

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN072618\
 Data File : BN002174.D
 Acq On : 26 Jul 2018 16:07
 Operator : JU/SJ
 Sample : PB111396BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB111396BL

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-BN072518.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	4.857	313	318	331	rBV	785168	1126261	10.19%	1.432%
2	5.310	389	395	414	rVB	5594977	7981806	72.23%	10.147%
3	6.881	653	662	675	rBV	5182340	8399793	76.01%	10.679%
4	7.234	714	722	733	rBV	5441135	8510108	77.01%	10.819%
5	7.693	793	800	810	rVB	1012388	1602405	14.50%	2.037%
6	8.010	846	854	863	rBV	4020079	6416564	58.06%	8.157%
7	8.840	988	995	1011	rBV	3144252	5298898	47.95%	6.737%
8	10.463	1263	1271	1282	rBV	1298943	2149795	19.45%	2.733%
9	12.939	1685	1692	1707	rBV	7416419	11050685	100.00%	14.049%
10	14.322	1920	1927	1937	rBV2	1685505	2533656	22.93%	3.221%
11	15.816	2174	2181	2198	rBV	4794559	6940333	62.80%	8.823%
12	17.069	2387	2394	2404	rBV2	1726120	2484127	22.48%	3.158%
13	19.710	2837	2843	2848	rBV	8065935	10172875	92.06%	12.933%
14	21.268	3102	3108	3114	rBV	1612582	2046199	18.52%	2.601%
15	23.492	3479	3486	3497	rVB2	1028429	1945791	17.61%	2.474%

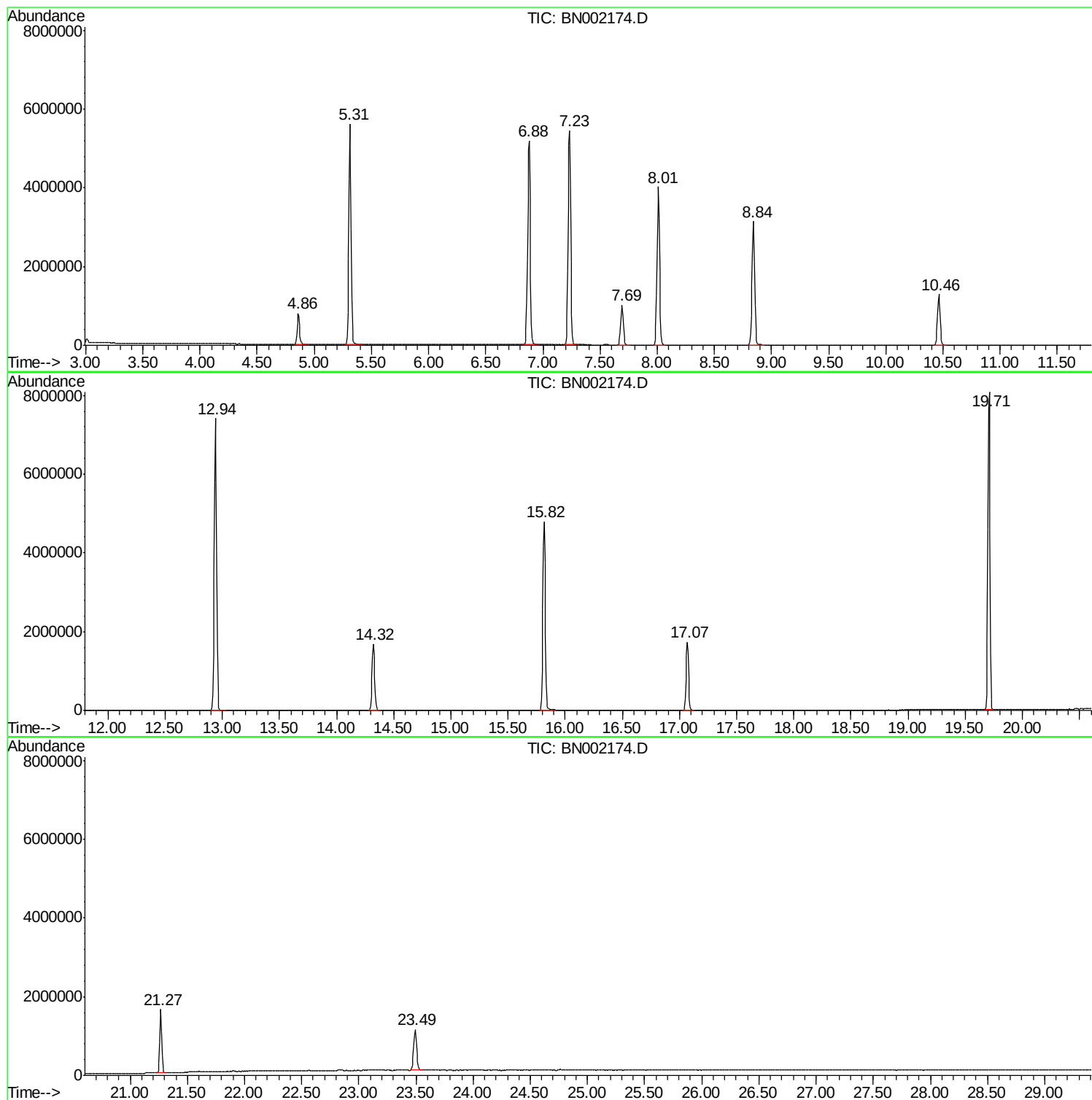
Sum of corrected areas: 78659296

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-BN072518.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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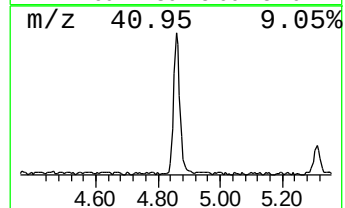
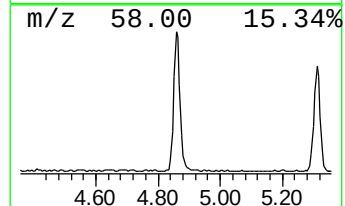
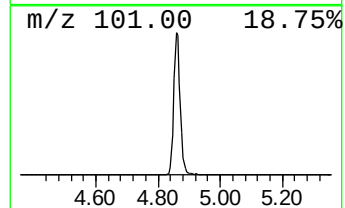
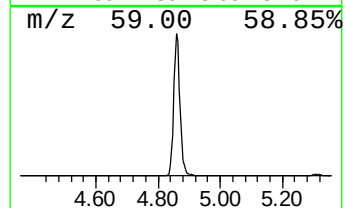
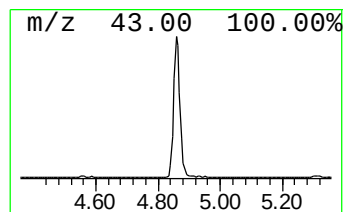
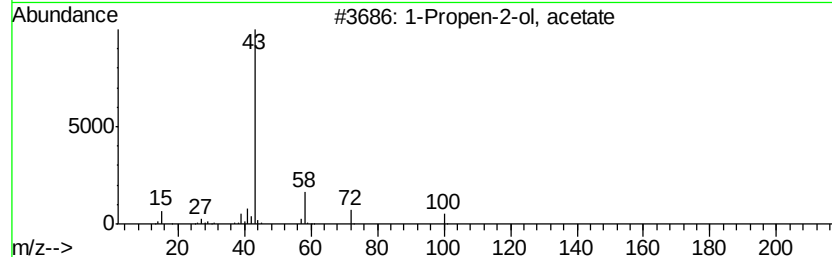
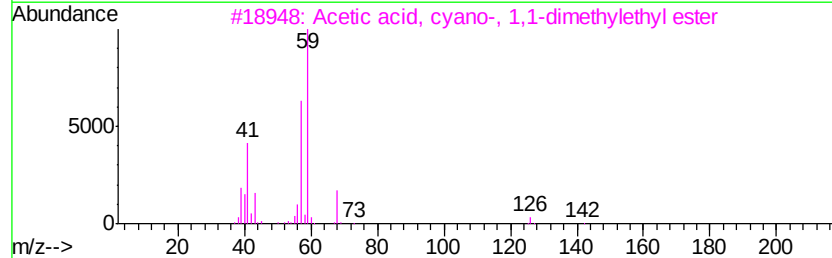
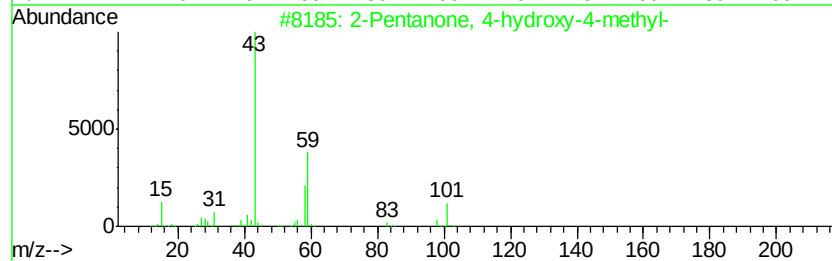
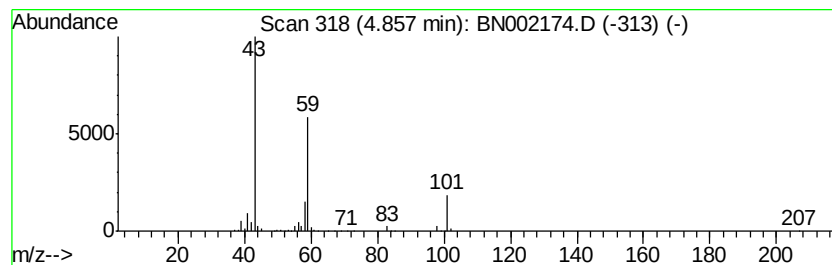
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.86	14.06 ng	1126260	1,4-Dichlorobenzene-d4	7.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
