

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110818\
 Data File : BG037886.D
 Acq On : 8 Nov 2018 15:22
 Operator : JU/SJ
 Sample : J5858-07
 Misc :
 ALS Vial : 5 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-BG101818.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

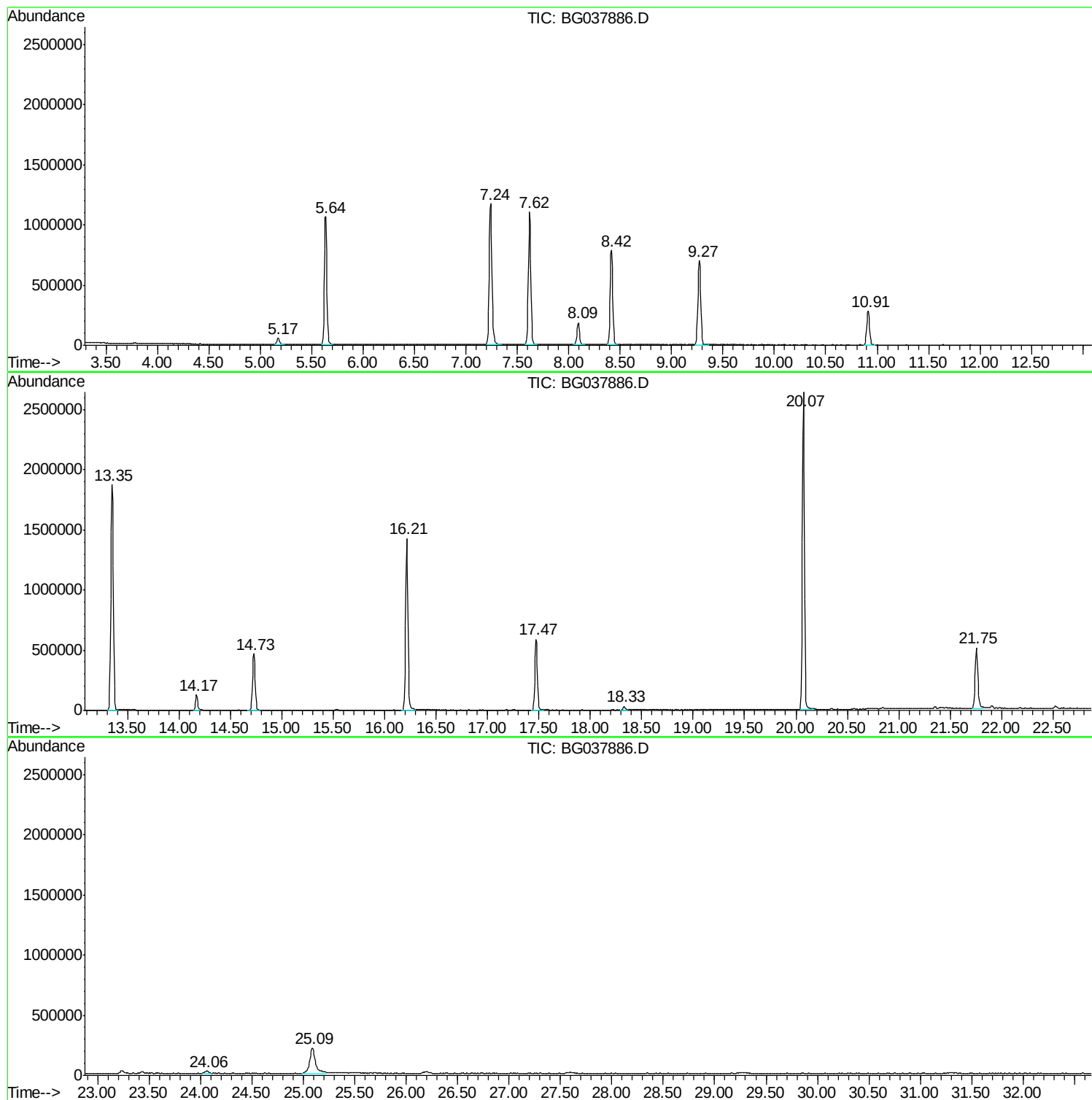
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.174	314	321	330	rVB	52563	86211	2.36%	0.384%
2	5.638	392	400	410	rBV	1061572	1809602	49.48%	8.070%
3	7.242	665	673	690	rBV	1171913	2212378	60.49%	9.867%
4	7.618	729	737	749	rBV	1100465	1953188	53.40%	8.711%
5	8.094	809	818	828	rBV	184079	330955	9.05%	1.476%
6	8.417	865	873	882	rVB	786452	1429978	39.10%	6.377%
7	9.269	1010	1018	1033	rBV	702177	1327141	36.29%	5.919%
8	10.914	1290	1298	1308	rBV	280315	518858	14.19%	2.314%
9	13.347	1704	1712	1722	rBV	1873960	3062791	83.74%	13.659%
10	14.169	1846	1852	1858	rBV	126343	186629	5.10%	0.832%
11	14.727	1937	1947	1956	rBV2	470618	799828	21.87%	3.567%
12	16.214	2192	2200	2214	rBV	1426604	2239589	61.24%	9.988%
13	17.471	2406	2414	2427	rBV2	589773	930458	25.44%	4.150%
14	18.329	2553	2560	2566	rBV3	29983	50279	1.37%	0.224%
15	20.074	2851	2857	2872	rBV	2637878	3657343	100.00%	16.311%
16	21.754	3136	3143	3154	rBV	504548	937193	25.62%	4.180%
17	24.058	3528	3535	3543	rVB9	20762	48647	1.33%	0.217%
18	25.086	3693	3710	3733	rBV2	207331	841494	23.01%	3.753%

