

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101218\
 Data File : BG037327.D
 Acq On : 13 Oct 2018 15:14
 Operator : JU/SJ
 Sample : PB113715BL
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB113715BL

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 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-BG101118.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	5.192	319	324	332	rBV	43376	71774	1.36%	0.290%
2	5.650	395	402	412	rBV	1115544	1836672	34.86%	7.426%
3	7.254	667	675	687	rBV	1172893	2042912	38.78%	8.260%
4	7.642	732	741	750	rBV	1040382	1878276	35.65%	7.594%
5	8.118	812	822	830	rBV	203327	364533	6.92%	1.474%
6	8.441	866	877	885	rBV	771257	1415046	26.86%	5.721%
7	9.293	1012	1022	1034	rBV	654909	1196924	22.72%	4.839%
8	10.938	1293	1302	1312	rBV	300130	539209	10.24%	2.180%
9	13.371	1708	1716	1724	rBV	1828226	2996013	56.87%	12.113%
10	14.751	1943	1951	1960	rVB2	486896	816081	15.49%	3.299%
11	16.232	2196	2203	2222	rBV	1513500	2414585	45.83%	9.762%
12	17.495	2410	2418	2426	rBV	747885	1148437	21.80%	4.643%
13	19.775	2801	2806	2811	rBV	84917	97692	1.85%	0.395%
14	20.098	2854	2861	2869	rBV	3701775	5268236	100.00%	21.300%
15	20.815	2979	2983	2987	rBV2	65362	76295	1.45%	0.308%
16	21.778	3140	3147	3159	rVB	772531	1307364	24.82%	5.286%
17	25.122	3706	3716	3731	rVB2	421226	1263638	23.99%	5.109%

