

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072718\
 Data File : BG035934.D
 Acq On : 27 Jul 2018 12:50
 Operator : JU/SJ
 Sample : PB111458BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

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 G\METHODS\8270-BG071218.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

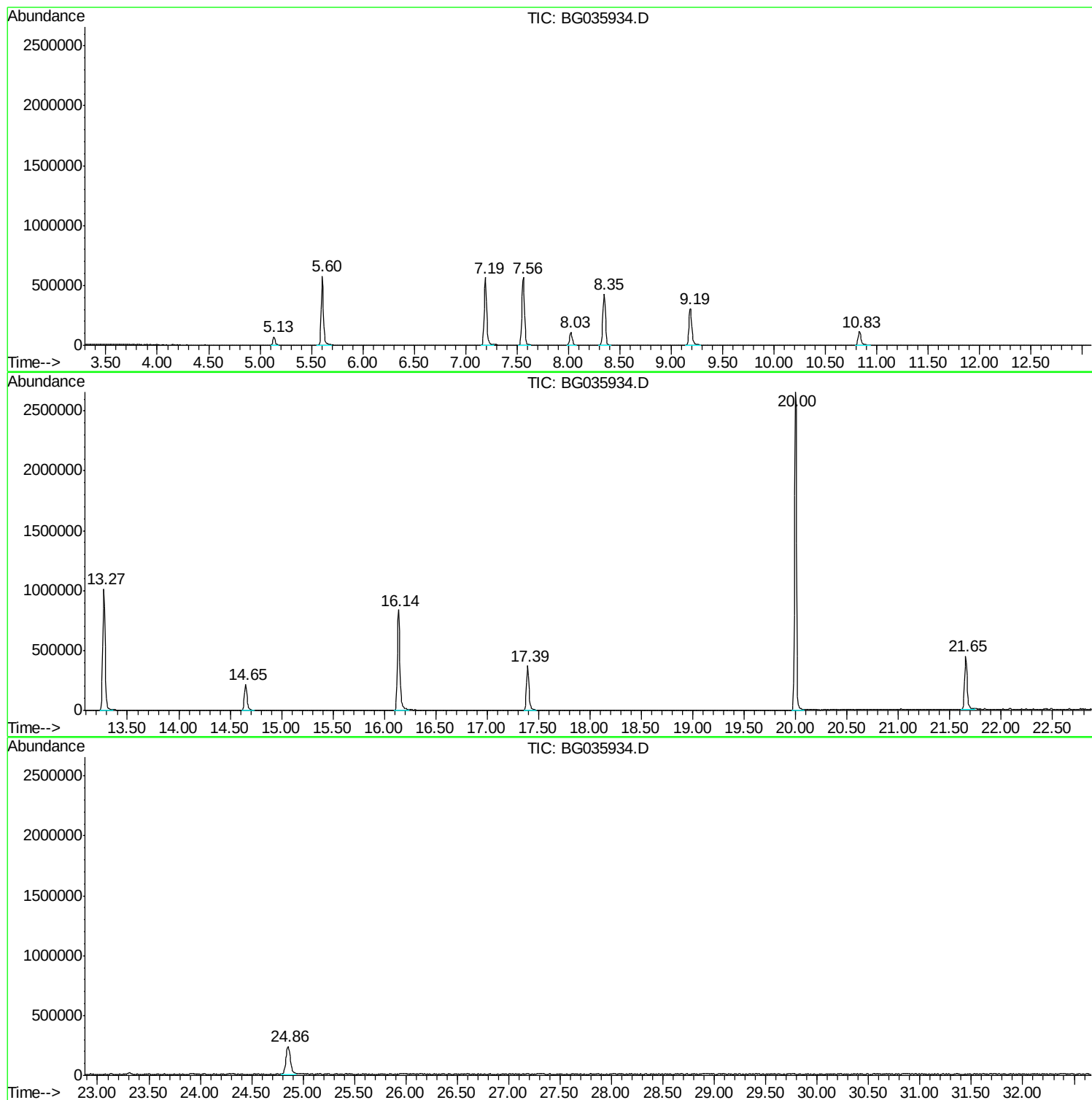
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.135	308	314	321	rBV	71291	110137	3.08%	0.786%
2	5.605	385	394	409	rBV	575381	937852	26.23%	6.691%
3	7.191	657	664	678	rBV	565909	1019164	28.50%	7.271%
4	7.561	719	727	738	rBV	564072	978639	27.37%	6.982%
5	8.025	798	806	814	rBV	109289	185489	5.19%	1.323%
6	8.348	852	861	871	rBV	428479	765471	21.41%	5.461%
7	9.188	995	1004	1020	rBV	304250	613206	17.15%	4.375%
8	10.827	1277	1283	1301	rBV	114603	238164	6.66%	1.699%
9	13.271	1692	1699	1714	rBV	1015383	1630193	45.59%	11.631%
10	14.651	1927	1934	1947	rBV3	220548	397467	11.12%	2.836%
11	16.138	2180	2187	2204	rBV	841348	1420549	39.73%	10.135%
12	17.395	2395	2401	2414	rBV	371936	627464	17.55%	4.477%
13	19.997	2838	2844	2858	rBV	2655704	3575913	100.00%	25.512%
14	21.654	3118	3126	3141	rBV2	445674	800440	22.38%	5.711%
15	24.855	3659	3671	3684	rBV2	233508	716208	20.03%	5.110%

