

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110918\
 Data File : BG037926.D
 Acq On : 9 Nov 2018 18:39
 Operator : JU/SJ
 Sample : PB114695BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB114695BL

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-BG101818.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	5.173	316	321	330	rBV	28505	48926	1.71%	0.324%
2	5.637	393	400	412	rBV	673709	1112615	38.88%	7.358%
3	7.235	664	672	689	rBV	708558	1278373	44.67%	8.454%
4	7.617	729	737	748	rBV	659877	1192209	41.66%	7.884%
5	8.093	811	818	826	rVB	153975	270781	9.46%	1.791%
6	8.416	865	873	882	rBV	496455	895539	31.29%	5.922%
7	9.268	1009	1018	1027	rBV	417782	788602	27.56%	5.215%
8	10.913	1290	1298	1310	rBV	234676	430261	15.03%	2.845%
9	13.346	1705	1712	1725	rBV	1266294	2065338	72.17%	13.658%
10	14.726	1940	1947	1960	rBV2	400827	673398	23.53%	4.453%
11	16.213	2192	2200	2218	rBV	824671	1342196	46.90%	8.876%
12	17.476	2406	2415	2425	rBV2	485844	797743	27.87%	5.275%
13	20.073	2851	2857	2874	rBV	2068047	2861900	100.00%	18.926%
14	21.753	3137	3143	3162	rBV	385342	814290	28.45%	5.385%
15	25.091	3699	3711	3726	rBV2	145319	549725	19.21%	3.635%