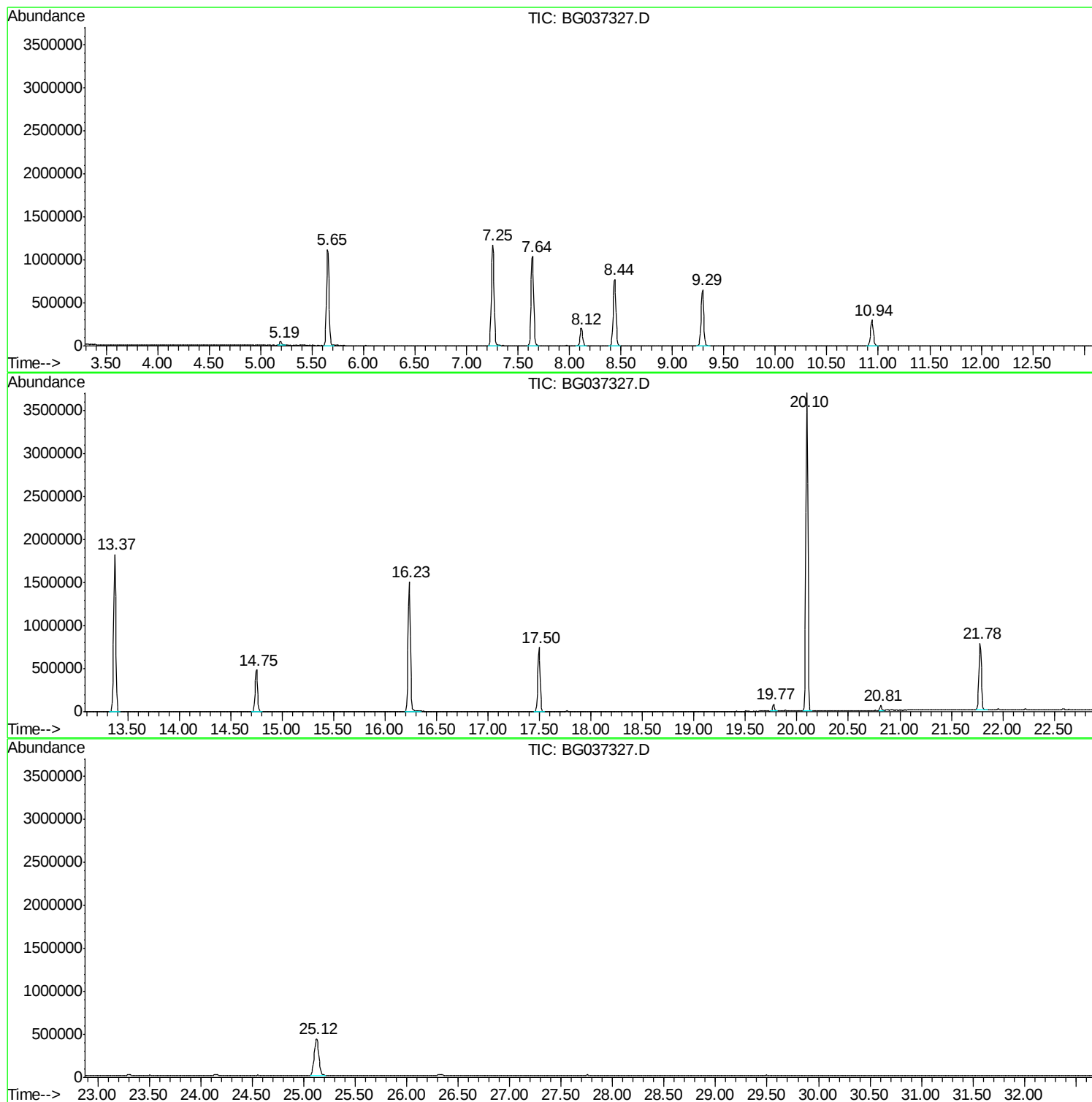
Sum of corrected areas: 24733687

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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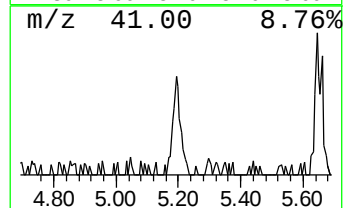
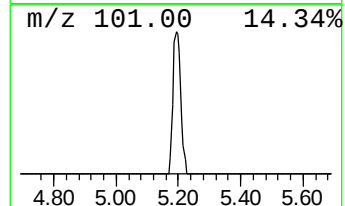
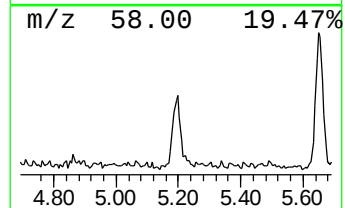
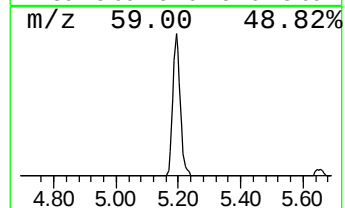
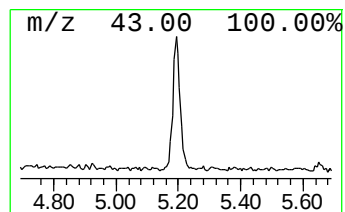
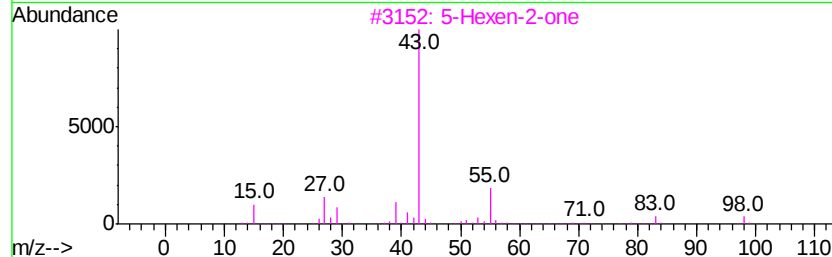
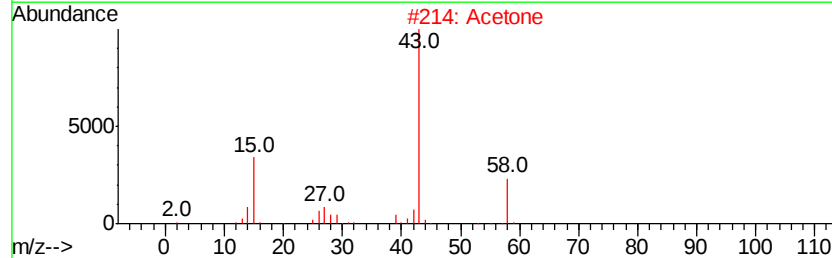
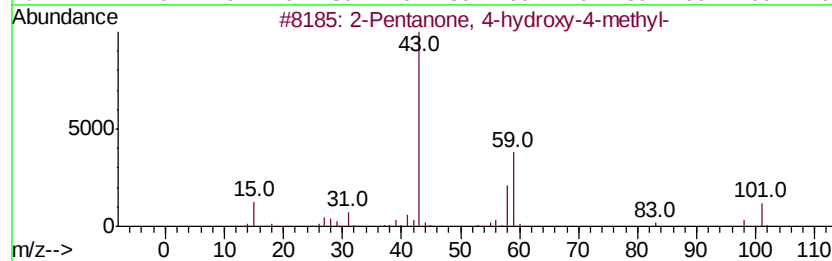
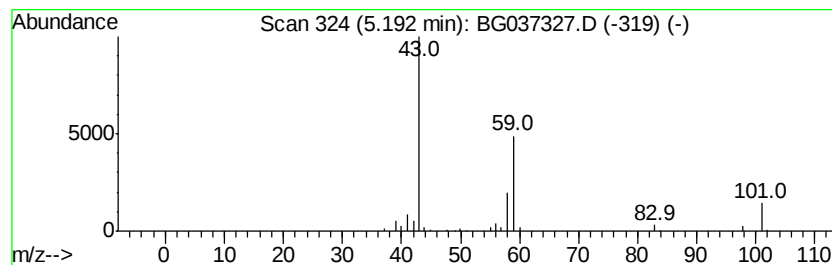
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	3.94 ng	71774	1,4-Dichlorobenzene-d4	8.12

