

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120418\
 Data File : BG038350.D
 Acq On : 5 Dec 2018 8:26
 Operator : JU/SJ
 Sample : PB115253BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB115253BL

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-BG120418.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	5.151	312	317	323	rBV	23929	36005	1.49%	0.342%
2	5.615	389	396	410	rVB	473965	789268	32.73%	7.493%
3	7.219	662	669	679	rBV	444310	785633	32.58%	7.458%
4	7.595	726	733	743	rBV	419750	739481	30.66%	7.020%
5	8.071	806	814	823	rBV	114114	204782	8.49%	1.944%
6	8.395	860	869	879	rVB	311211	570915	23.67%	5.420%
7	9.246	1004	1014	1024	rBV	261090	493171	20.45%	4.682%
8	10.892	1285	1294	1308	rBV	183148	360453	14.95%	3.422%
9	13.330	1700	1709	1721	rBV	703604	1159807	48.09%	11.011%
10	14.711	1936	1944	1955	rBV3	212888	373742	15.50%	3.548%
11	16.197	2190	2197	2212	rBV2	587874	925562	38.38%	8.787%
12	17.454	2403	2411	2420	rBV	319927	521018	21.60%	4.946%
13	20.057	2847	2854	2862	rBV	1734241	2411621	100.00%	22.895%
14	21.738	3130	3140	3151	rBV	347666	596123	24.72%	5.659%
15	25.051	3693	3704	3715	rVB2	187278	566033	23.47%	5.374%