Sum of corrected areas: 22422562

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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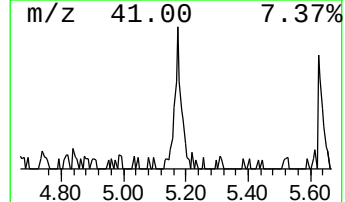
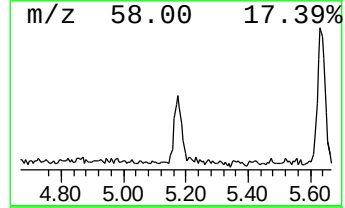
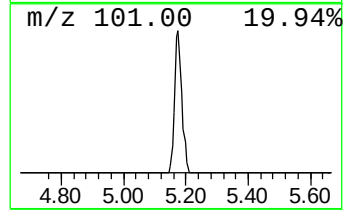
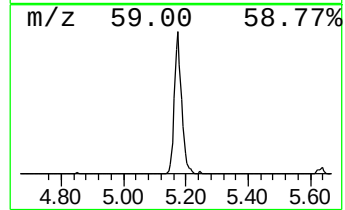
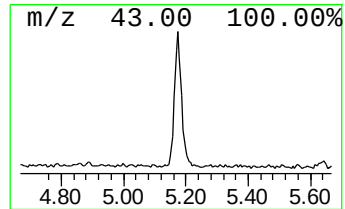
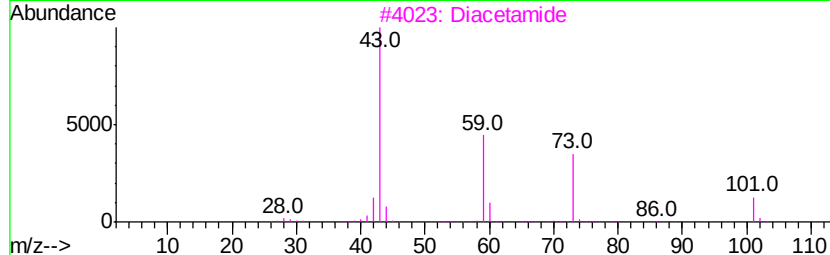
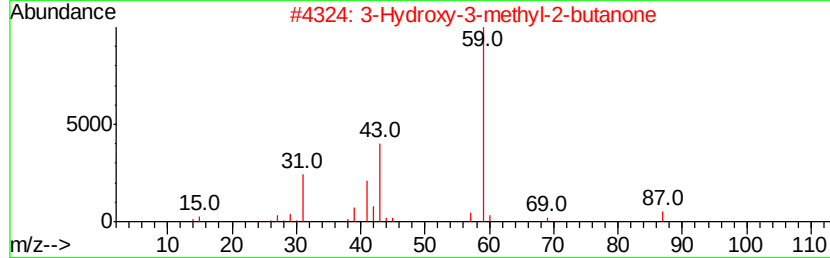
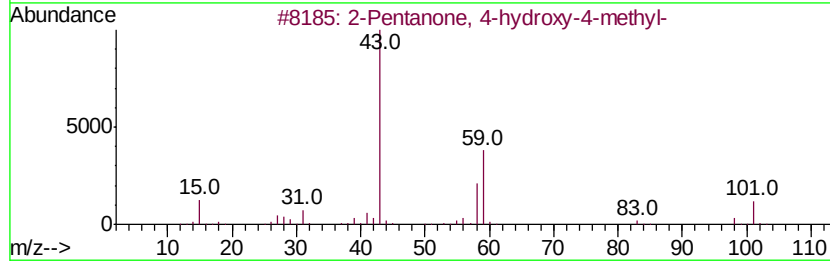
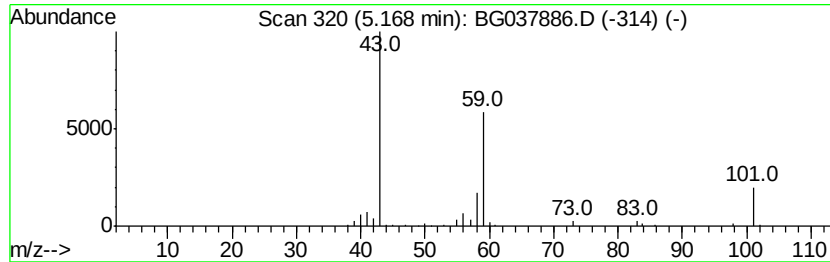
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	5.21 ng	86211	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	40
2			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	17
3			Diacetamide	101	C4H7NO2	000625-77-4	9
4			Pentanal, oxime	101	C5H11NO	000628-79-5	9
5			2-Butanone, 3-ethoxy-3-methyl-	130	C7H14O2	036687-99-7	9