Sum of corrected areas: 14016356

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



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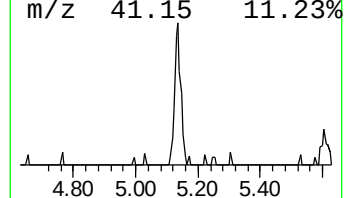
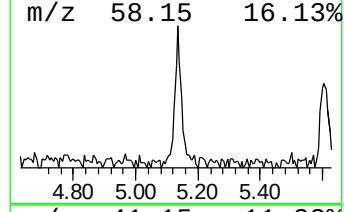
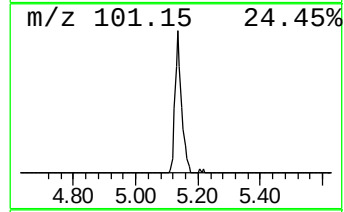
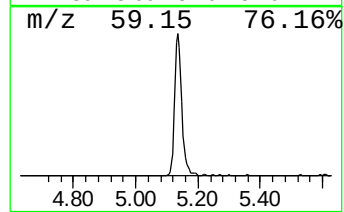
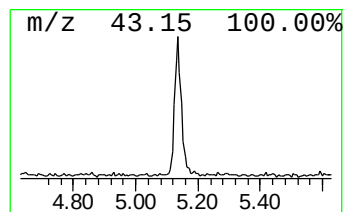
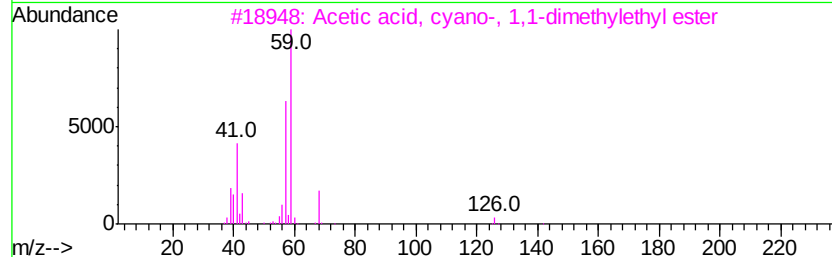
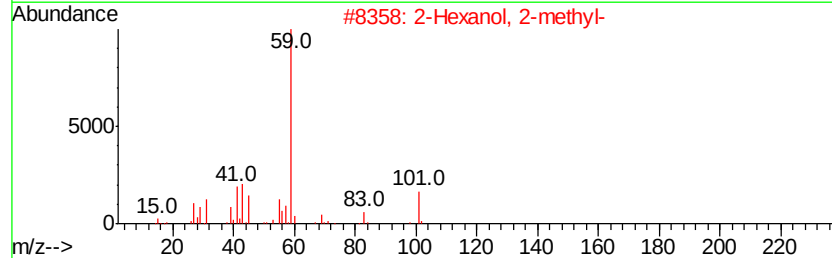
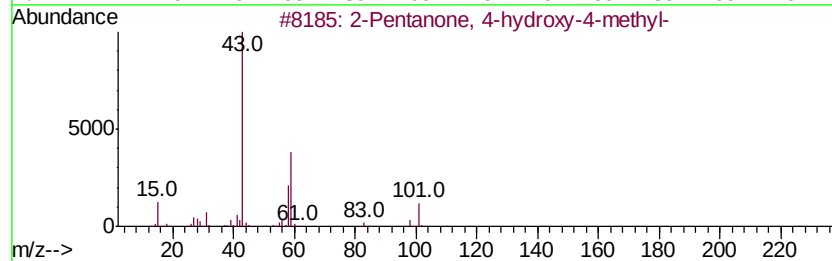
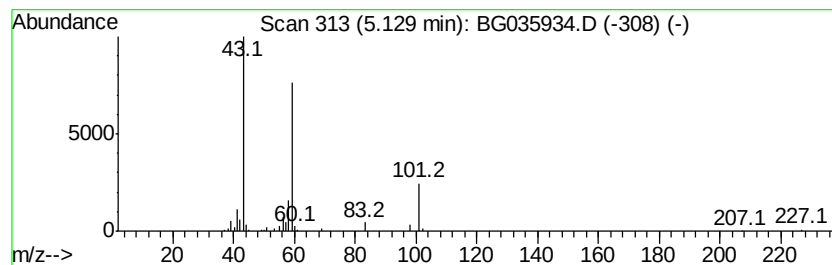
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.13	11.88 ng	110137	1,4-Dichlorobenzene-d4	8.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	25
3	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
4	5-Hexen-2-one	98	C6H10O	000109-49-9	9
5	4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9



