

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072718\
 Data File : BG035932.D
 Acq On : 27 Jul 2018 11:33
 Operator : JU/SJ
 Sample : PB111401BL
 Misc :
 ALS Vial : 4 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

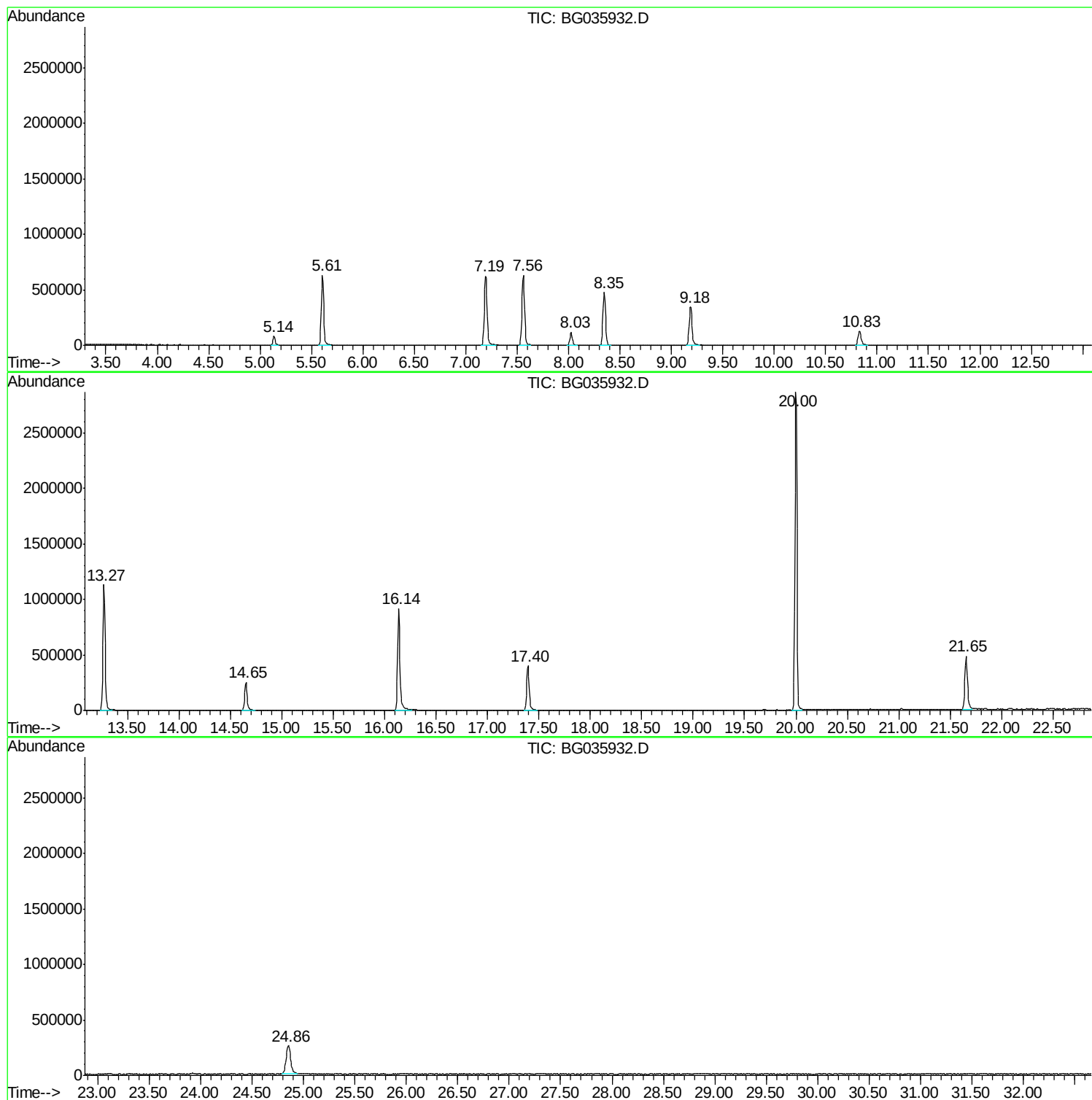
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.135	309	314	321	rBV2	80544	126932	3.22%	0.816%
2	5.605	387	394	408	rBV	631779	1039109	26.36%	6.678%
3	7.191	656	664	681	rBV	624084	1147183	29.10%	7.373%
4	7.561	720	727	739	rBV	629351	1102122	27.96%	7.083%
5	8.025	799	806	815	rBV	117603	205833	5.22%	1.323%
6	8.349	853	861	871	rBV	478265	855031	21.69%	5.495%
7	9.183	996	1003	1018	rBV	345592	682080	17.30%	4.384%
8	10.828	1276	1283	1297	rBV	126791	269230	6.83%	1.730%
9	13.271	1692	1699	1712	rBV	1132329	1812411	45.97%	11.648%
10	14.652	1927	1934	1949	rBV2	252028	441579	11.20%	2.838%
11	16.138	2180	2187	2209	rBV	916677	1615886	40.99%	10.385%
12	17.395	2394	2401	2416	rBV	400803	681612	17.29%	4.381%
13	19.997	2838	2844	2855	rBV	2863469	3942210	100.00%	25.336%
14	21.654	3120	3126	3135	rBV2	480064	851225	21.59%	5.471%
15	24.856	3660	3671	3685	rBV	255122	787580	19.98%	5.062%

