

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG121217\
 Data File : BG031489.D
 Acq On : 12 Dec 2017 13:15
 Operator : SJ/JU
 Sample : PB104888BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB104888BL

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	5.179	314	322	332	rBV	139606	221134	4.09%	1.166%
2	5.667	399	405	420	rBV	528516	945043	17.49%	4.984%
3	7.277	668	679	692	rBV	525233	983030	18.19%	5.185%
4	7.641	734	741	762	rVB	562674	1001995	18.54%	5.285%
5	8.111	813	821	832	rBV	161199	285704	5.29%	1.507%
6	8.434	868	876	885	rBV	665601	1169989	21.65%	6.171%
7	9.281	1012	1020	1033	rBV	397840	829095	15.34%	4.373%
8	10.926	1293	1300	1320	rBV	165951	394684	7.30%	2.082%
9	13.352	1706	1713	1725	rBV	1346963	2334649	43.21%	12.313%
10	14.733	1940	1948	1964	rBV2	310233	593586	10.99%	3.131%
11	16.219	2194	2201	2219	rBV	664304	1231800	22.80%	6.497%
12	17.477	2408	2415	2431	rBV2	540153	925283	17.12%	4.880%
13	19.698	2788	2793	2797	rBV	37831	57663	1.07%	0.304%
14	20.068	2850	2856	2867	rBV	4011588	5403598	100.00%	28.499%
15	21.748	3134	3142	3158	rBV2	707175	1359772	25.16%	7.172%
16	25.033	3689	3701	3714	rBV3	388387	1223733	22.65%	6.454%

