

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG121117\
 Data File : BG031466.D
 Acq On : 11 Dec 2017 21:44
 Operator : SJ/JU
 Sample : I6762-04
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 120417-FOPI-F-U-S

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:\HPCHEM1\BNA_G\METHODS\8270-BG121117.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	5.186	317	323	334	rBV	53364	86790	1.63%	0.483%
2	5.668	398	405	424	rBV	255758	479457	9.01%	2.667%
3	7.277	671	679	698	rBV2	157723	328026	6.16%	1.825%
4	7.642	733	741	761	rBV	492061	901329	16.93%	5.014%
5	8.112	815	821	833	rBV	164517	297236	5.58%	1.654%
6	8.435	868	876	888	rBV	681575	1248885	23.46%	6.948%
7	9.275	1011	1019	1034	rBV	466524	953614	17.91%	5.305%
8	10.926	1293	1300	1323	rBV	178679	415952	7.81%	2.314%
9	13.353	1706	1713	1736	rBV	1607625	2657467	49.92%	14.784%
10	14.733	1941	1948	1960	rBV	375577	652472	12.26%	3.630%
11	16.220	2190	2201	2215	rBV	774995	1339657	25.17%	7.453%
12	17.471	2408	2414	2431	rBV2	557399	927445	17.42%	5.160%
13	20.062	2849	2855	2871	rBV	3844772	5323147	100.00%	29.614%
14	21.743	3133	3141	3155	rBV	716845	1244283	23.37%	6.922%
15	25.021	3689	3699	3717	rBV3	360482	1119547	21.03%	6.228%