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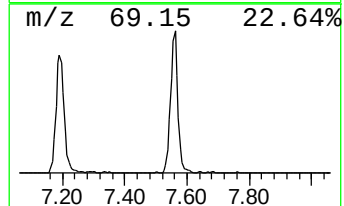
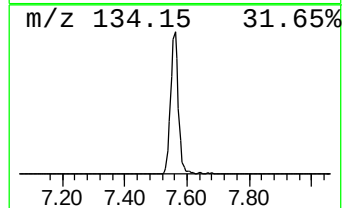
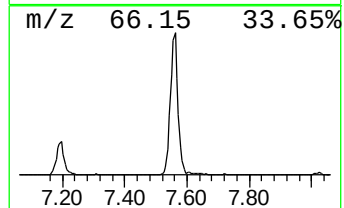
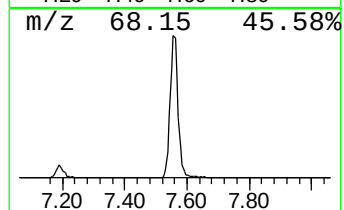
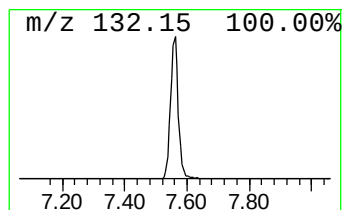
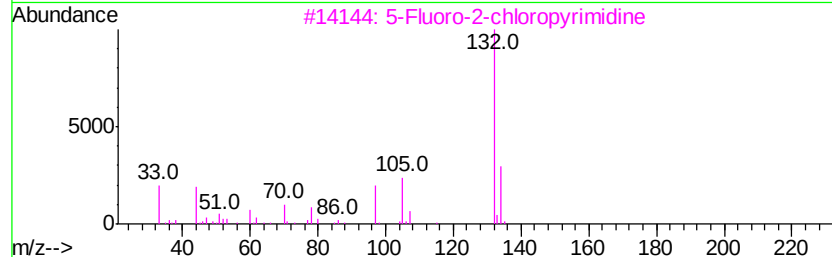
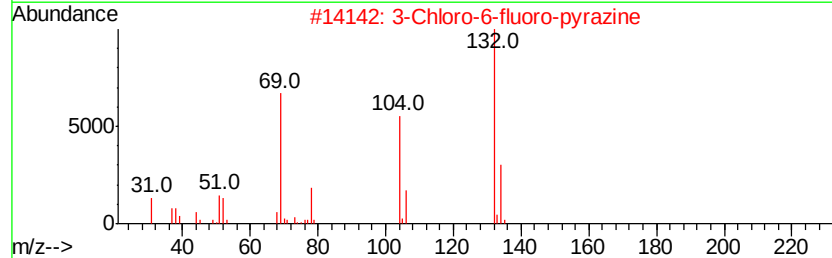
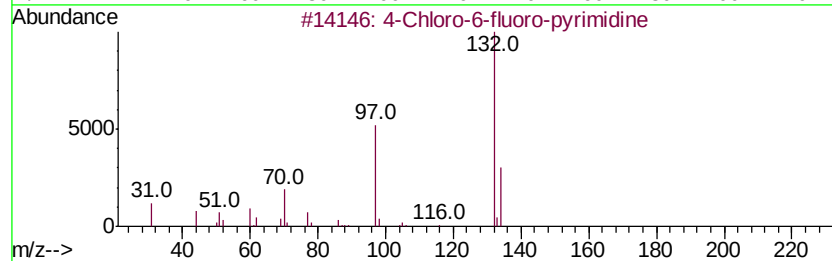
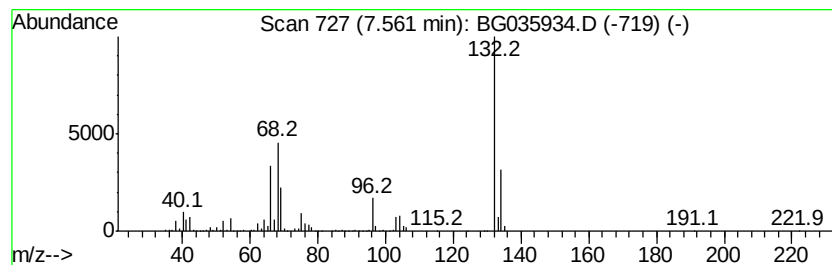
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Peak Number 2 unknown7.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	105.52 ng	978639	1,4-Dichlorobenzene-d4	8.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.13	11.9	ng	110137	1	8.03	185489	20.0
unknown7.56	7.56	105.5	ng	978639	1	8.03	185489	20.0