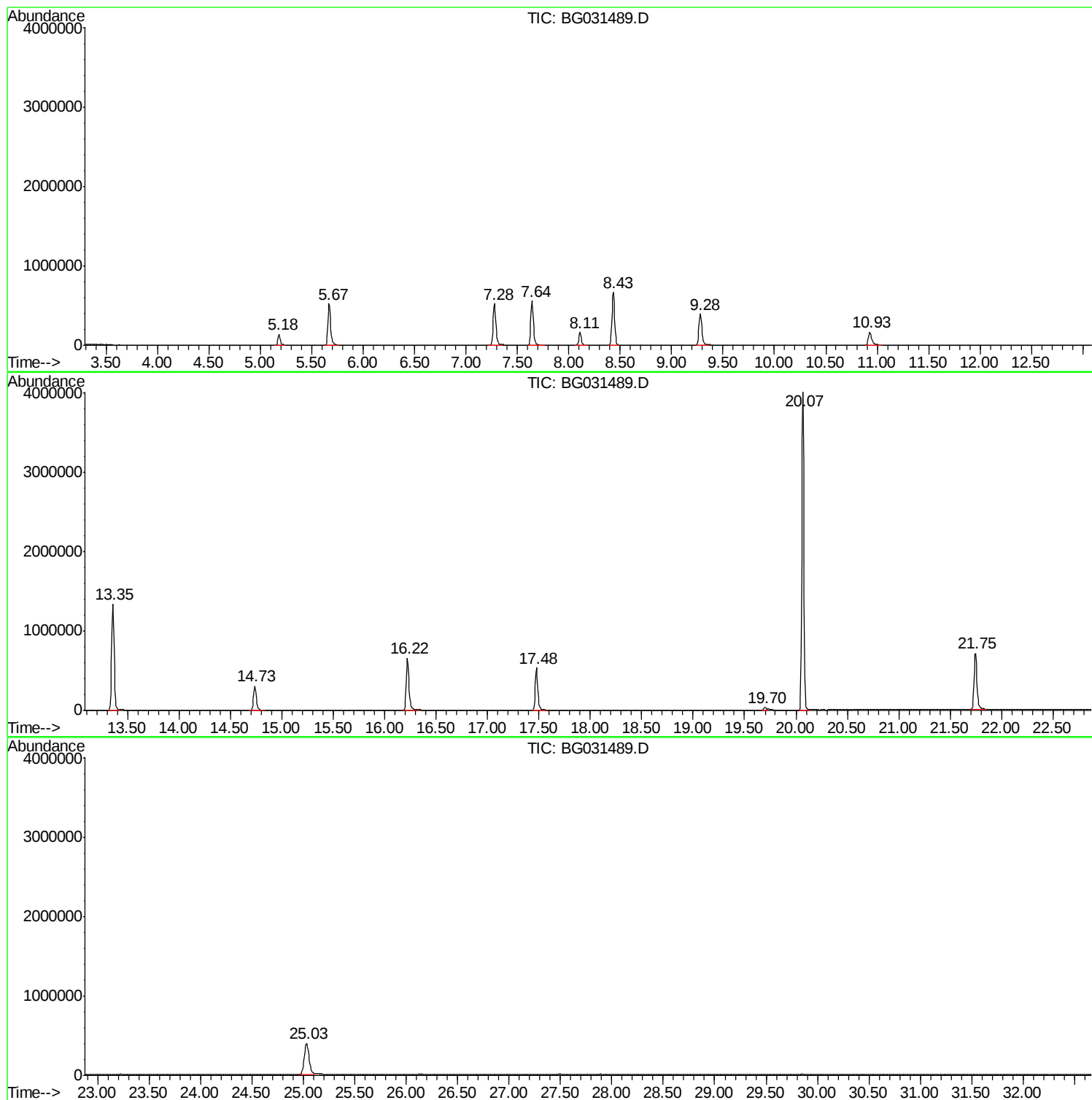
Sum of corrected areas: 18960758

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



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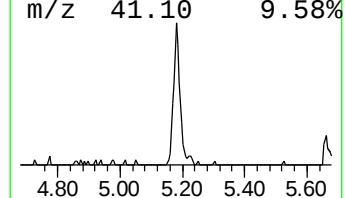
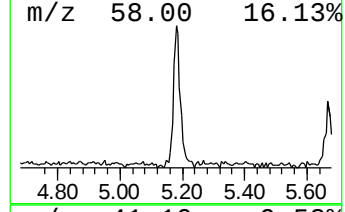
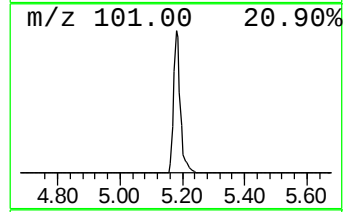
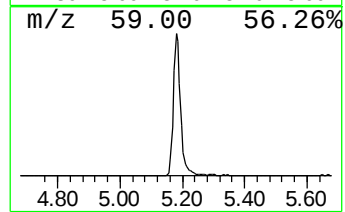
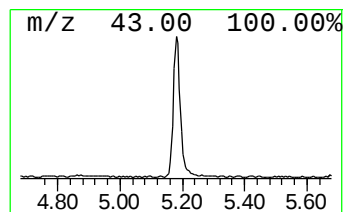
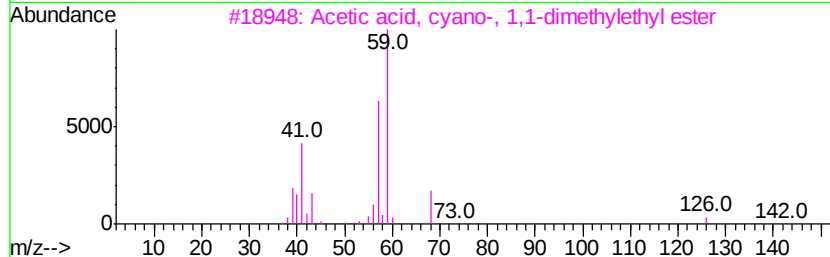
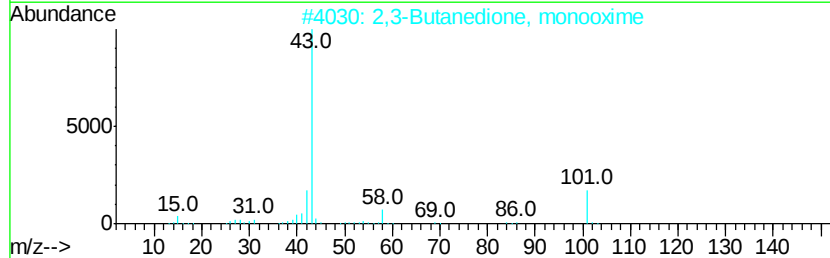
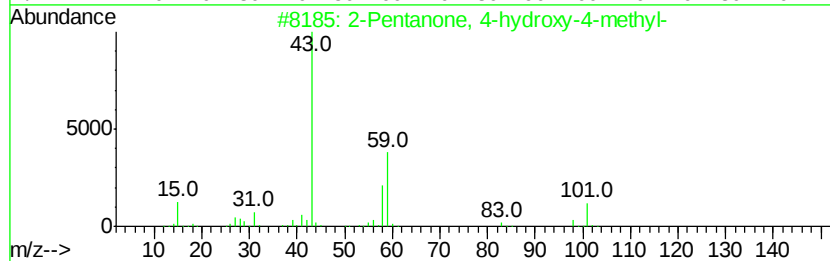
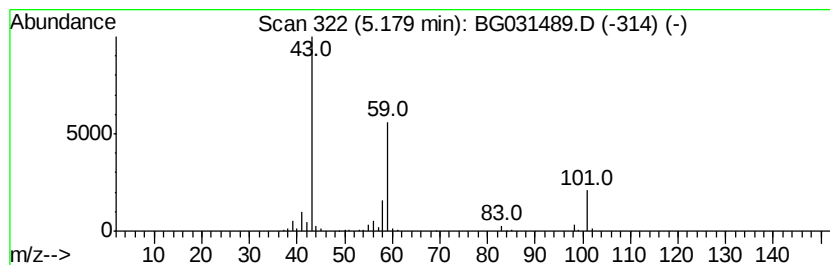
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	15.48 ng	221134	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	45
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
4			Butane	58	C4H10	000106-97-8	9