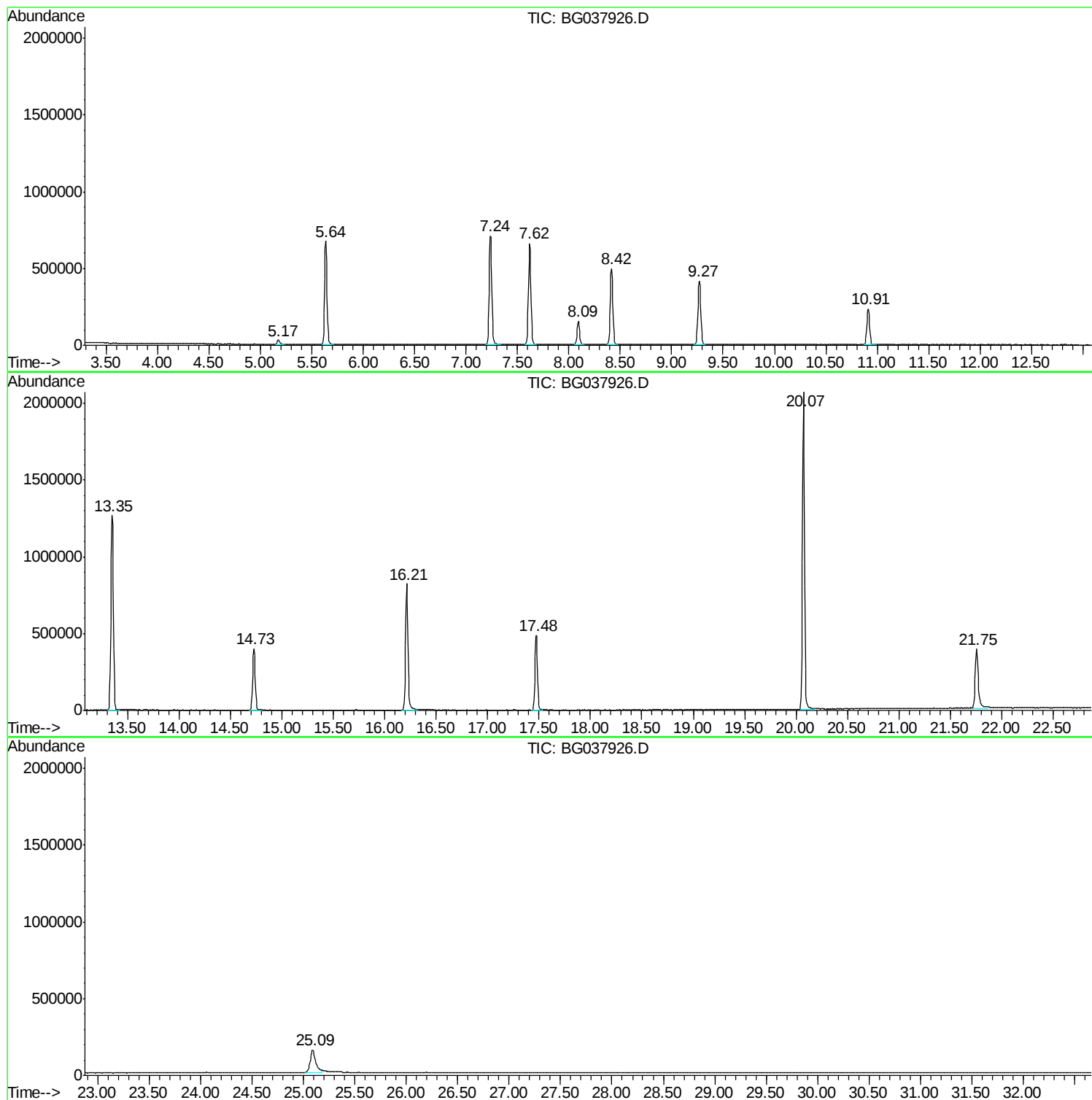
Sum of corrected areas: 15121896

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



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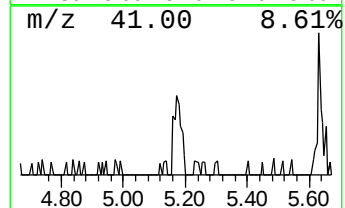
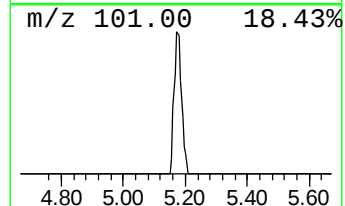
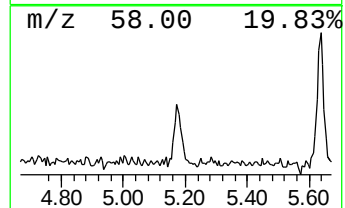
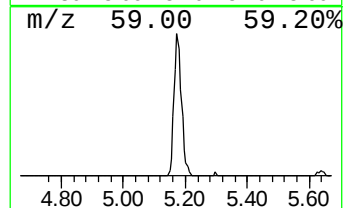
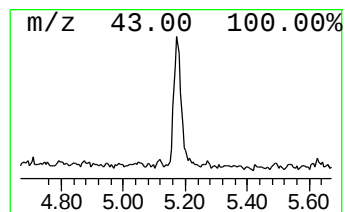
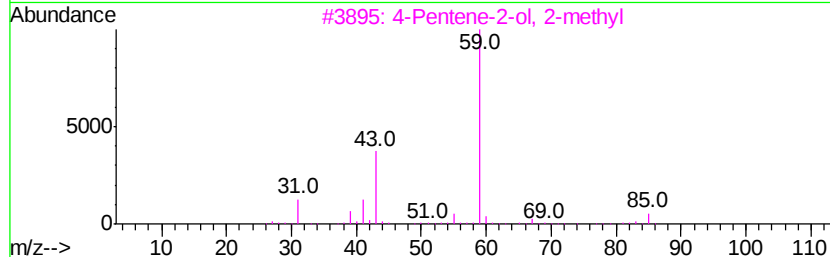
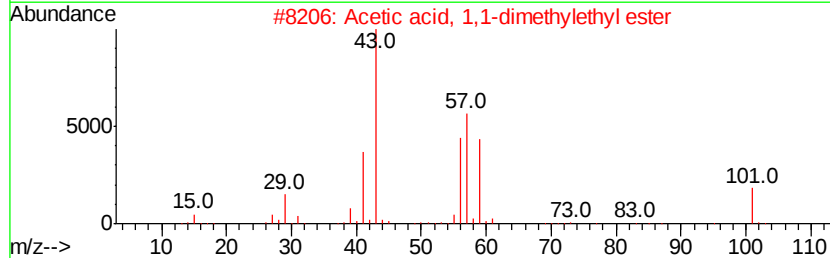
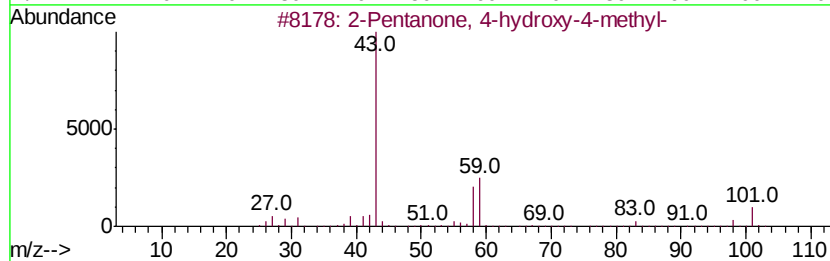
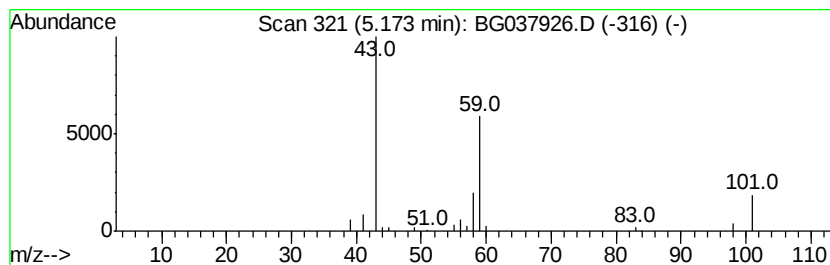
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	3.61 ng	48926	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	42
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	33
3			4-Pentene-2-ol, 2-methyl	100	C6H12O	000624-97-5	9
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	9
5			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	9



