

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG121217\
 Data File : BG031519.D
 Acq On : 13 Dec 2017 8:41
 Operator : SJ/JU
 Sample : I6763-09
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 120517-TIDO-F-U-B

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_G\METHODS\8270-BG121117.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	5.185	318	323	332	rBV	58684	89939	1.46%	0.435%
2	5.672	400	406	421	rBV	329232	593045	9.62%	2.868%
3	7.282	673	680	698	rBV	207005	437957	7.10%	2.118%
4	7.646	734	742	763	rBV	606237	1114972	18.08%	5.393%
5	8.111	814	821	829	rBV	191029	336351	5.45%	1.627%
6	8.434	867	876	888	rBV	832885	1492915	24.21%	7.221%
7	9.280	1011	1020	1033	rBV	568012	1125949	18.26%	5.446%
8	10.925	1292	1300	1320	rBV	223419	494722	8.02%	2.393%
9	13.352	1706	1713	1735	rBV	2019831	3341855	54.19%	16.163%
10	14.732	1939	1948	1961	rVB2	460654	785860	12.74%	3.801%
11	16.219	2196	2201	2222	rBV	438322	875340	14.19%	4.234%
12	17.476	2408	2415	2430	rBV2	653193	1069735	17.34%	5.174%
13	20.067	2849	2856	2864	rBV	4515285	6167450	100.00%	29.829%
14	21.748	3134	3142	3151	rBV	800191	1413704	22.92%	6.837%
15	25.032	3690	3701	3715	rBV2	458058	1336135	21.66%	6.462%

