

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG112618\
 Data File : BG038152.D
 Acq On : 27 Nov 2018 1:59
 Operator : JU/SJ
 Sample : PB115042BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB115042BL

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-BG111318.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	5.168	315	320	331	rVB	40207	69151	2.16%	0.411%
2	5.632	392	399	410	rBV	820234	1374852	42.88%	8.171%
3	7.236	664	672	682	rBV	852143	1546110	48.22%	9.189%
4	7.618	725	737	749	rBV	799166	1448139	45.17%	8.606%
5	8.088	807	817	825	rBV	154084	268125	8.36%	1.593%
6	8.411	864	872	881	rBV	596513	1070682	33.39%	6.363%
7	9.263	1009	1017	1028	rBV	513060	966988	30.16%	5.747%
8	10.908	1287	1297	1305	rBV	209897	392854	12.25%	2.335%
9	13.346	1704	1712	1721	rBV	1387571	2257015	70.39%	13.414%
10	14.721	1938	1946	1955	rBV2	337384	564462	17.61%	3.355%
11	16.213	2192	2200	2215	rBV3	958285	1590226	49.60%	9.451%
12	17.471	2407	2414	2424	rBV2	429169	677219	21.12%	4.025%
13	20.074	2850	2857	2867	rBV	2269187	3206241	100.00%	19.055%
14	21.754	3137	3143	3153	rBV	426815	736764	22.98%	4.379%
15	25.080	3696	3709	3731	rBV	192117	657601	20.51%	3.908%