Sum of corrected areas: 15560023

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



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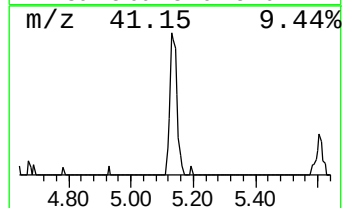
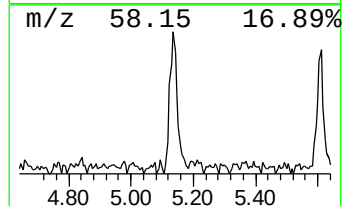
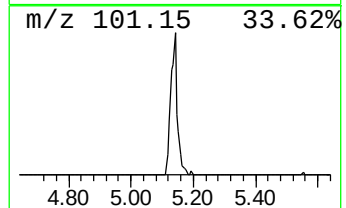
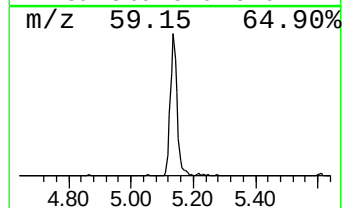
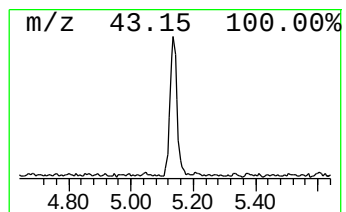
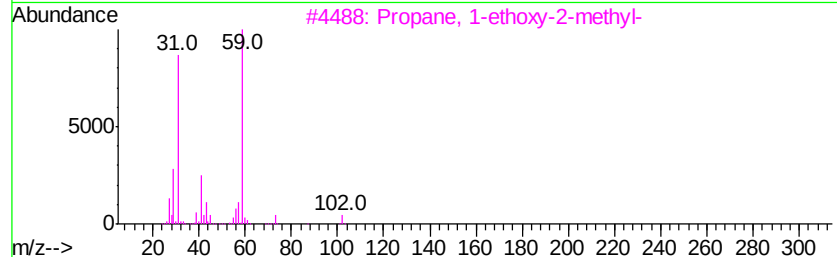
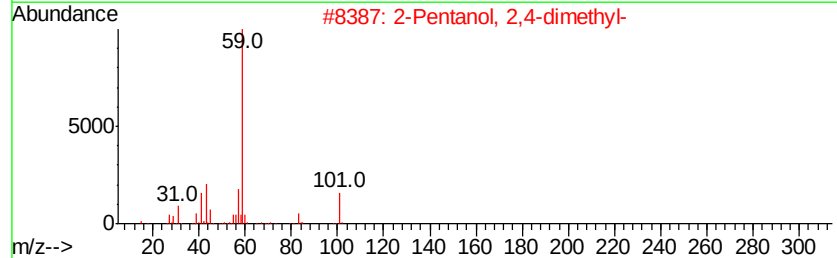
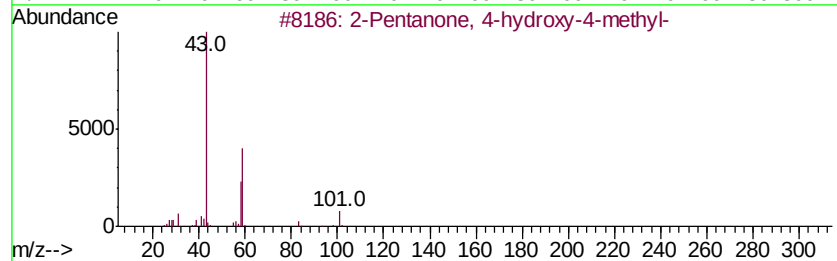
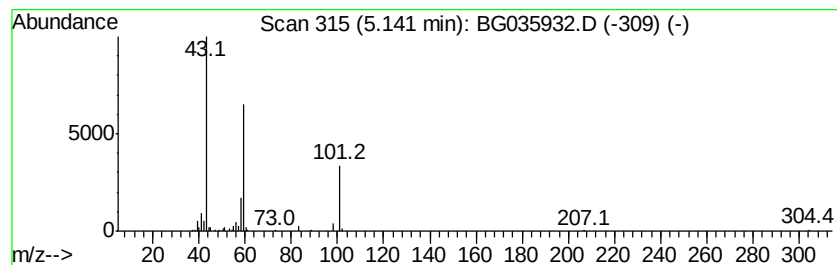
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	12.33 ng	126932	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	56
2			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	40
3			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	25
4			1,2-Butanediol	90	C4H10O2	000584-03-2	23
5			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	17