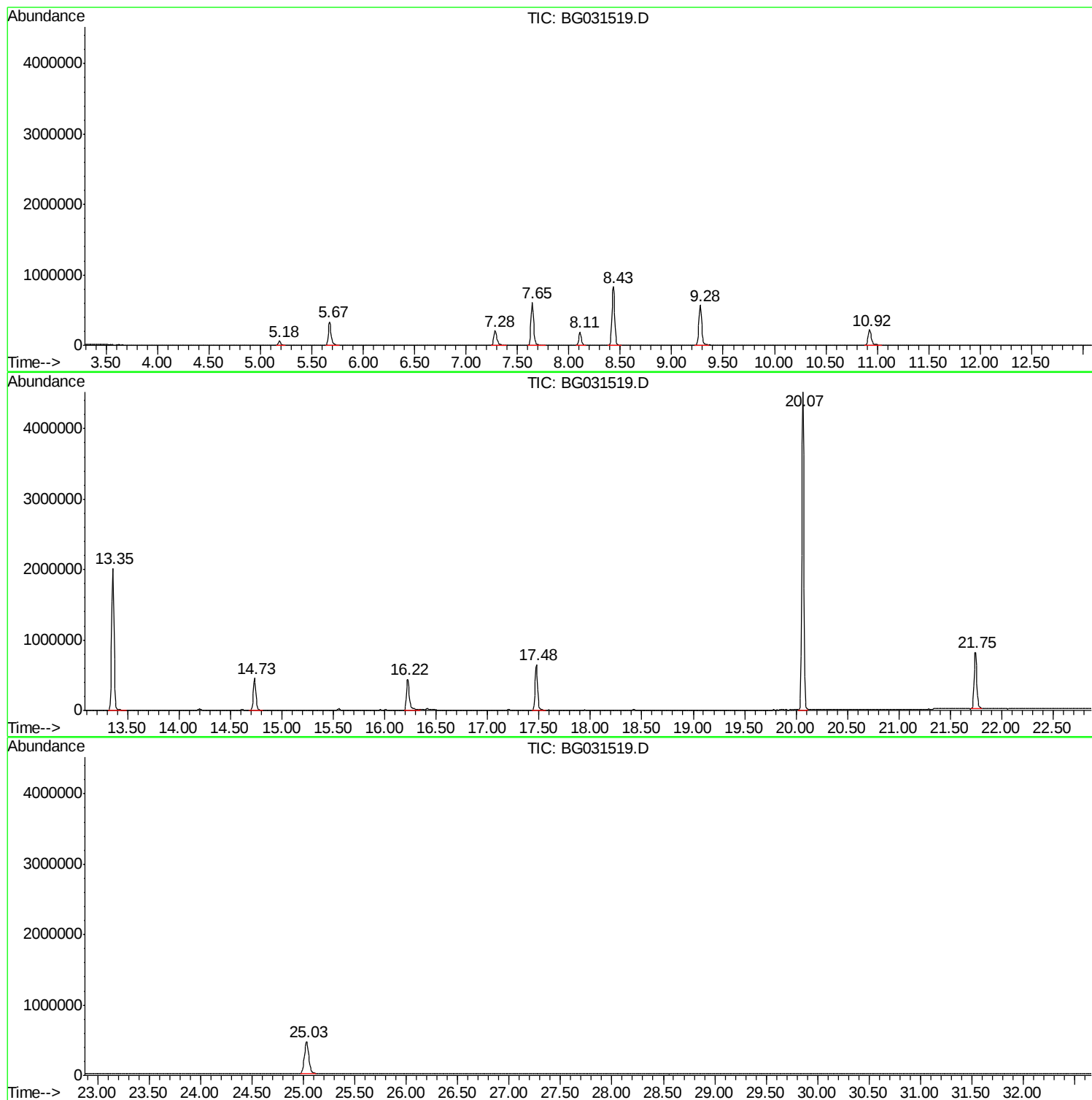
Sum of corrected areas: 20675929

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Instrument :
BNA_G
ClientSampleId :
120517-TIDO-F-U-B

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG121117.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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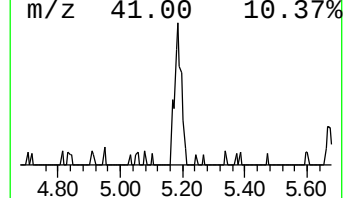
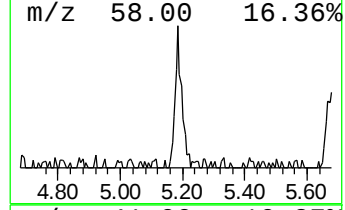
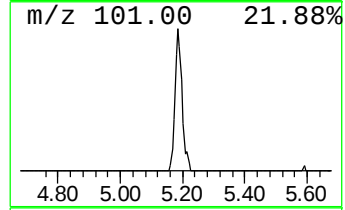
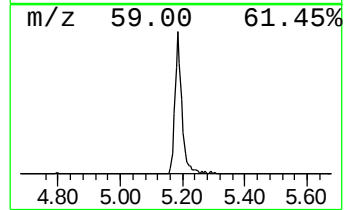
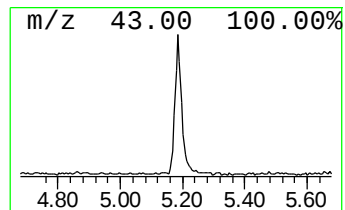
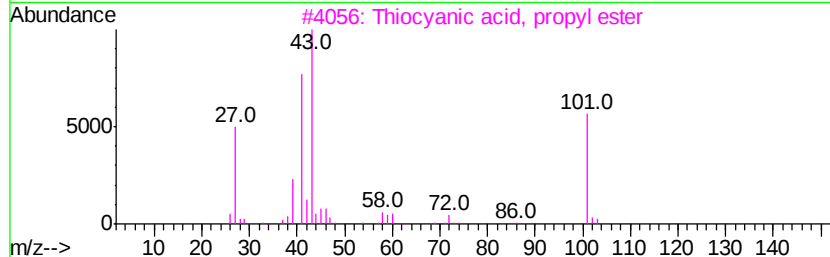
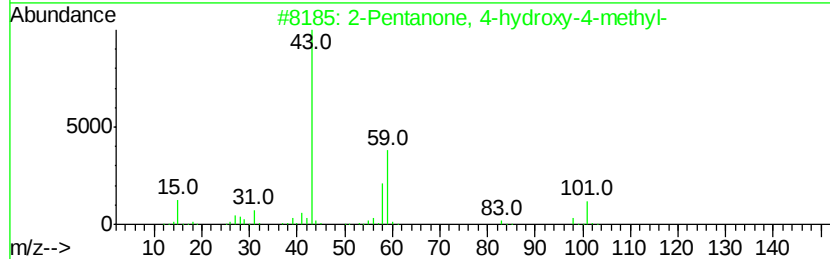
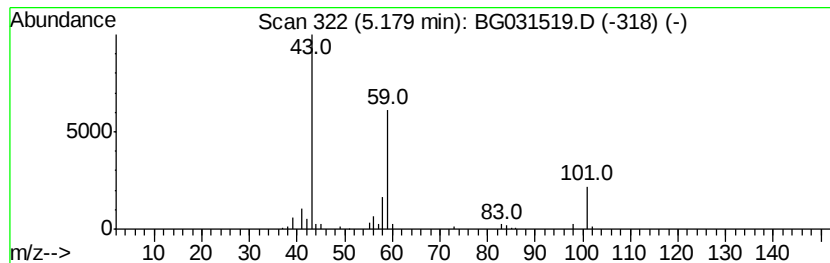
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	5.35 ng	89939	1,4-Dichlorobenzene-d4	8.11