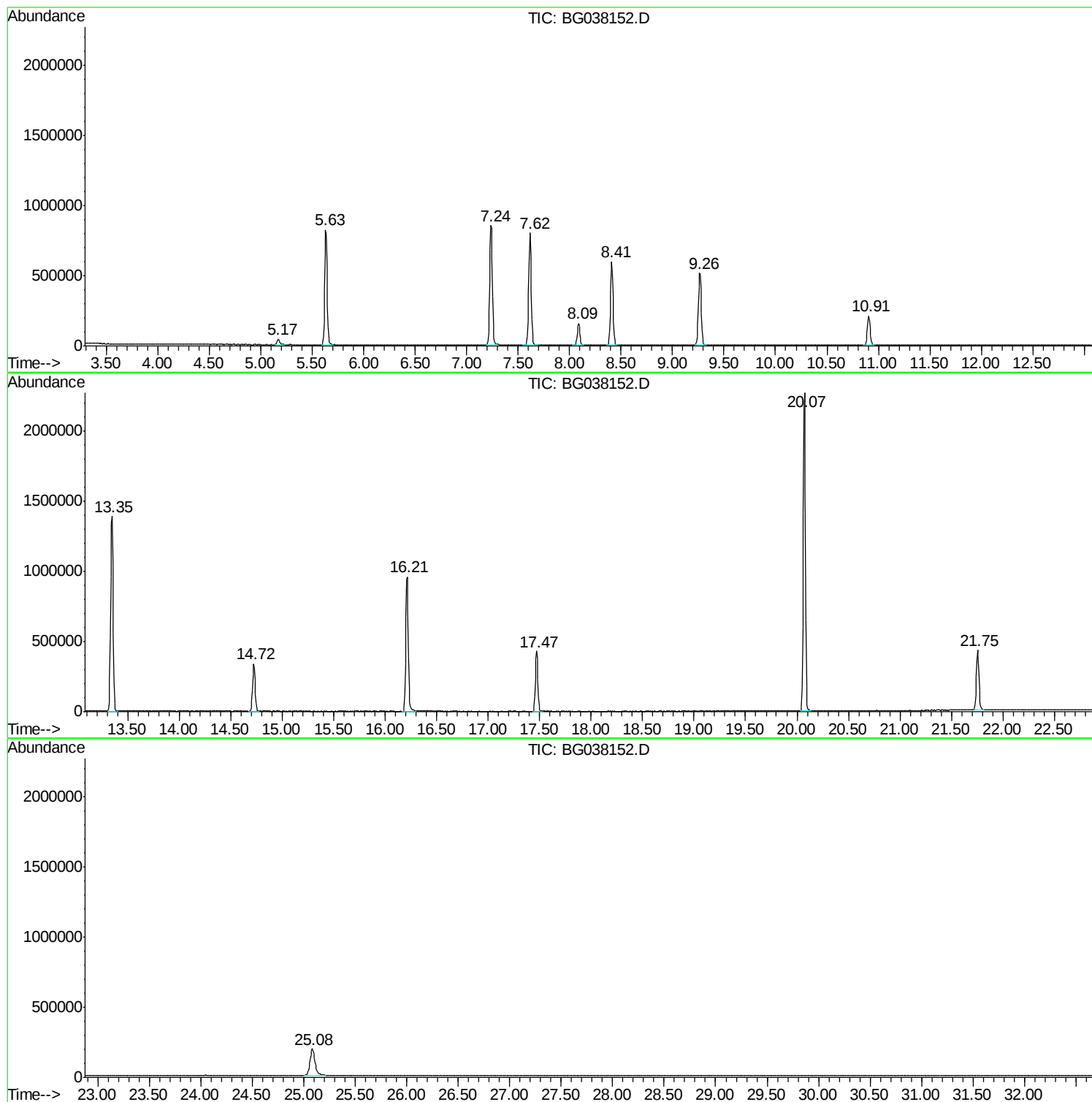
Sum of corrected areas: 16826429

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



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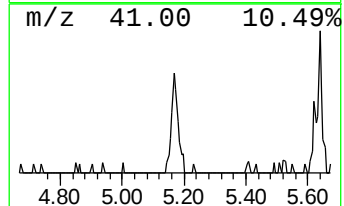
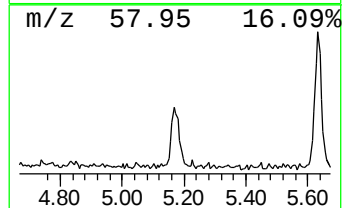
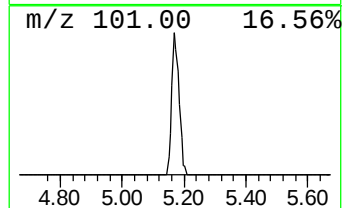
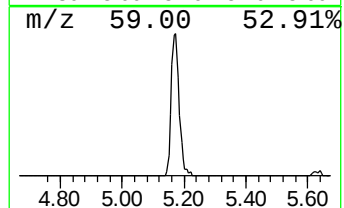
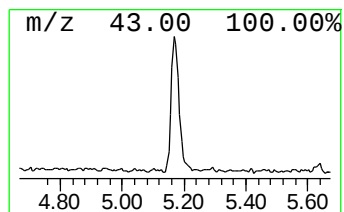
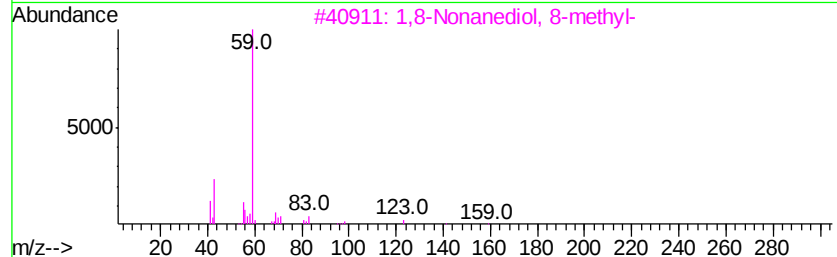
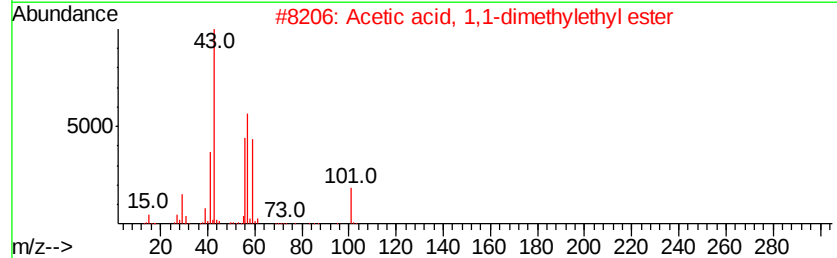
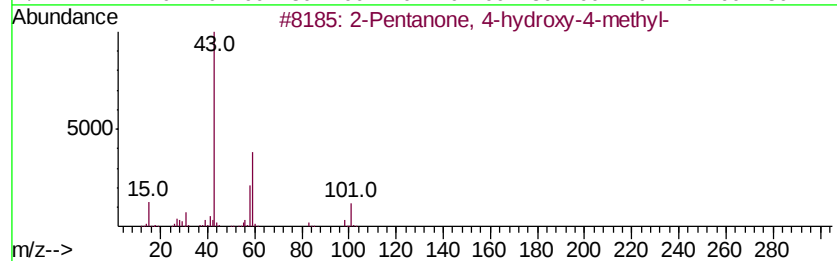
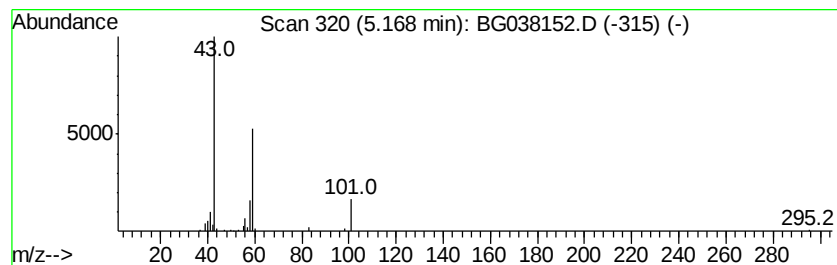
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	5.16 ng	69151	1,4-Dichlorobenzene-d4	8.09

