

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111618\
 Data File : BG038042.D
 Acq On : 16 Nov 2018 20:37
 Operator : JU/SJ
 Sample : PB114874BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB114874BL

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	5.169	315	320	328	rVB	31984	51101	1.82%	0.355%
2	5.627	391	398	410	rBV	686560	1159287	41.30%	8.064%
3	7.231	663	671	684	rBV	716049	1264635	45.06%	8.797%
4	7.613	728	736	747	rVB	674070	1199629	42.74%	8.345%
5	8.083	809	816	826	rVB	129747	230039	8.20%	1.600%
6	8.406	861	871	880	rBV	506748	917936	32.70%	6.385%
7	9.264	1009	1017	1026	rBV	422074	782666	27.88%	5.444%
8	10.903	1289	1296	1305	rVB	176829	332649	11.85%	2.314%
9	13.341	1703	1711	1719	rBV	1202503	1860391	66.28%	12.941%
10	14.722	1938	1946	1956	rVB2	291492	479332	17.08%	3.334%
11	16.209	2192	2199	2211	rBV4	844128	1317073	46.92%	9.162%
12	17.466	2406	2413	2421	rBV2	379115	610774	21.76%	4.249%
13	20.069	2850	2856	2867	rBV	2129239	2806781	100.00%	19.524%
14	21.749	3135	3142	3152	rBV	399183	682426	24.31%	4.747%
15	25.075	3691	3708	3723	rVB2	215275	681100	24.27%	4.738%

