

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG121117\
 Data File : BG031472.D
 Acq On : 12 Dec 2017 2:35
 Operator : SJ/JU
 Sample : I6762-08
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 120417-TIDO-F-U-S

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	5.181	318	322	330	rBV	42254	73020	1.25%	0.385%
2	5.669	399	405	416	rBV	251852	464362	7.93%	2.446%
3	7.279	673	679	698	rBV	149364	312292	5.34%	1.645%
4	7.643	734	741	755	rBV	493448	891192	15.22%	4.694%
5	8.107	813	820	829	rBV	162569	291525	4.98%	1.535%
6	8.431	867	875	891	rBV	677558	1237533	21.14%	6.518%
7	9.277	1012	1019	1032	rBV	479488	947502	16.19%	4.990%
8	10.928	1292	1300	1318	rBV	178519	413291	7.06%	2.177%
9	13.354	1706	1713	1725	rBV	1639593	2724537	46.54%	14.350%
10	14.735	1941	1948	1964	rVB2	366523	667998	11.41%	3.518%
11	16.216	2194	2200	2217	rBV	862331	1456555	24.88%	7.671%
12	17.473	2406	2414	2429	rBV2	616680	1042957	17.82%	5.493%
13	20.064	2849	2855	2871	rBV	4259660	5853635	100.00%	30.830%
14	21.744	3133	3141	3156	rBV	773387	1380654	23.59%	7.272%
15	25.017	3687	3698	3714	rBV2	395919	1229899	21.01%	6.478%