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	37
4			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25



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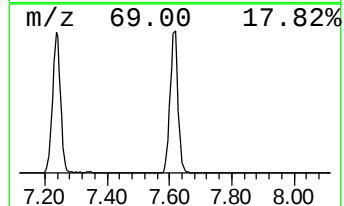
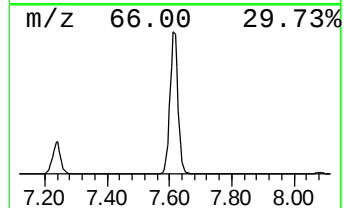
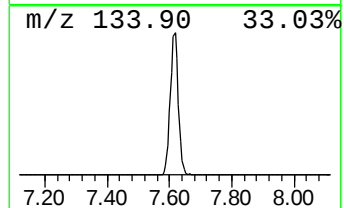
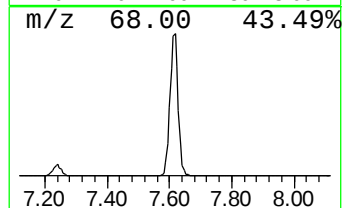
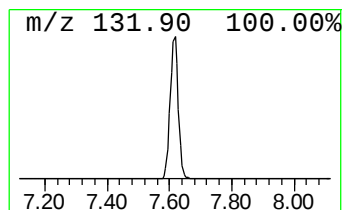
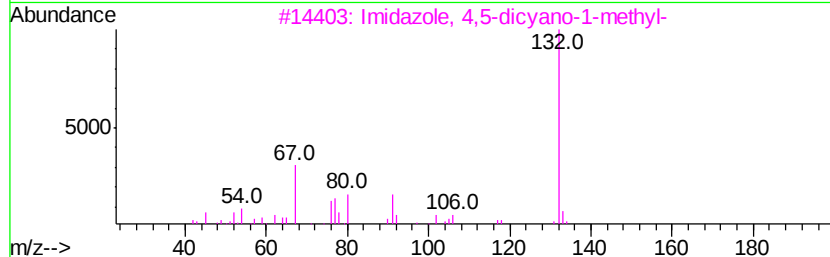
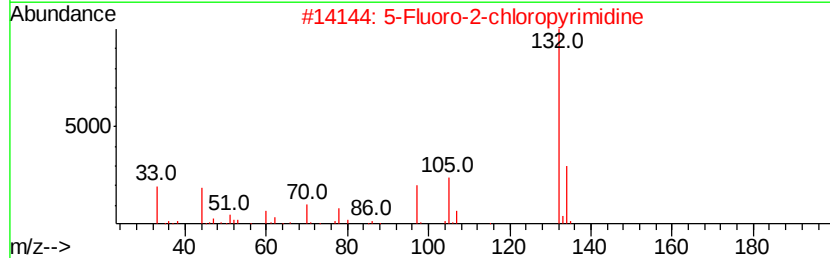
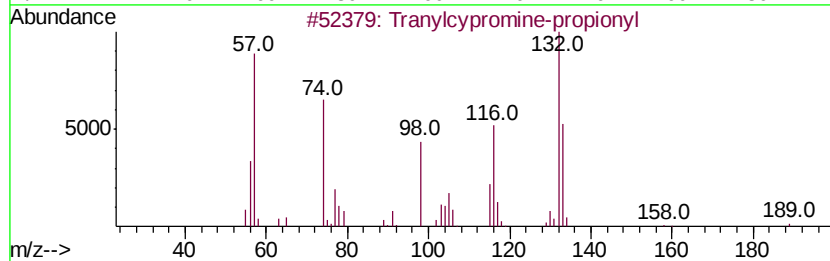
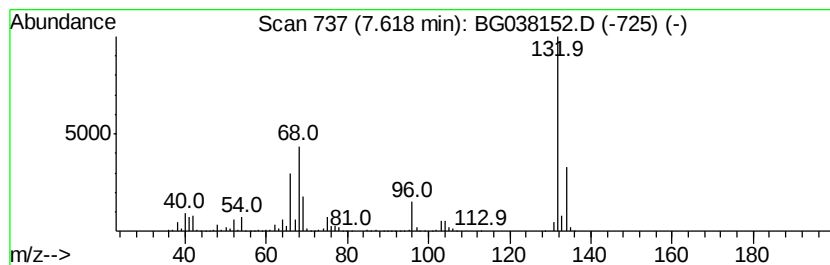
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Peak Number 2 unknown7.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	108.02 ng	1448140	1,4-Dichlorobenzene-d4	8.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	9
4		Xenon	132	Xe	007440-63-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.17	5.2	ng	69151	1	8.09	268125	20.0
unknown7.62	7.62	108.0	ng	1448140	1	8.09	268125	20.0