

Data Path : Z:\HPCHEM1\BNA N\DATA\BN020618\
 Data File : BN000023.D
 Acq On : 06 Feb 2018 09:36
 Operator :
 Sample : PB106290BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB106290BL

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:\HPCHEM1\BNA_N\METHODS\8270-BN020118.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	3.075	10	15	33	rVB	270913	495911	1.39%	0.261%
2	5.952	496	504	522	rBV	9160461	15060498	42.15%	7.913%
3	7.599	775	784	800	rBV	9937499	16464079	46.07%	8.650%
4	7.993	843	851	863	rBV	9695973	16281685	45.56%	8.554%
5	8.475	923	933	942	rBV	1841754	3053769	8.55%	1.604%
6	8.804	981	989	1000	rVB	7108911	11969956	33.50%	6.289%
7	9.652	1123	1133	1154	rBV	5680755	9386241	26.27%	4.931%
8	11.316	1408	1416	1425	rBV	2841448	4835029	13.53%	2.540%
9	13.716	1817	1824	1839	rBV	16750203	25024263	70.03%	13.148%
10	15.081	2049	2056	2065	rBV2	4370970	6767232	18.94%	3.555%
11	16.551	2298	2306	2335	rBV	11067330	16005337	44.79%	8.409%
12	17.804	2512	2519	2532	rBV2	5855971	8546551	23.92%	4.490%
13	20.369	2949	2955	2972	rBV	27156011	35734122	100.00%	18.774%
14	21.927	3214	3220	3234	rBV2	7672608	10279555	28.77%	5.401%
15	24.427	3637	3645	3662	rVB2	4885536	10429410	29.19%	5.480%