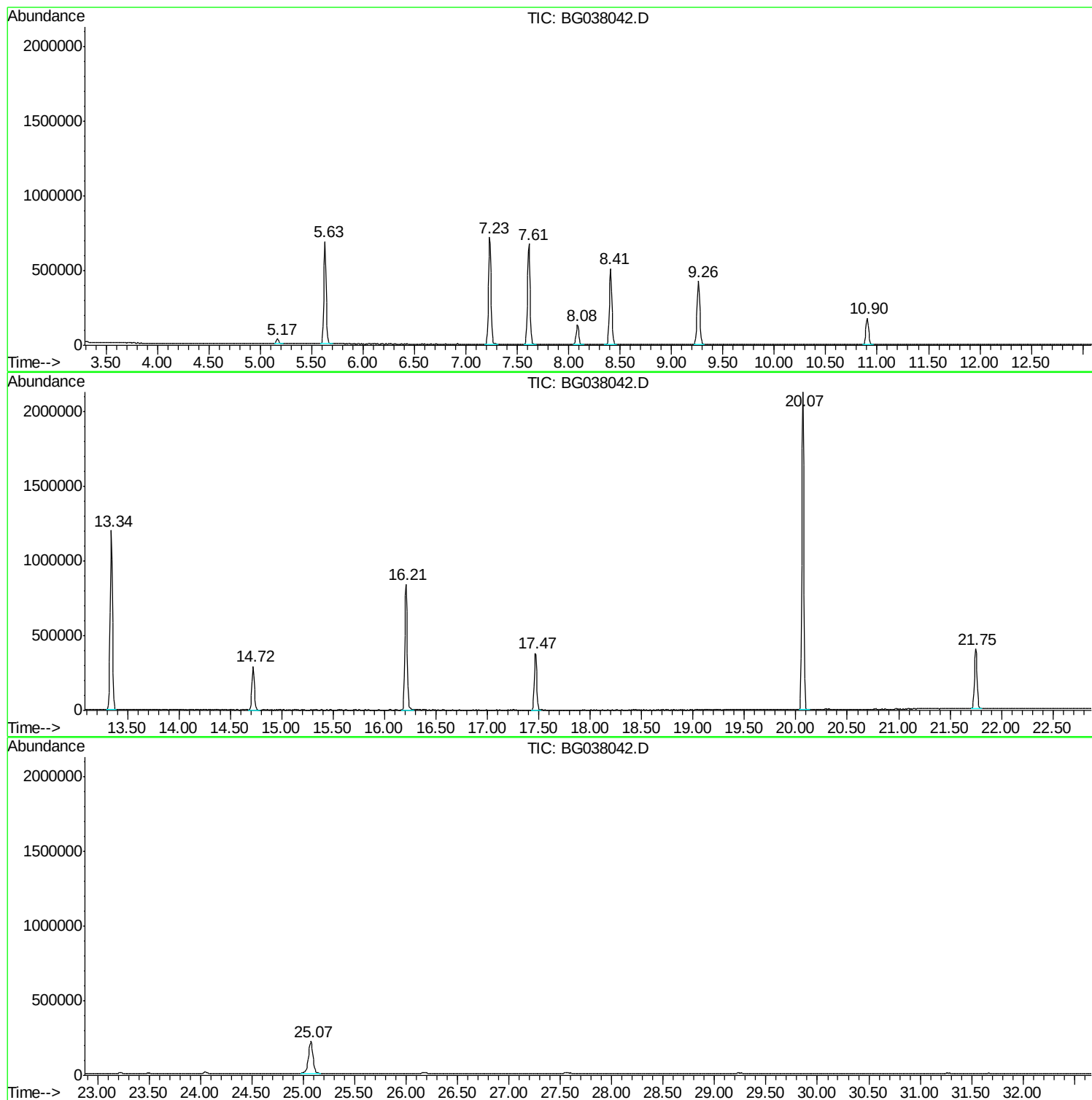
Sum of corrected areas: 14375819

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



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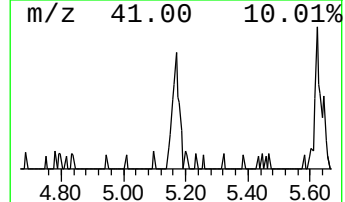
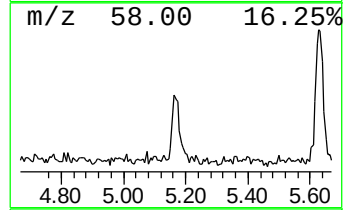
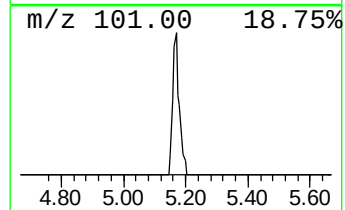
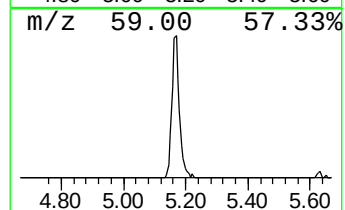
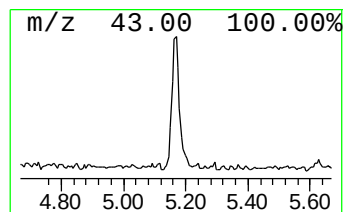
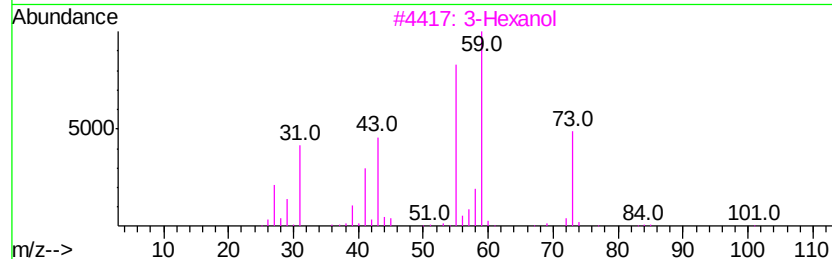
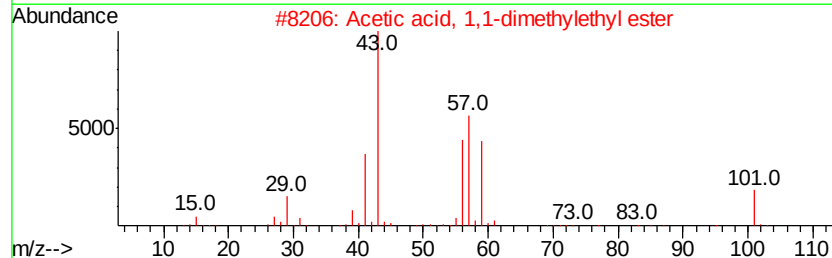
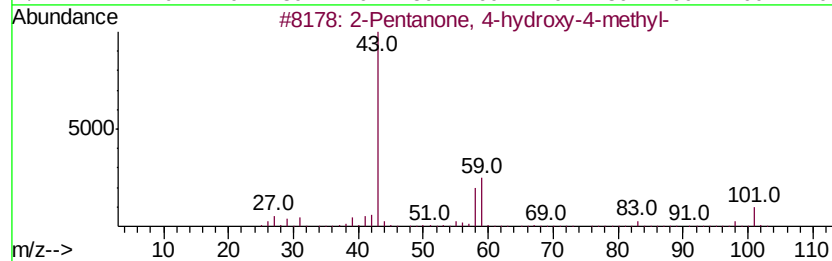
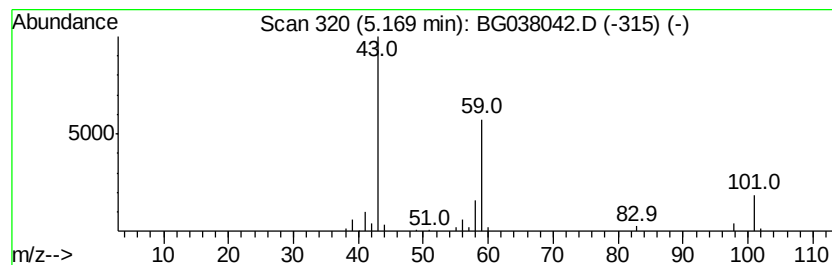
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	4.44 ng	51101	1,4-Dichlorobenzene-d4	8.08