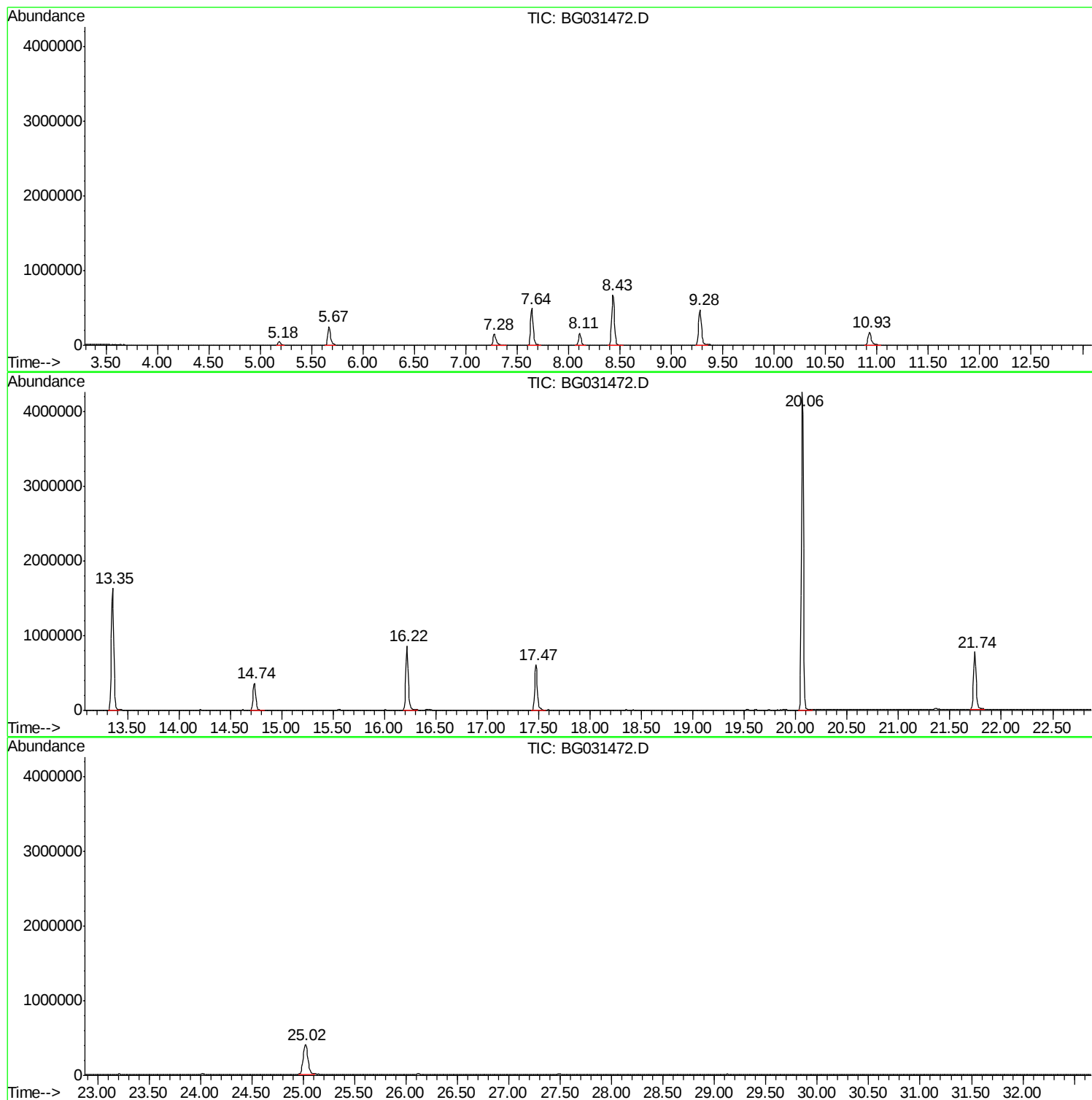
Sum of corrected areas: 18986952

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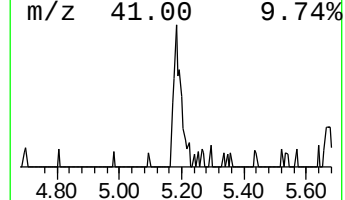
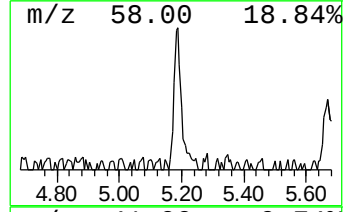
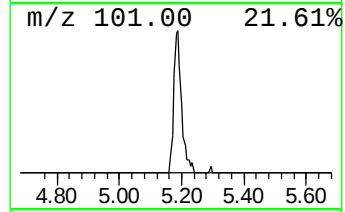
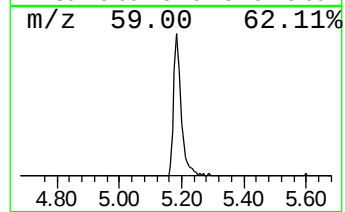
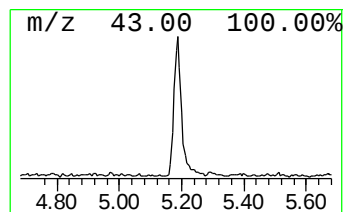
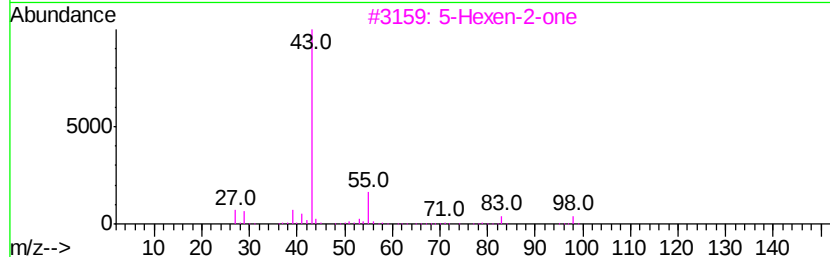
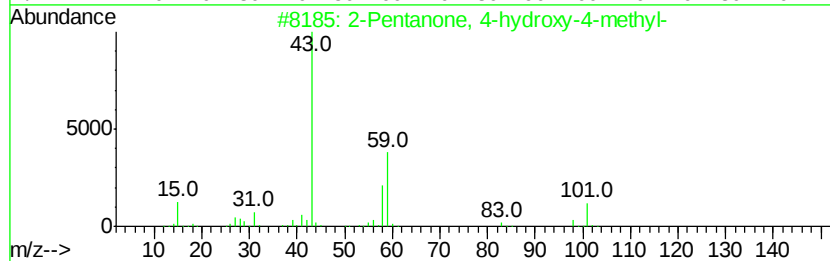
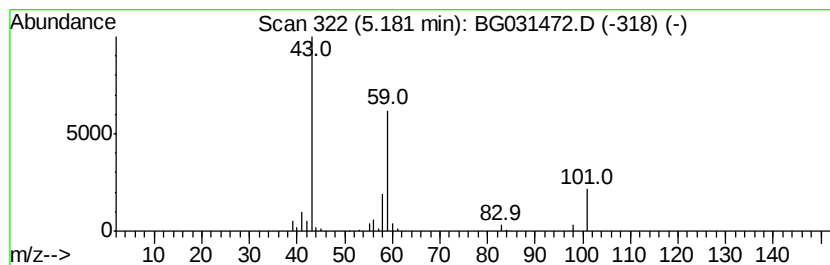
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	5.01 ng	73020	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	40
2			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
3			5-Hexen-2-one	98	C6H10O	000109-49-9	9
4			Acetone	58	C3H6O	000067-64-1	9
5			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9



