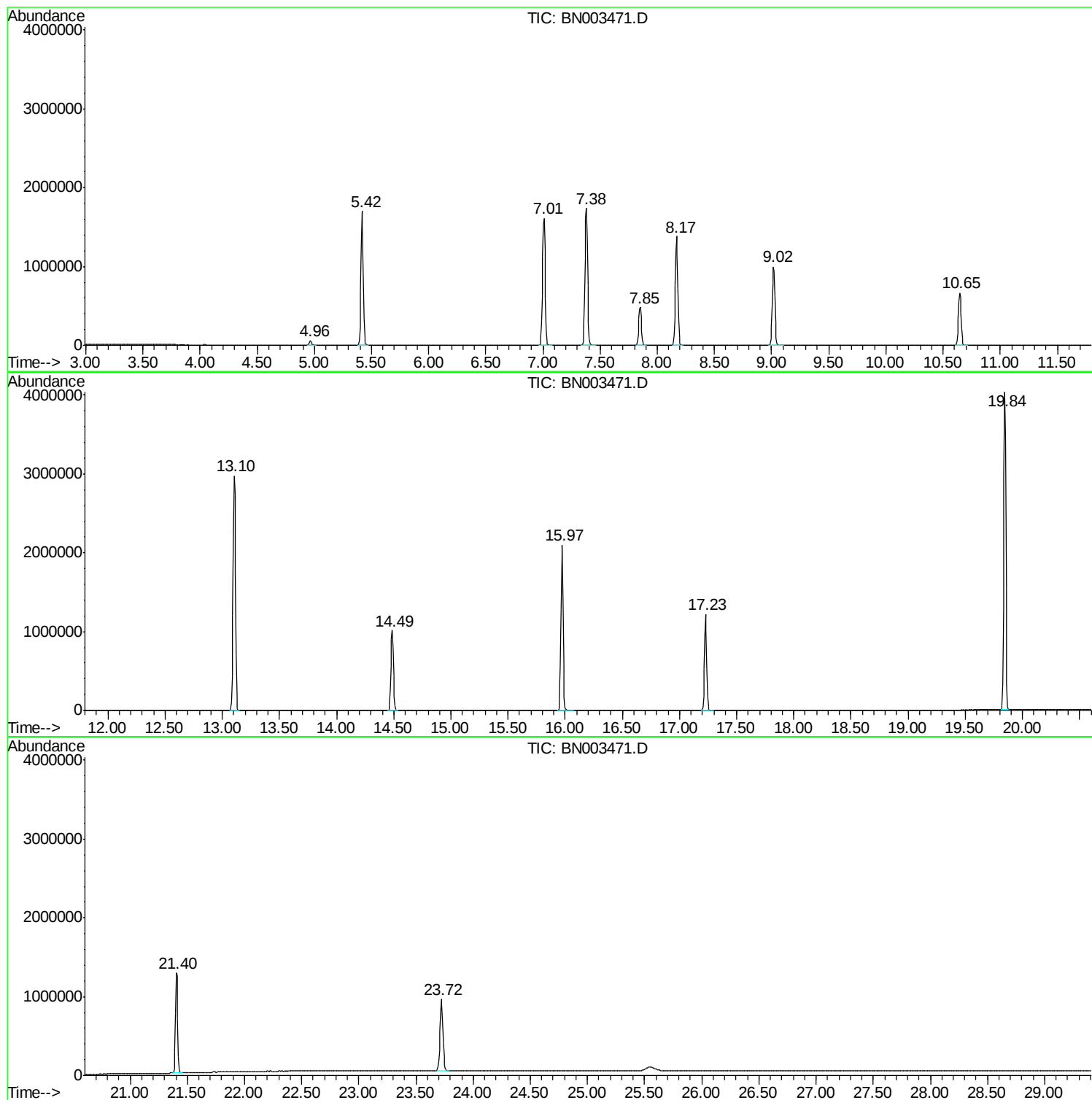
Sum of corrected areas: 32147332

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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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Instrument :
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ClientSampled :
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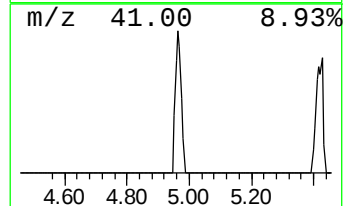
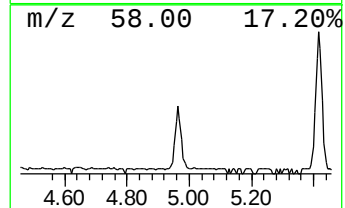
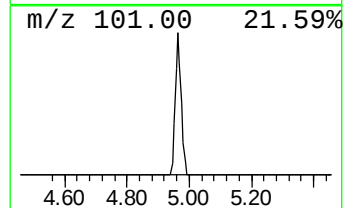
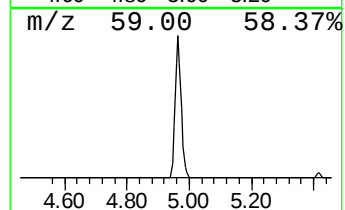
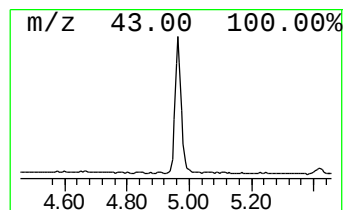
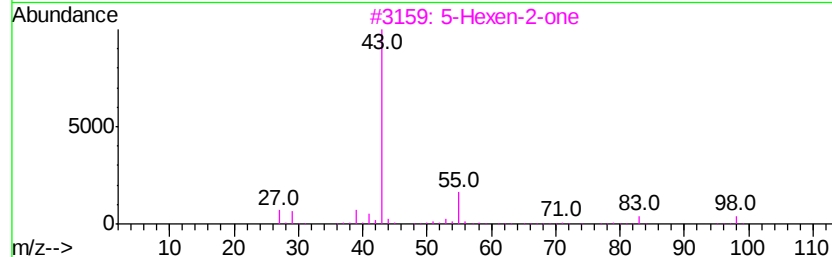