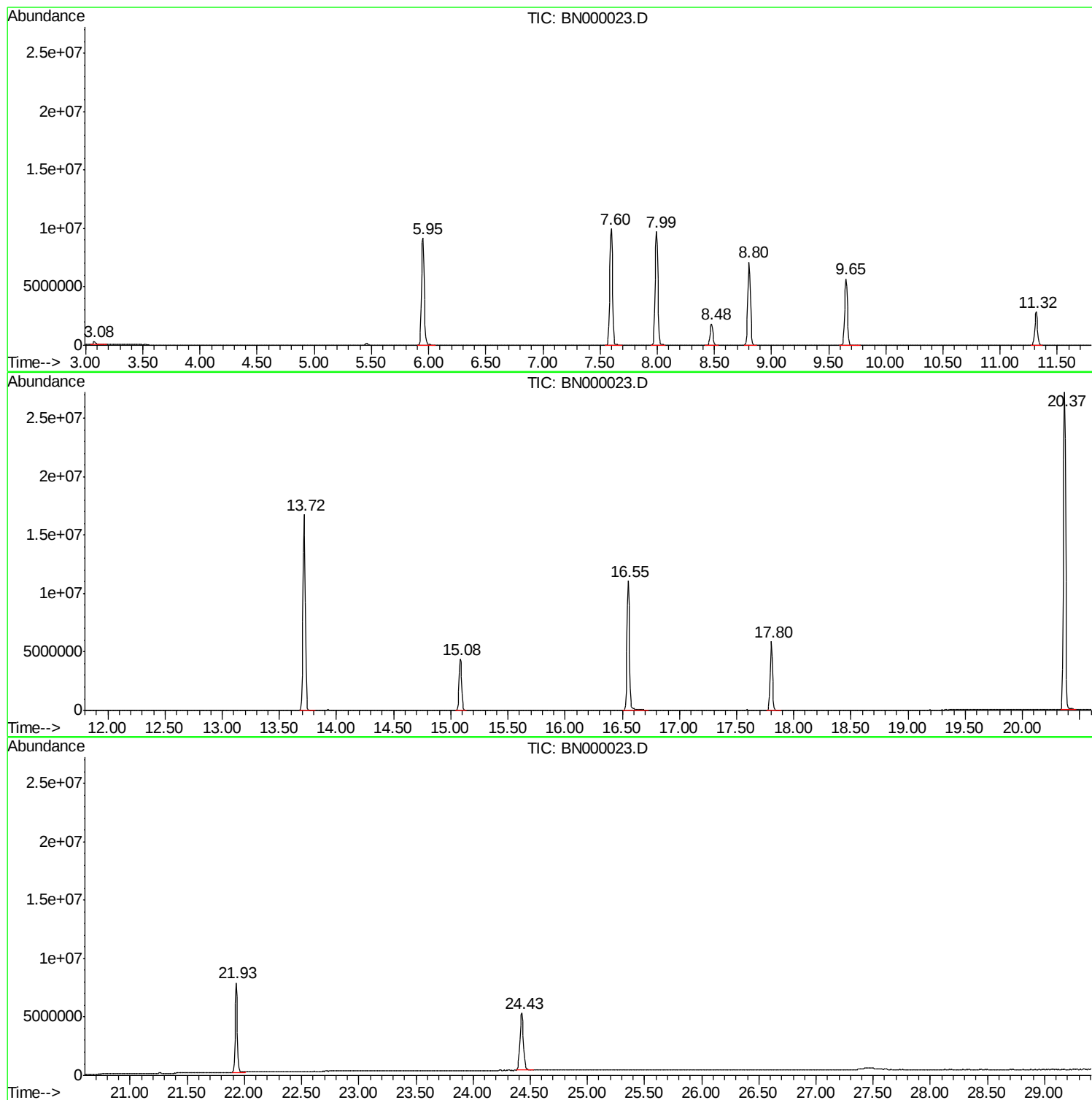
Sum of corrected areas: 190333638

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Quant Method : Z:\HPCHEM1\BNA N\METHODS\8270-BN020118.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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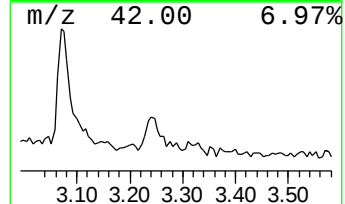
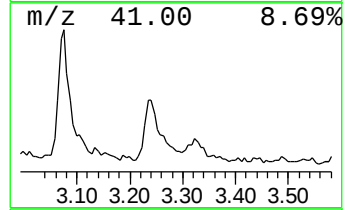
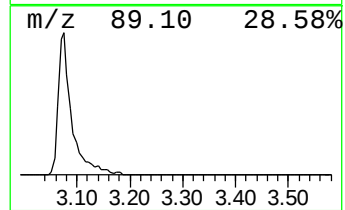
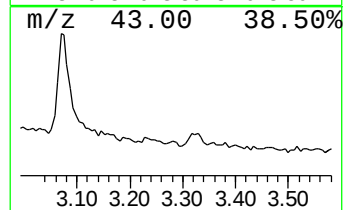
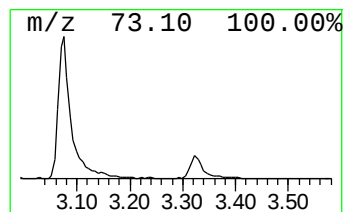
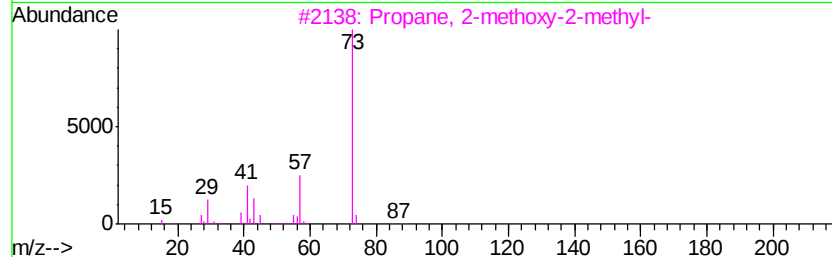
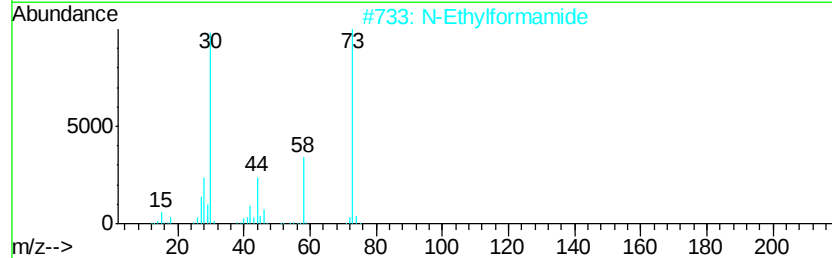
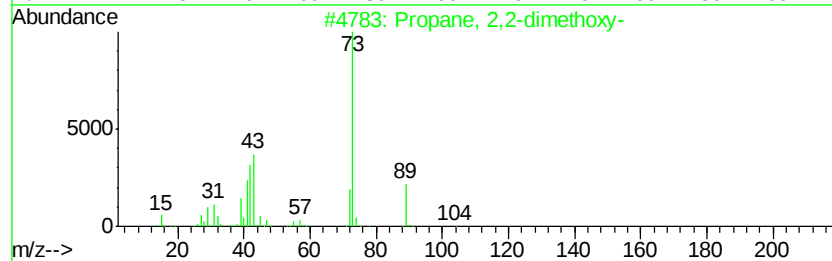
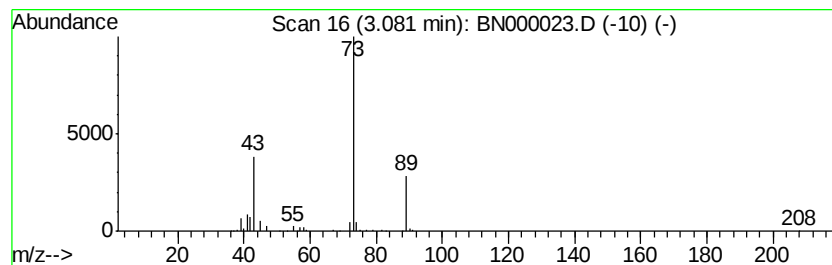
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Peak Number 1 Propane, 2,2-dimethoxy- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
3.08	3.25 ng	495911	1,4-Dichlorobenzene-d4	8.48	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	50
2	N-Ethylformamide	73	C3H7NO	000627-45-2	37
3	Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	12
4	1,3-Dioxolane, 2-pentadecyl-	284	C18H36O2	004360-57-0	9
5	2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	9