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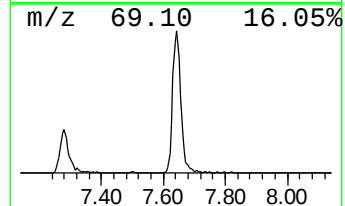
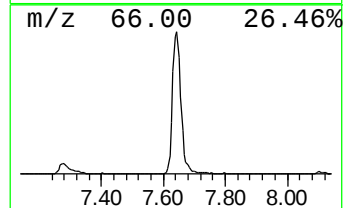
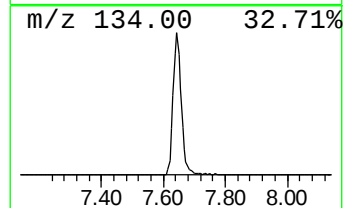
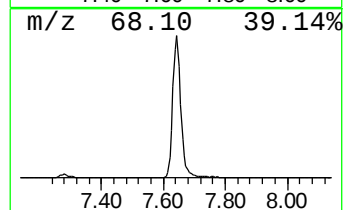
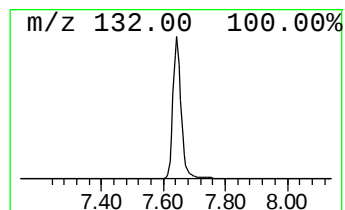
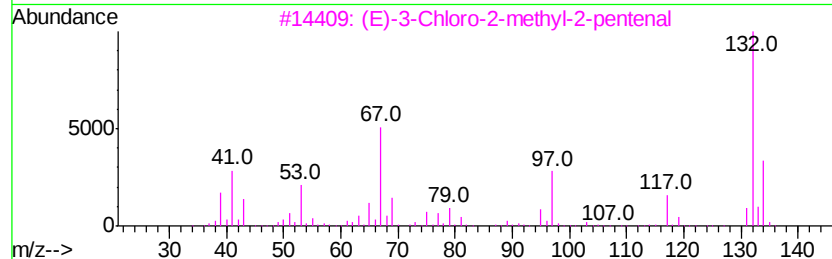
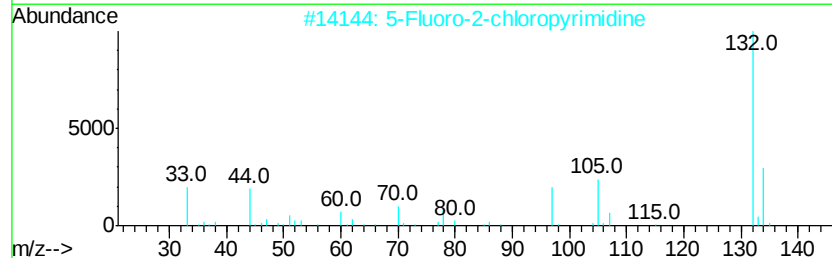
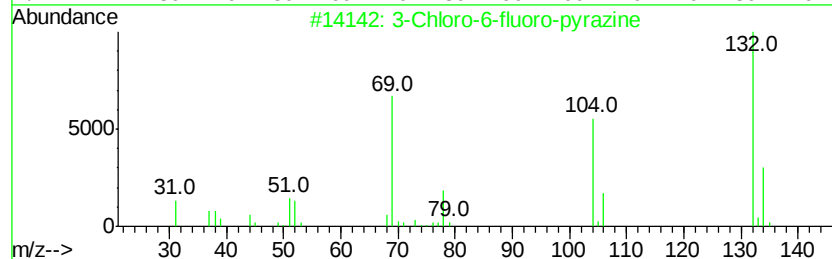
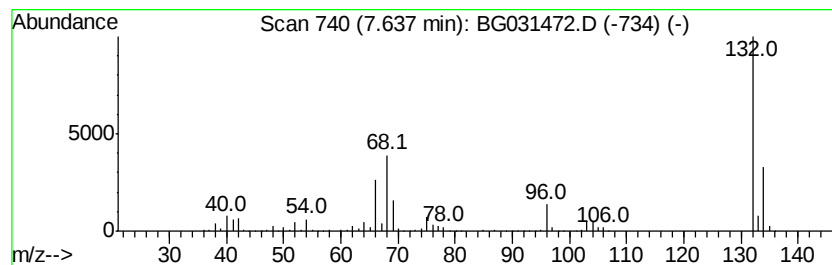
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Peak Number 2 unknown7.64 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.64	61.14 ng	891192	1,4-Dichlorobenzene-d4	8.11

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	25
3			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
4			Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
5			2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.18	5.0	ng	73020	1	8.11	291525	20.0
unknown7.64	7.64	61.1	ng	891192	1	8.11	291525	20.0