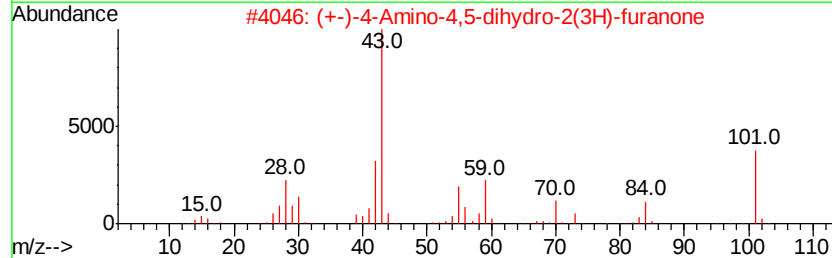
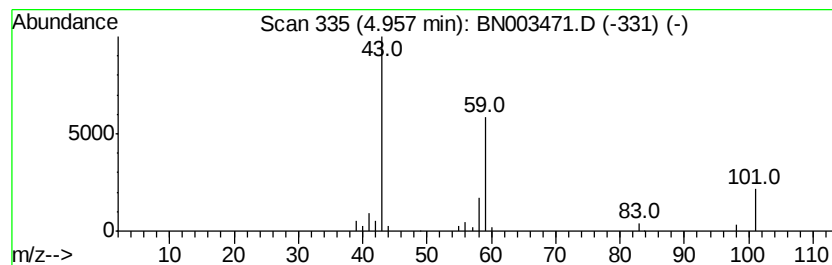
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.96	2.18 ng	82331	1,4-Dichlorobenzene-d4	7.85

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			(+)-4-Amino-4,5-dihydro-2(3H)-furanone	101	C4H7NO2	016504-58-8	9
3			5-Hexen-2-one	98	C6H10O	000109-49-9	9
4			Acetone	58	C3H6O	000067-64-1	9
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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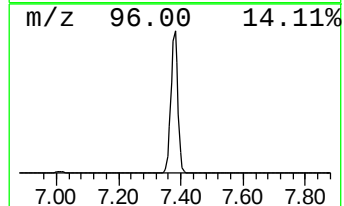
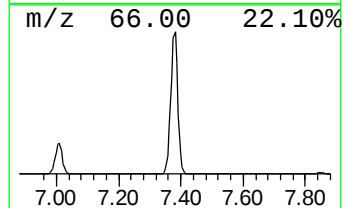
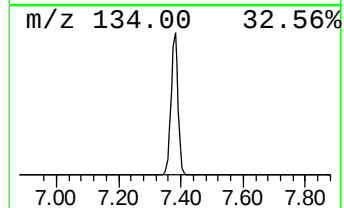
