

Data Path : Z:\HPCHEM1\BNA G\DATA\BG113017\
 Data File : BG031303.D
 Acq On : 1 Dec 2017 00:45
 Operator : SJ/JU
 Sample : I6635-03
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 112817-JETT-E-U-B

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-BG111017.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	5.209	321	327	333	rBV	52156	79444	1.31%	0.373%
2	5.691	403	409	426	rBV	272639	494595	8.15%	2.325%
3	7.295	676	682	695	rBV	154496	326649	5.38%	1.535%
4	7.665	738	745	757	rBV	558848	1014944	16.72%	4.770%
5	8.129	817	824	834	rBV	212804	370179	6.10%	1.740%
6	8.452	871	879	891	rBV	827449	1528376	25.17%	7.183%
7	9.298	1015	1023	1036	rBV	568881	1126447	18.55%	5.294%
8	10.949	1295	1304	1327	rBV	240358	526279	8.67%	2.474%
9	13.376	1709	1717	1736	rBV	1985774	3266899	53.81%	15.354%
10	14.751	1944	1951	1967	rBV3	465575	843139	13.89%	3.963%
11	16.237	2198	2204	2227	rBV	996409	1700760	28.01%	7.994%
12	17.494	2411	2418	2435	rBV2	688215	1120828	18.46%	5.268%
13	20.085	2852	2859	2871	rBV	4538431	6071388	100.00%	28.535%
14	21.766	3138	3145	3154	rBV	844692	1446475	23.82%	6.798%
15	25.062	3695	3706	3718	rBV2	452811	1360231	22.40%	6.393%