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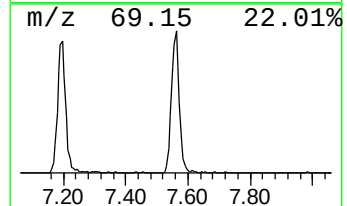
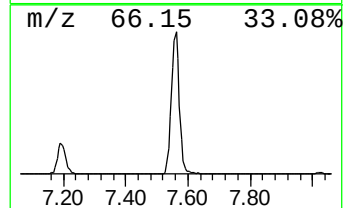
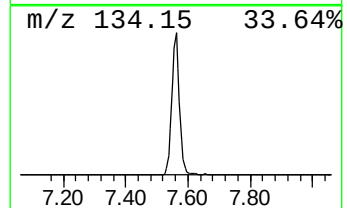
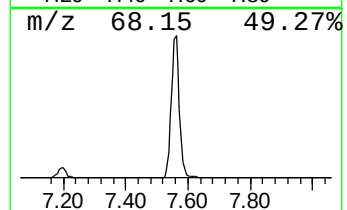
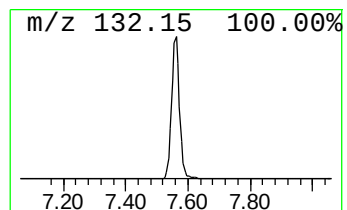
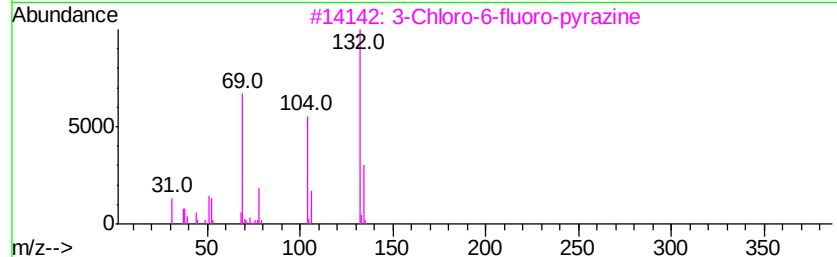
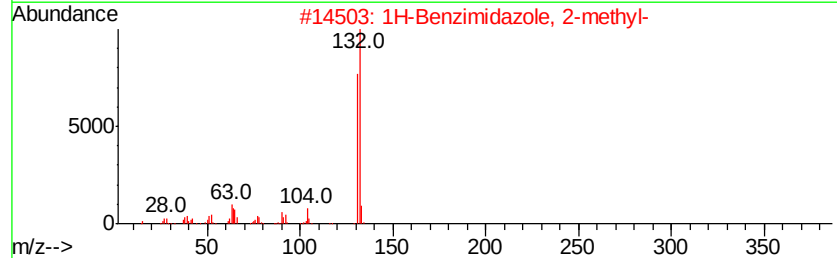
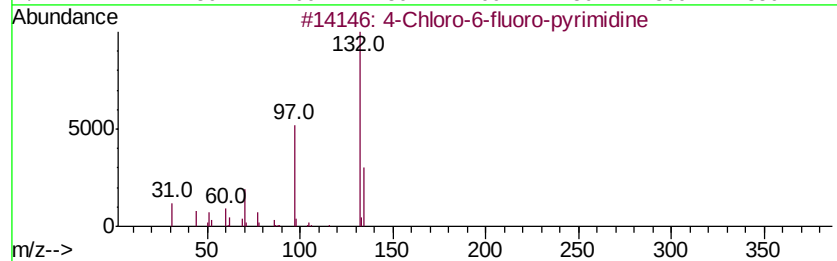
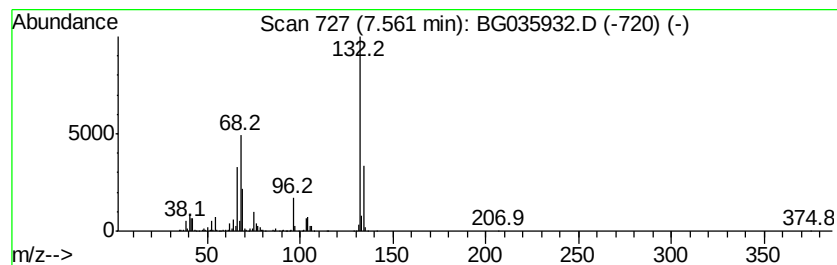
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Peak Number 2 unknown7.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	107.09 ng	1102120	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		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
4		N-(-3,4-Methylenedioxyphenyl)isopropylamine	267	C17H17NO2	1000378-96-6	10
5		Tranlylcypromine-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...	5.14	12.3	ng	126932	1	8.03	205833	20.0
unknown7.56	7.56	107.1	ng	1102120	1	8.03	205833	20.0