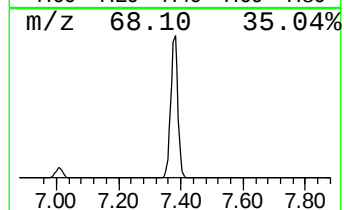
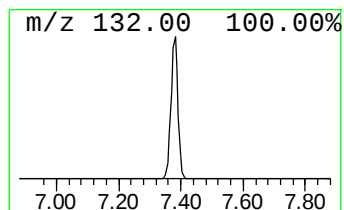
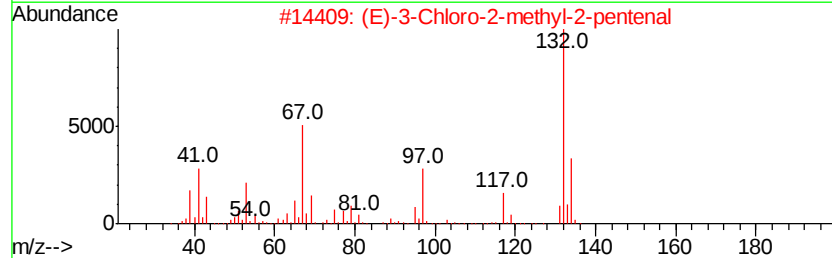
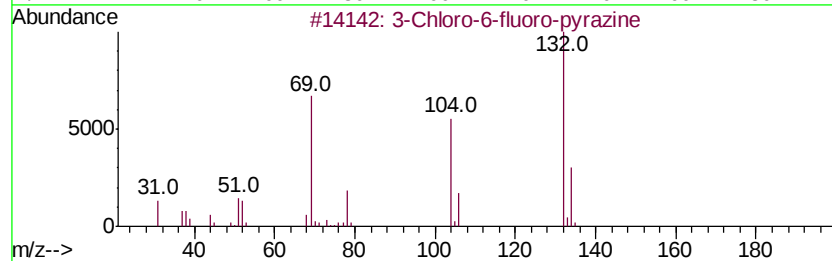
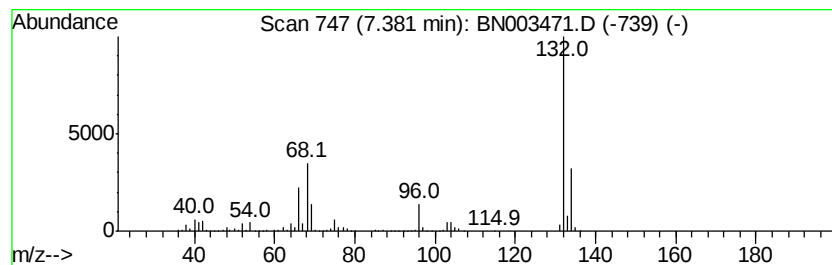
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Peak Number 2 unknown7.38 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.38	73.06 ng	2758400	1,4-Dichlorobenzene-d4	7.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	16
4		Amphetaminil	250	C17H18N2	017590-01-1	12
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.96	2.2	ng	82331	1	7.85	755127	20.0
unknown7.38	7.38	73.1	ng	2758400	1	7.85	755127	20.0