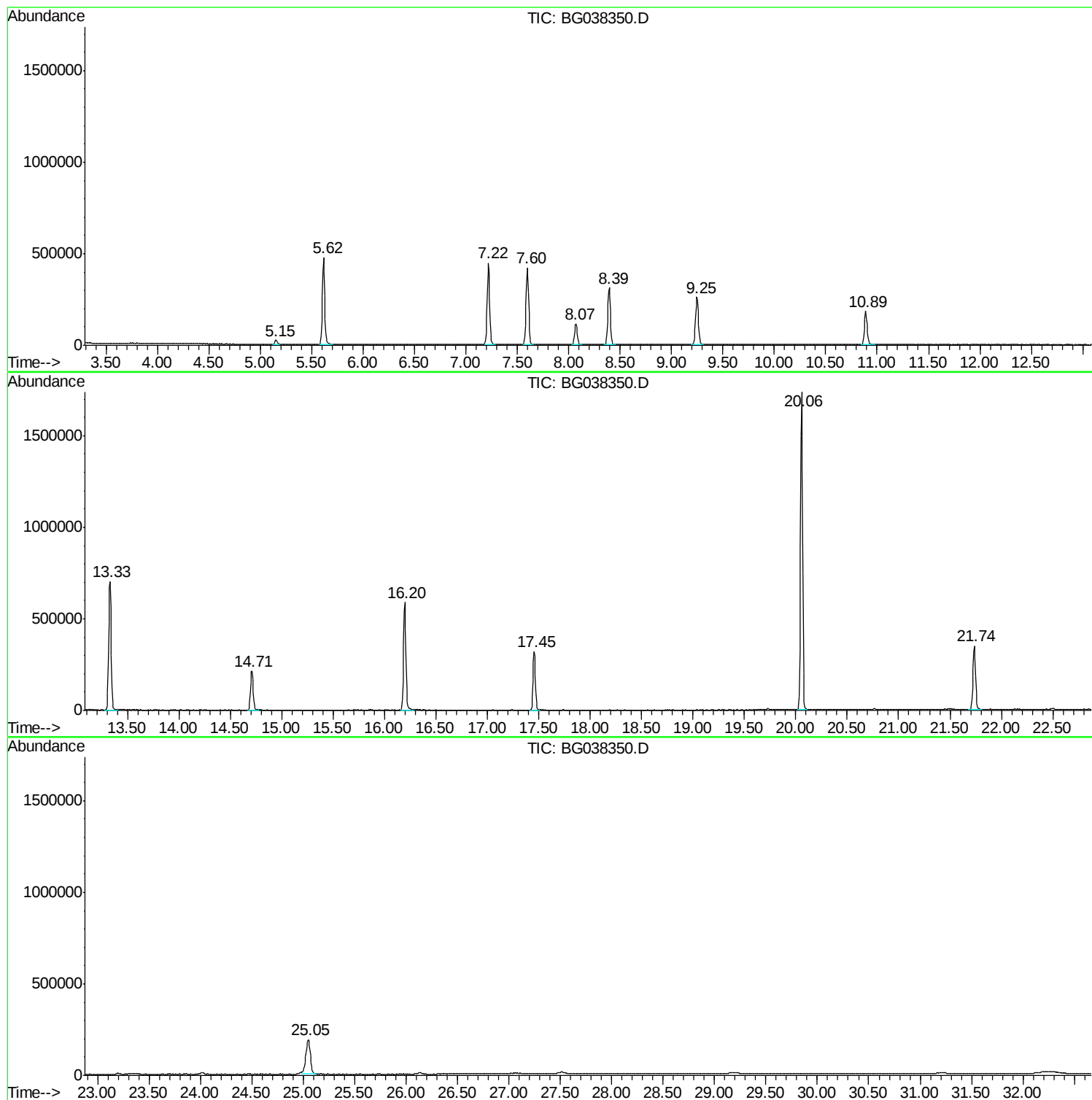
Sum of corrected areas: 10533614

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



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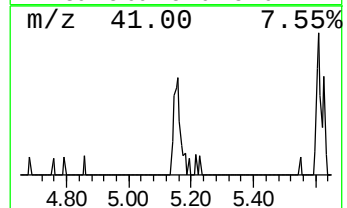
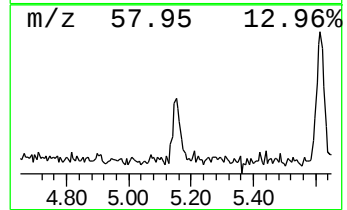
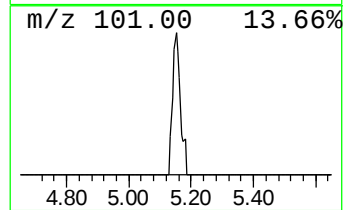
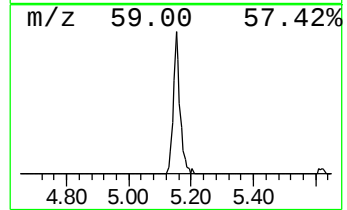
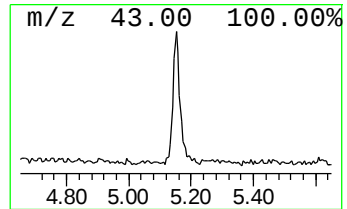
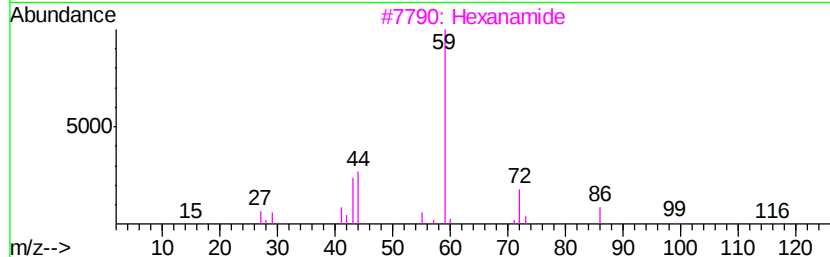
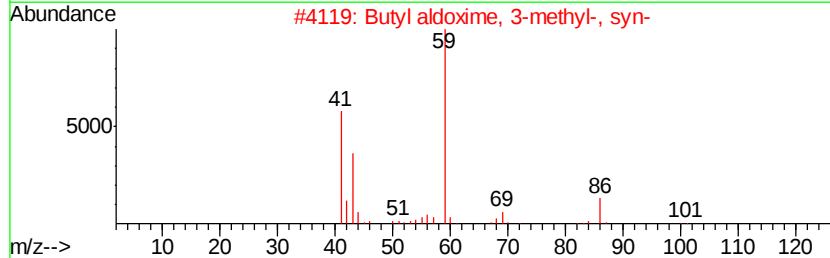
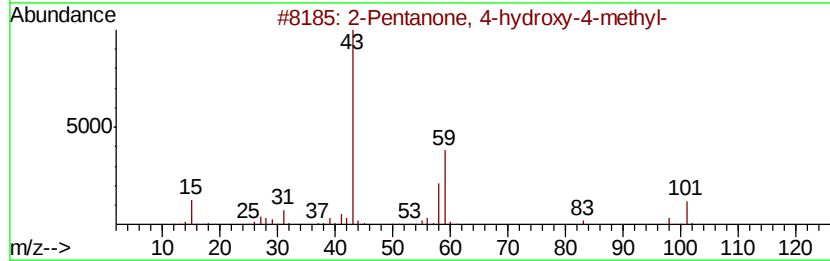
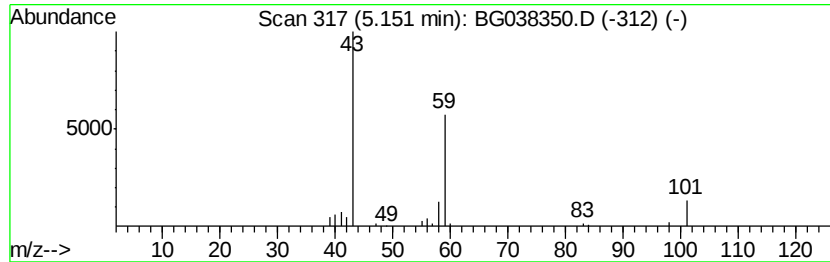
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.15	3.52 ng	36005	1,4-Dichlorobenzene-d4	8.07