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3			Thiocyanic acid, propyl ester	101	C4H7NS	004251-16-5	9
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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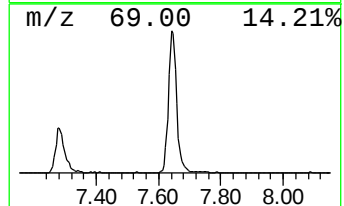
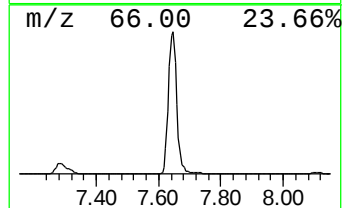
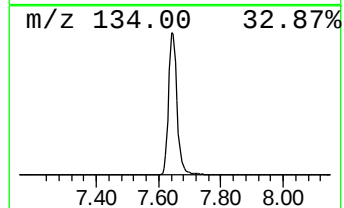
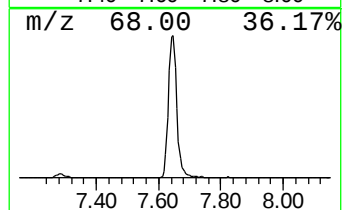
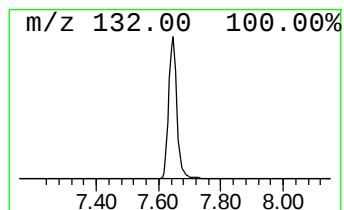
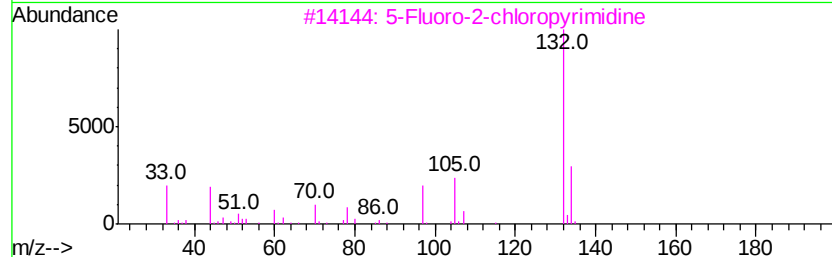
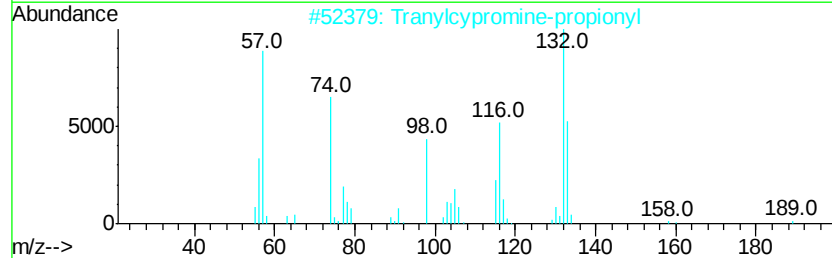
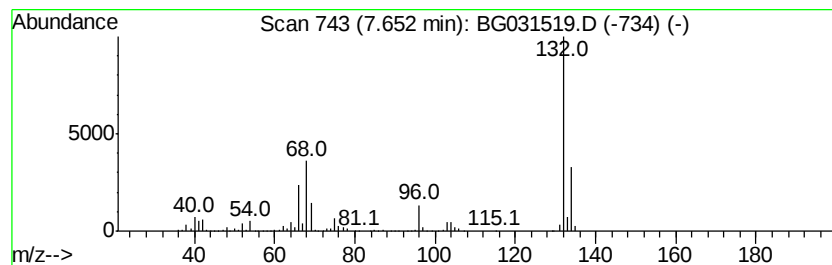
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Peak Number 2 unknown7.65 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.65	66.30 ng	1114970	1,4-Dichlorobenzene-d4	8.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.18	5.3	ng	89939	1	8.11	336351	20.0
unknown7.65	7.65	66.3	ng	1114970	1	8.11	336351	20.0