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	83
2			Acetone	58	C3H6O	000067-64-1	9
3			5-Hexen-2-one	98	C6H10O	000109-49-9	9
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9



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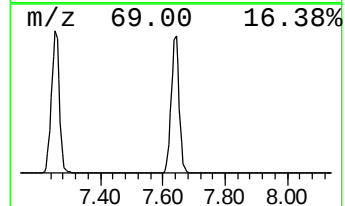
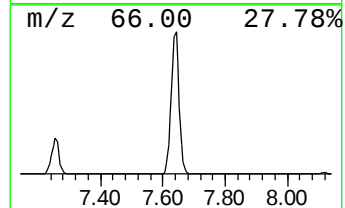
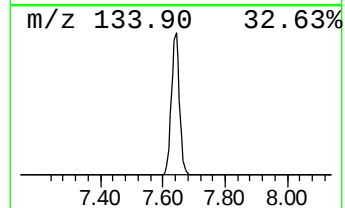
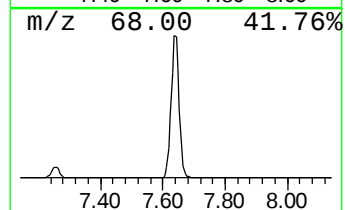
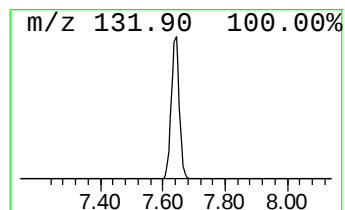
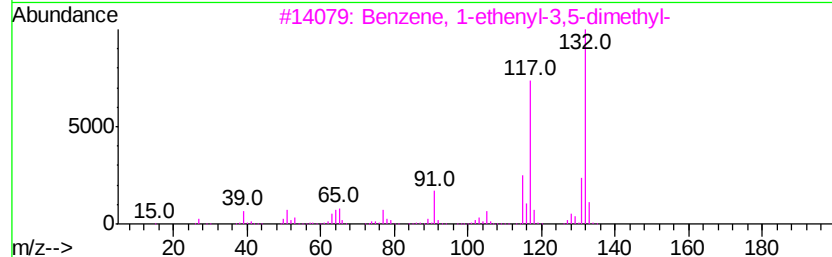
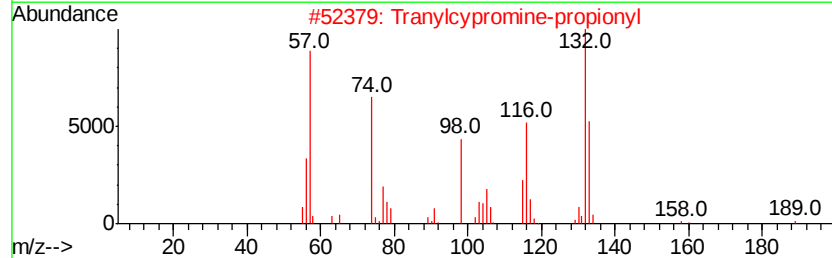
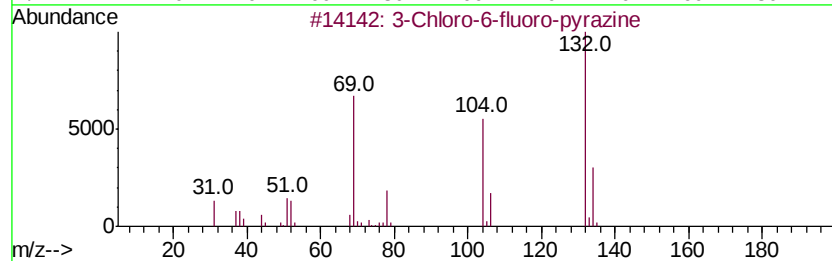
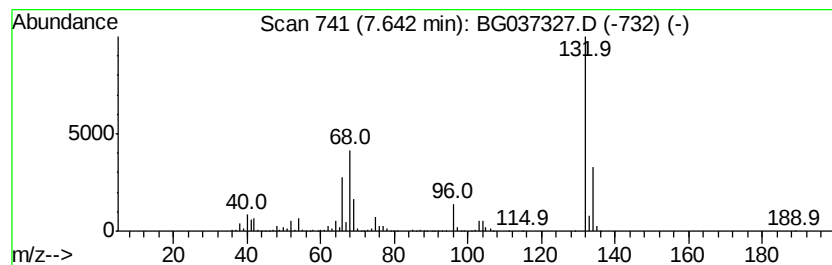
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Peak Number 2 unknown7.64 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.64	103.05 ng	1878280	1,4-Dichlorobenzene-d4	8.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
4		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.19	3.9	ng	71774	1	8.12	364533	20.0
unknown7.64	7.64	103.1	ng	1878280	1	8.12	364533	20.0