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	23
3			Hexanamide	115	C6H13NO	000628-02-4	9
4			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	9
5			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	9



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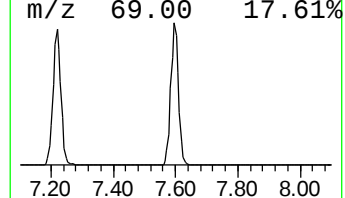
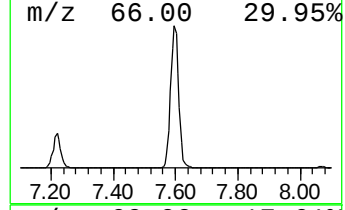
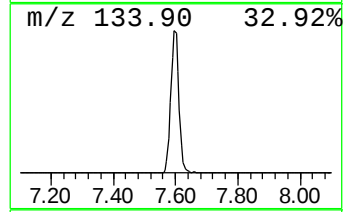
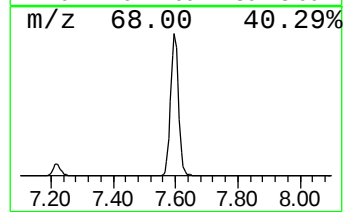
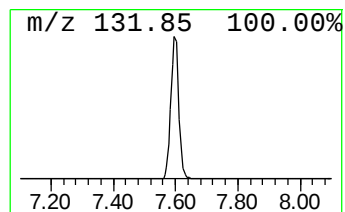
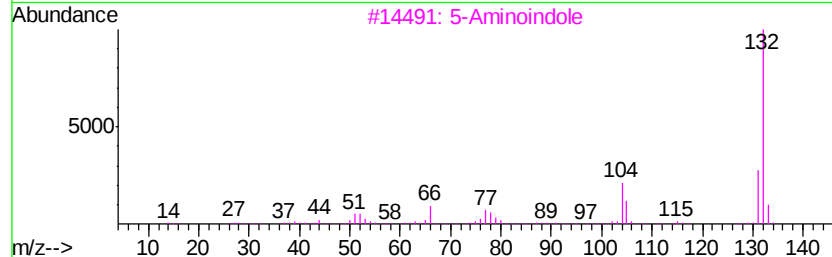
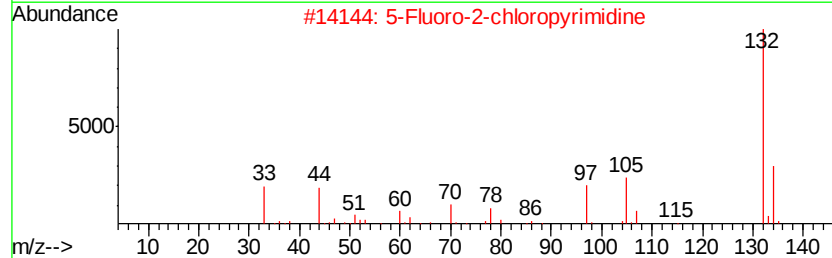
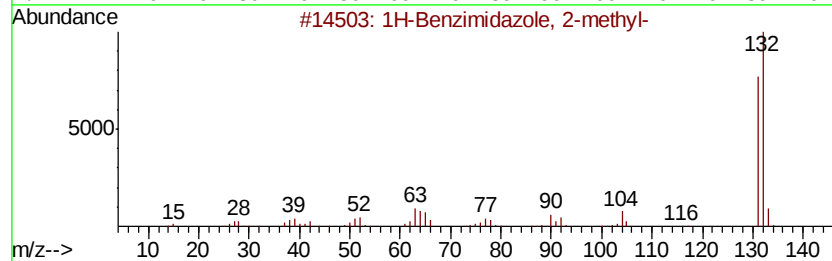
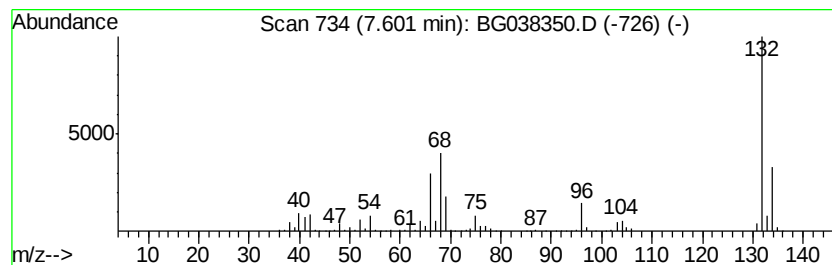
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Peak Number 2 unknown7.60 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.60	72.22 ng	739481	1,4-Dichlorobenzene-d4	8.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.15	3.5	ng	36005	1	8.07	204782	20.0
unknown7.60	7.60	72.2	ng	739481	1	8.07	204782	20.0