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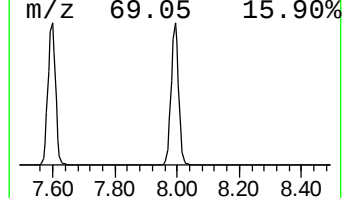
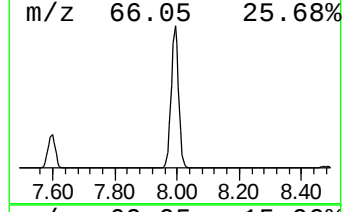
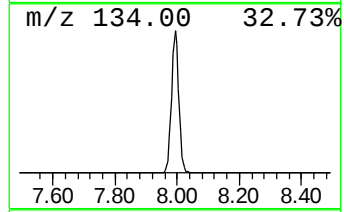
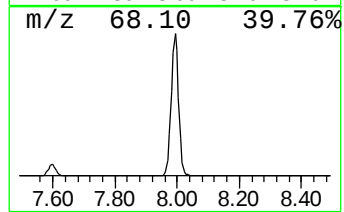
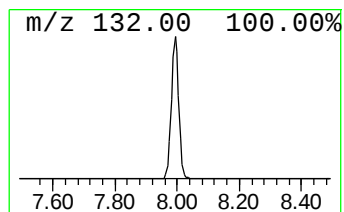
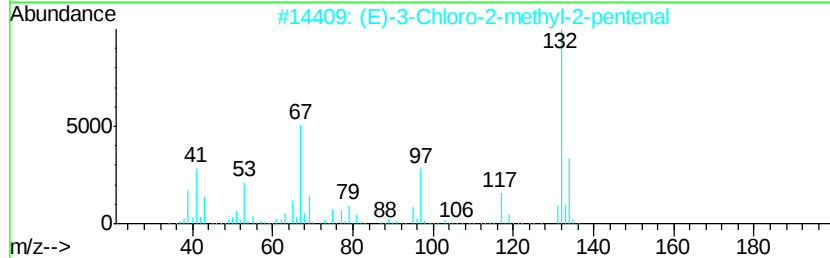
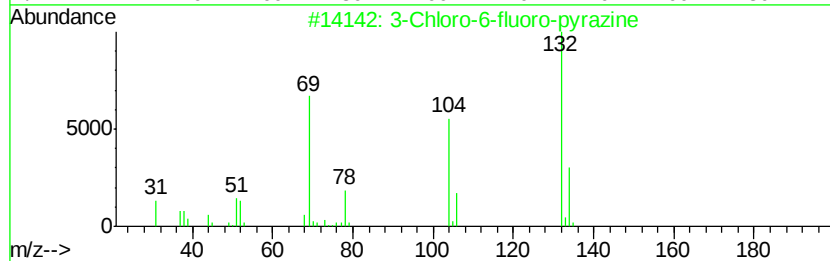
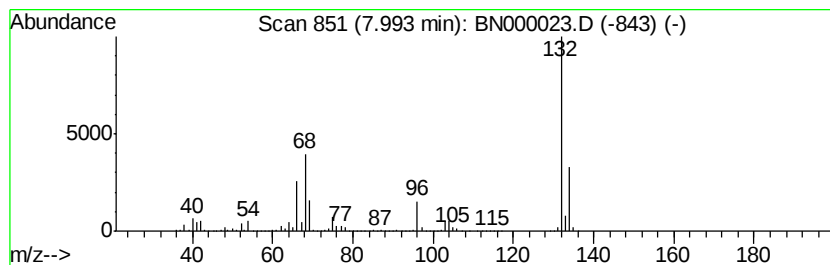
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Peak Number 2 unknown7.99 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.99	106.63 ng	16281700	1,4-Dichlorobenzene-d4	8.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
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
Propane, 2,2-dime...	3.08	3.3	ng	495911	1	8.48	3053770	20.0
unknown7.99	7.99	106.6	ng	16281700	1	8.48	3053770	20.0