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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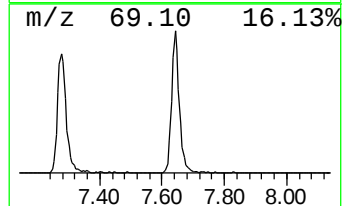
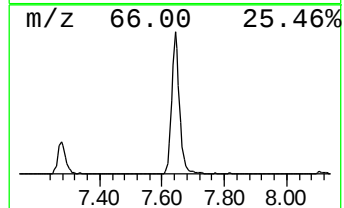
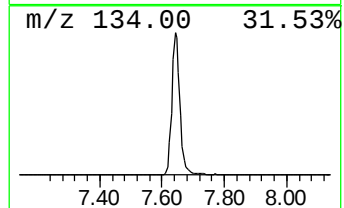
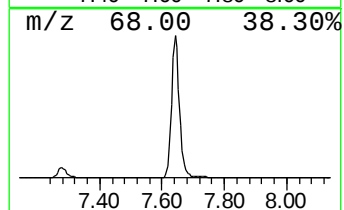
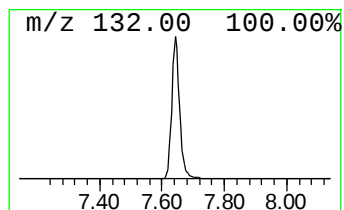
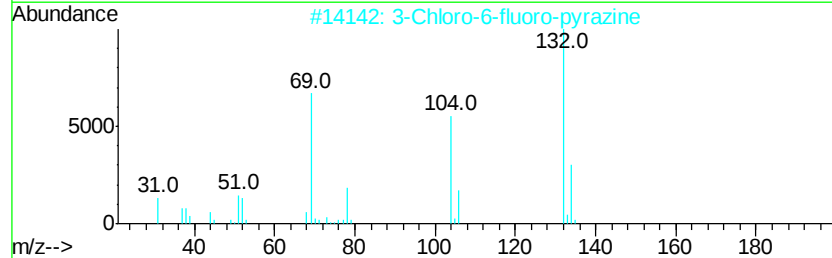
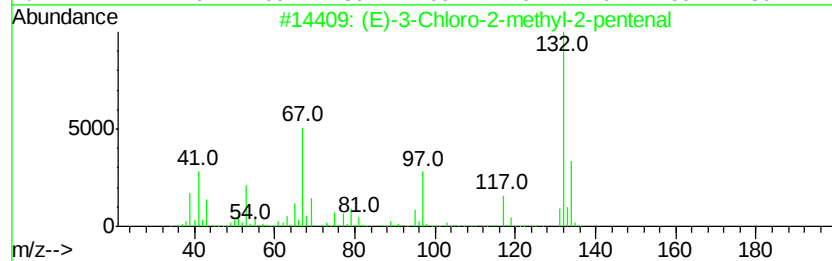
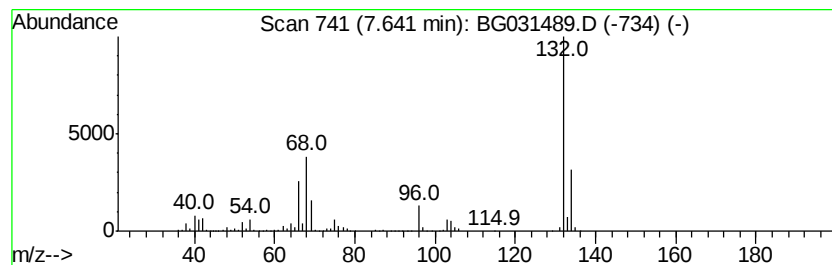
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Peak Number 2 unknown7.64 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.64	70.14 ng	1002000	1,4-Dichlorobenzene-d4	8.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.18	15.5	ng	221134	1	8.11	285704	20.0
unknown7.64	7.64	70.1	ng	1002000	1	8.11	285704	20.0