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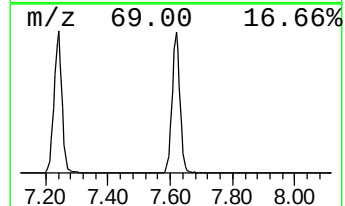
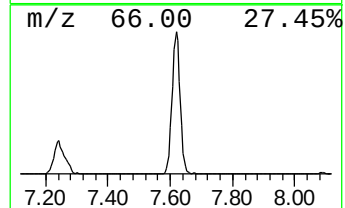
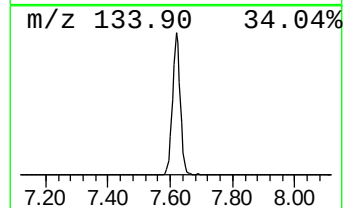
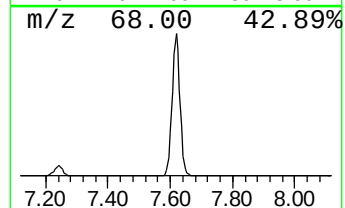
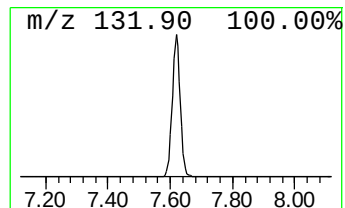
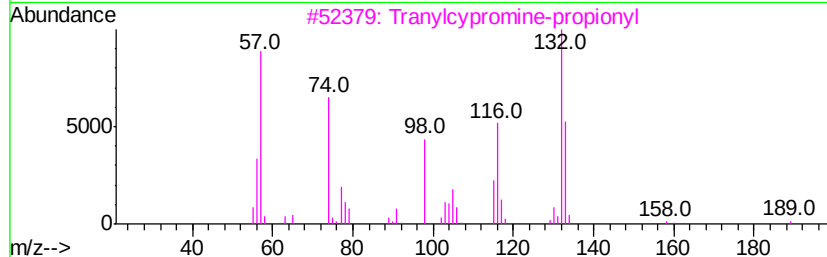
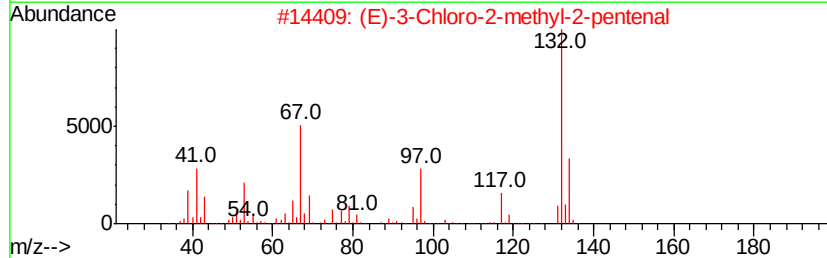
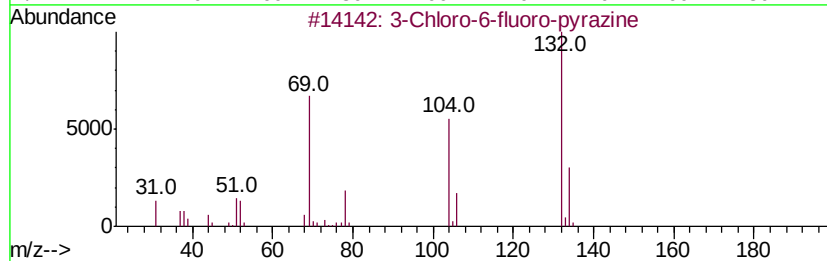
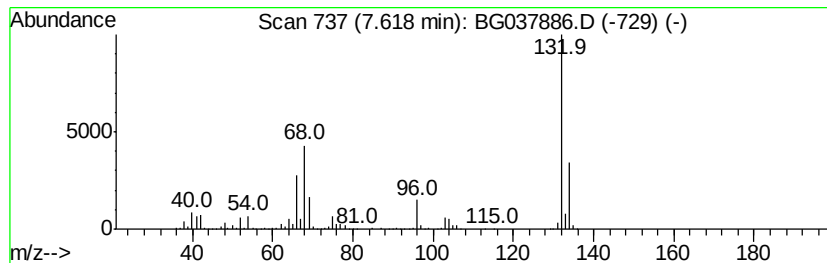
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Peak Number 2 unknown7.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	118.03 ng	1953190	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	12
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
5		Xenon	132	Xe	007440-63-3	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	86211	1	8.09	330955	20.0
unknown7.62	7.62	118.0	ng	1953190	1	8.09	330955	20.0