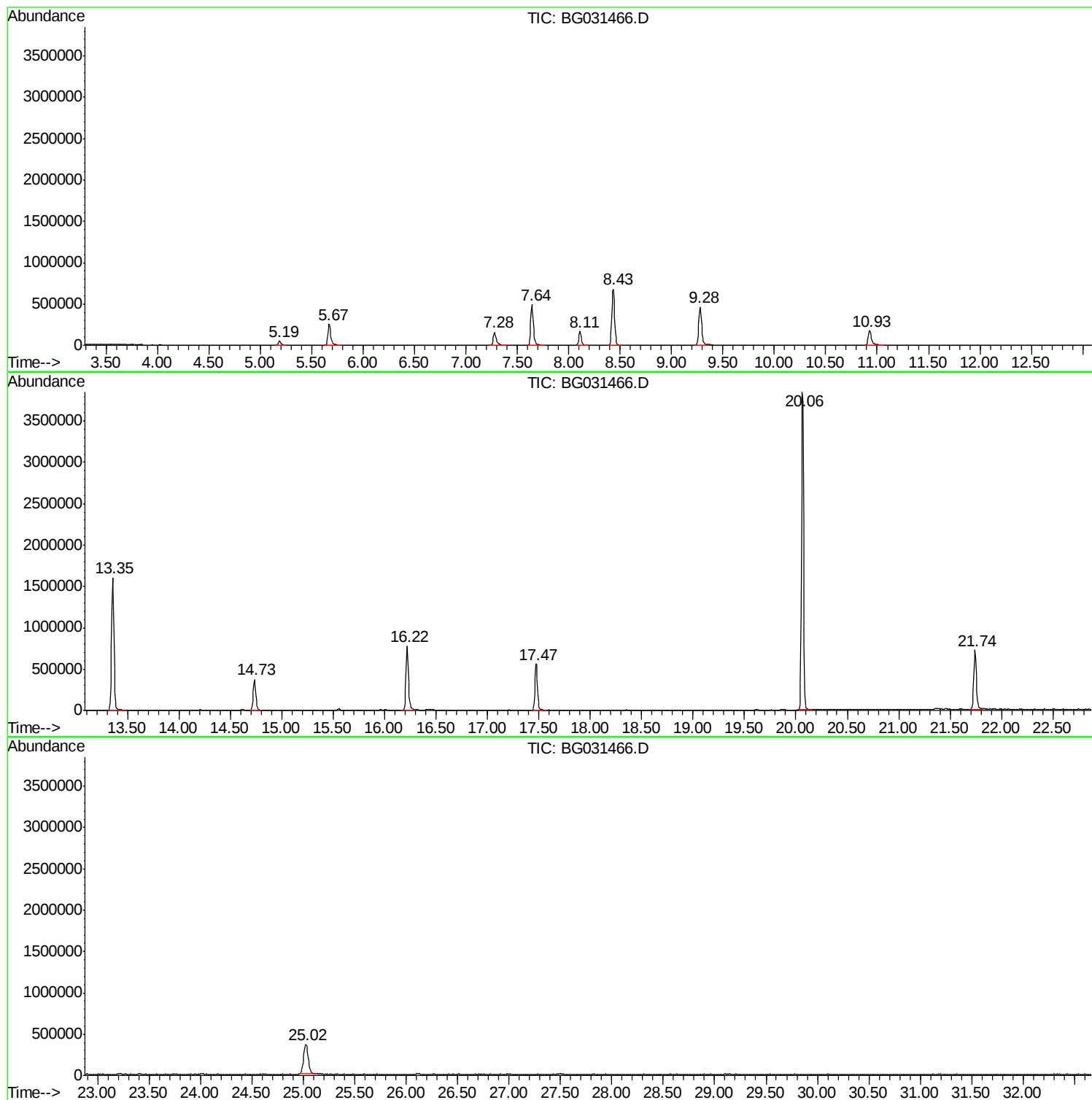
Sum of corrected areas: 17975307

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG121117.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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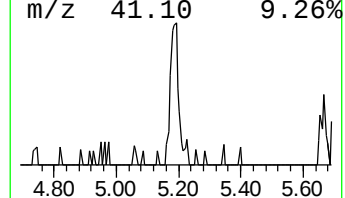
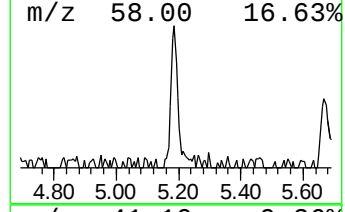
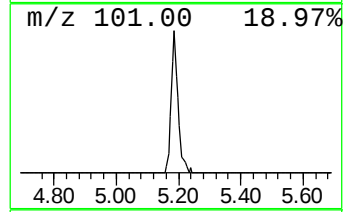
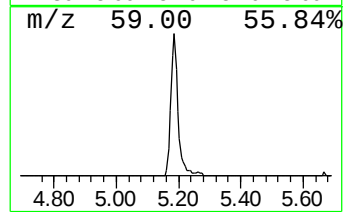
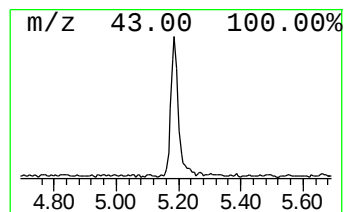
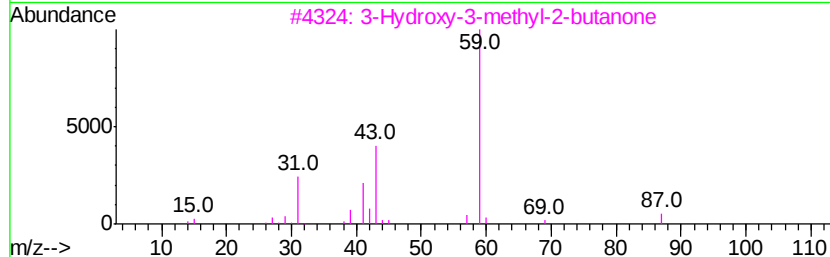
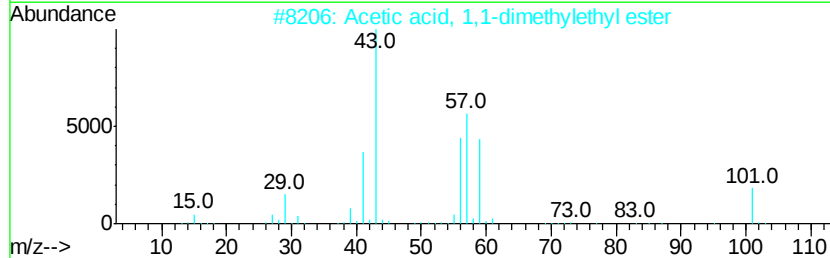
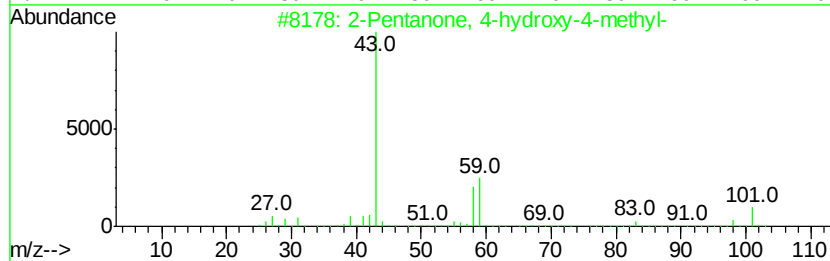
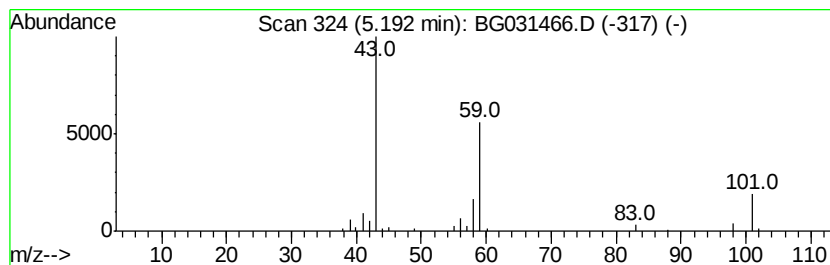
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG121117.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	5.84 ng	86790	1,4-Dichlorobenzene-d4	8.11