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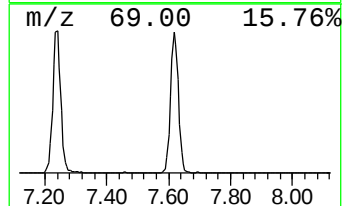
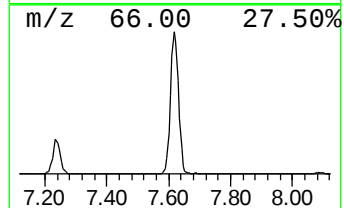
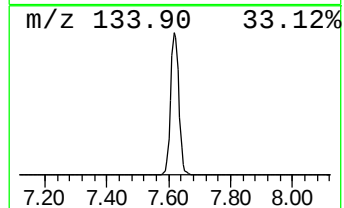
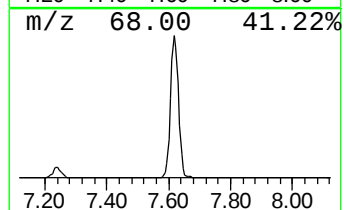
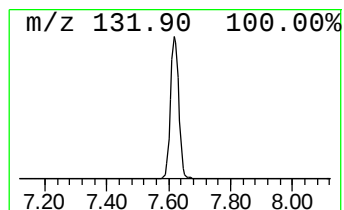
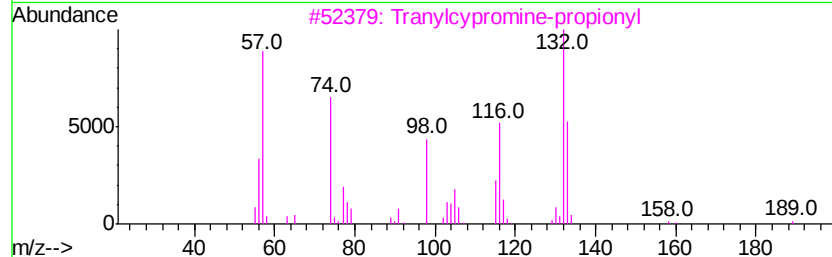
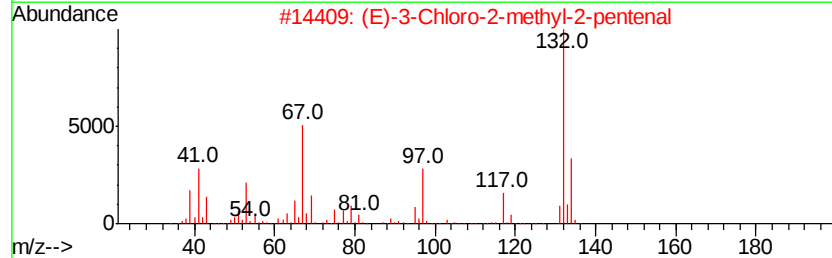
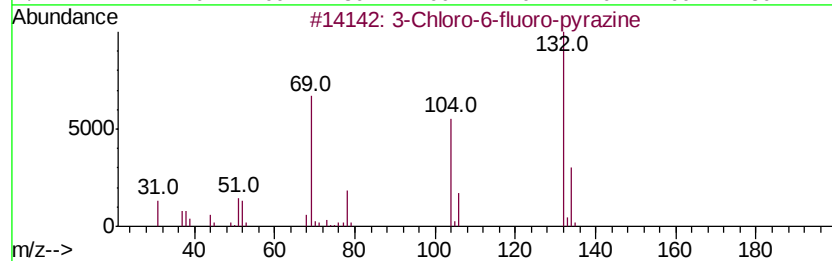
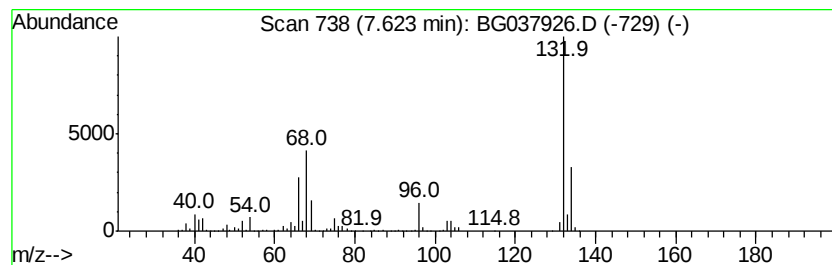
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Peak Number 2 unknown7.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	88.06 ng	1192210	1,4-Dichlorobenzene-d4	8.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	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	3.6	ng	48926	1	8.09	270781	20.0
unknown7.62	7.62	88.1	ng	1192210	1	8.09	270781	20.0