3		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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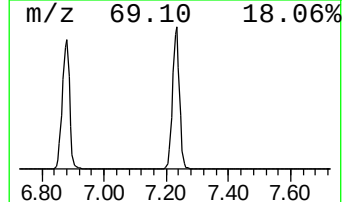
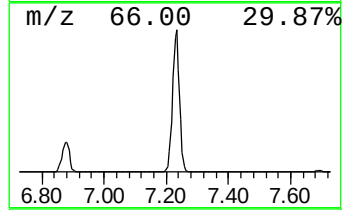
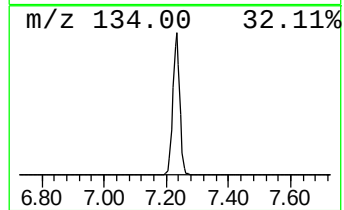
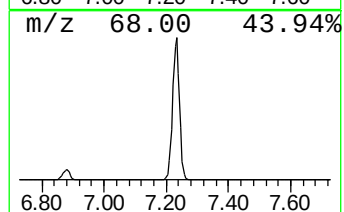
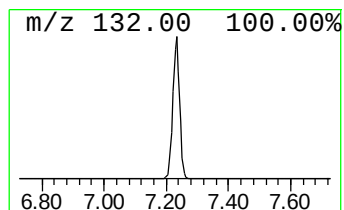
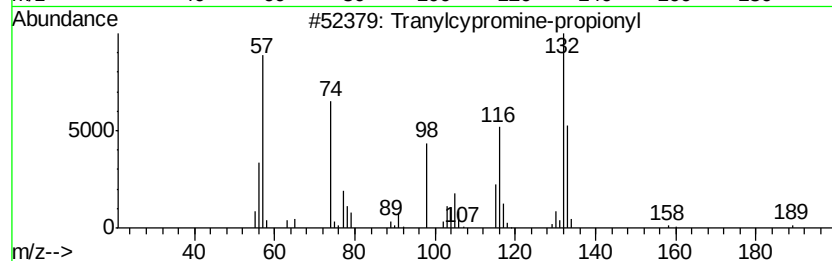
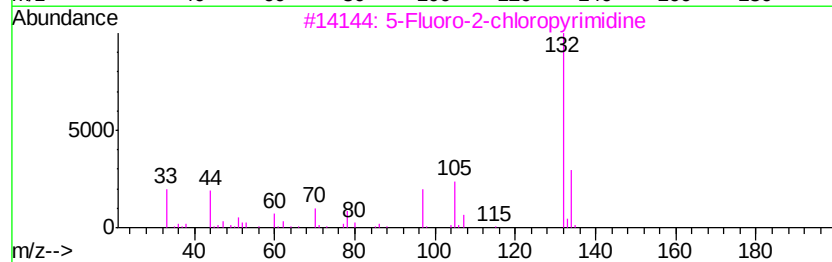
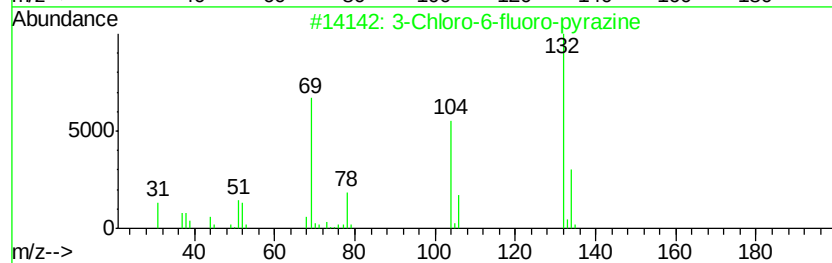
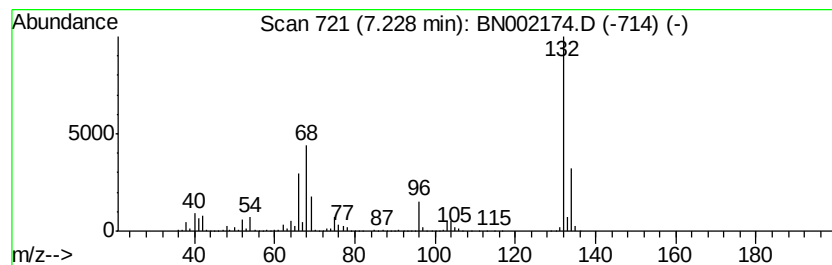
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Peak Number 2 unknown7.23 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.23	106.22 ng	8510110	1,4-Dichlorobenzene-d4	7.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		Xenon	132	Xe	007440-63-3	9
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.86	14.1	ng	1126260	1	7.69	1602410	20.0
unknown7.23	7.23	106.2	ng	8510110	1	7.69	1602410	20.0