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	25
4			Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	17
5			Acetic acid, chloro-, 1,1-dimeth...	150	C6H11ClO2	000107-59-5	17



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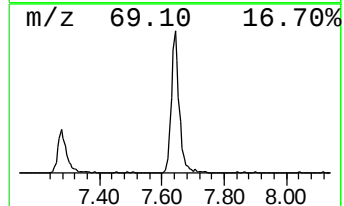
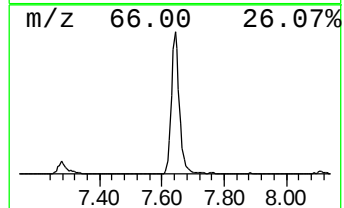
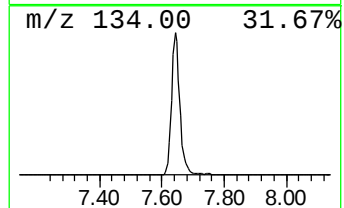
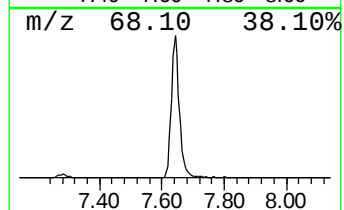
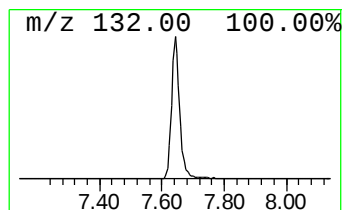
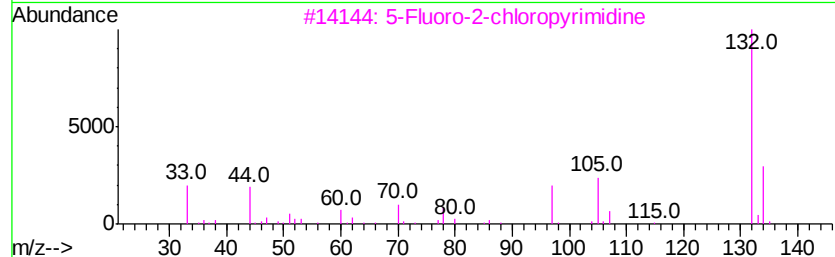
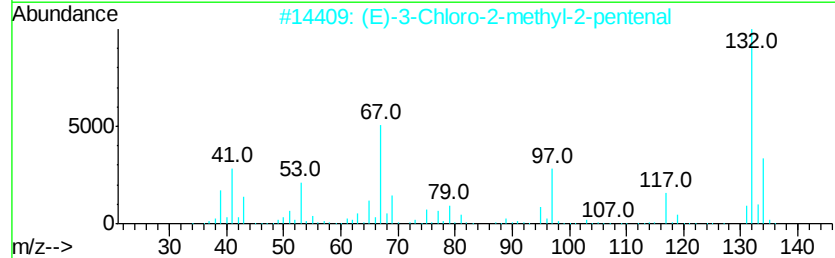
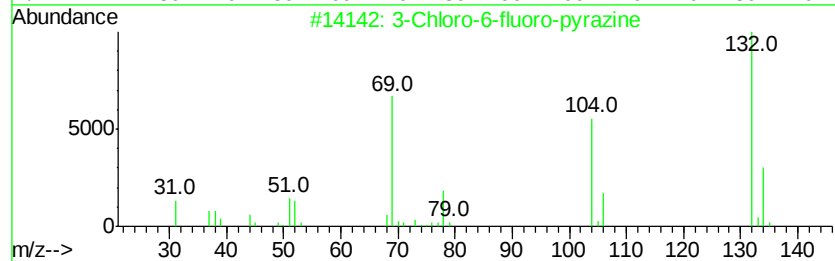
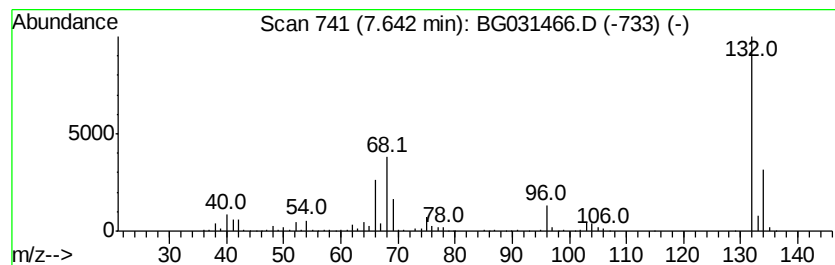
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Peak Number 2 unknown7.64 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.64	60.65 ng	901329	1,4-Dichlorobenzene-d4	8.11

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	17
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.19	5.8	ng	86790	1	8.11	297236	20.0
unknown7.64	7.64	60.6	ng	901329	1	8.11	297236	20.0