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3			3-Hexanol	102	C6H14O	000623-37-0	33
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



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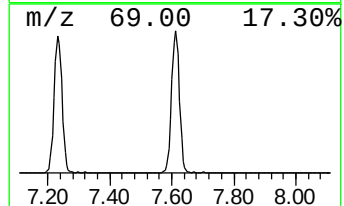
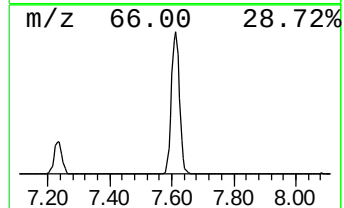
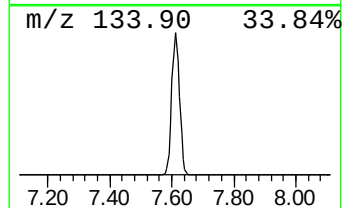
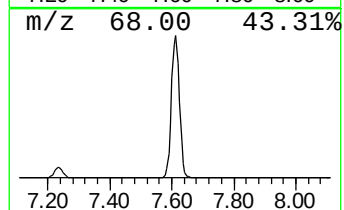
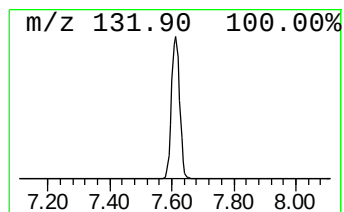
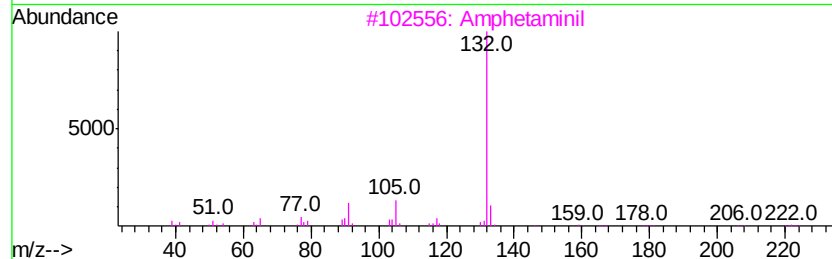
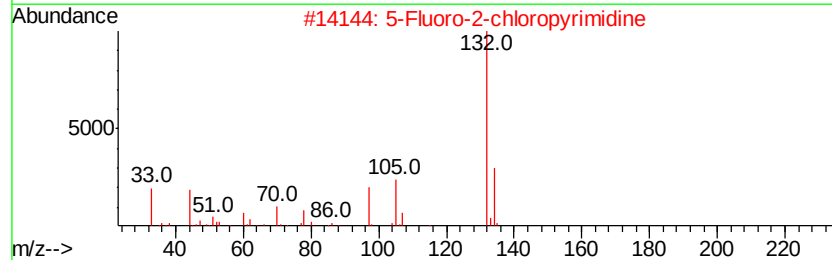
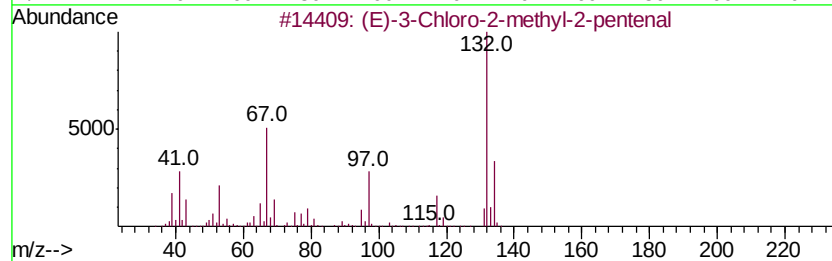
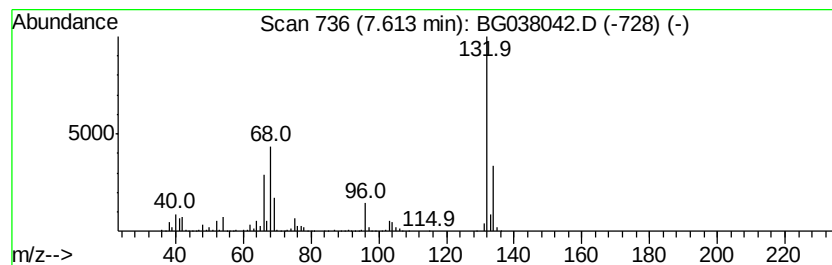
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Peak Number 2 unknown7.61 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.61	104.30 ng	1199630	1,4-Dichlorobenzene-d4	8.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		Amphetaminil	250	C17H18N2	017590-01-1	12
4		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	12
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	4.4	ng	51101	1	8.08	230039	20.0
unknown7.61	7.61	104.3	ng	1199630	1	8.08	230039	20.0