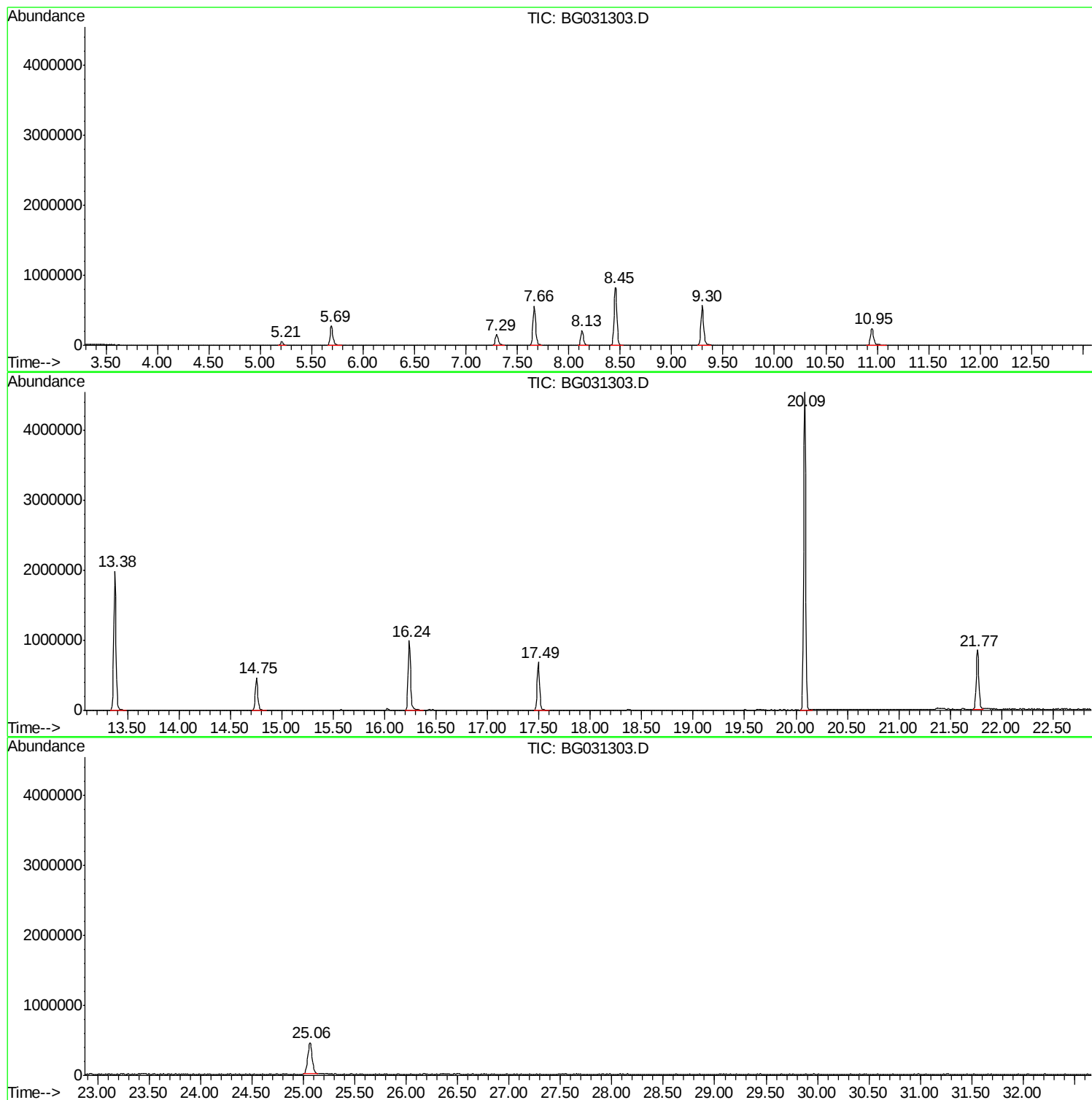
Sum of corrected areas: 21276633

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG111017.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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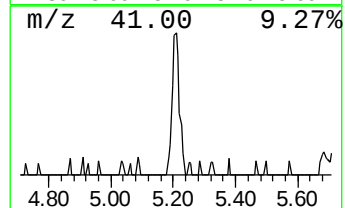
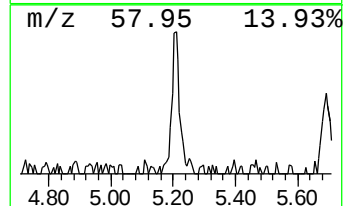
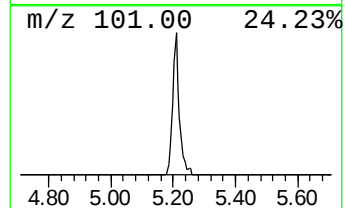
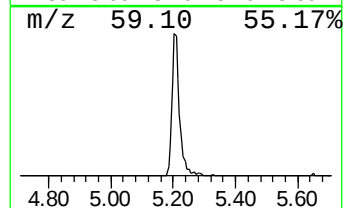
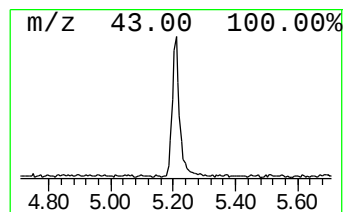
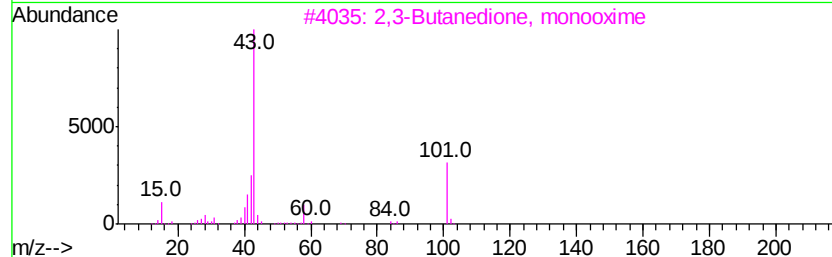
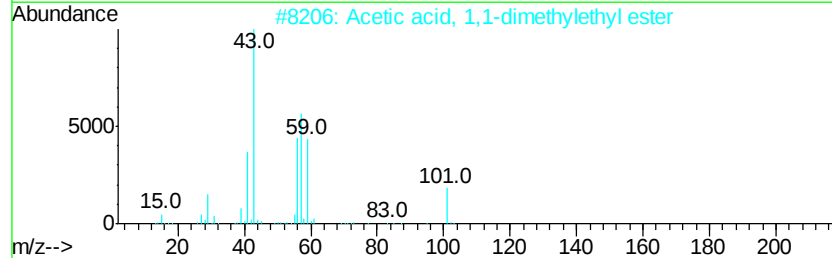
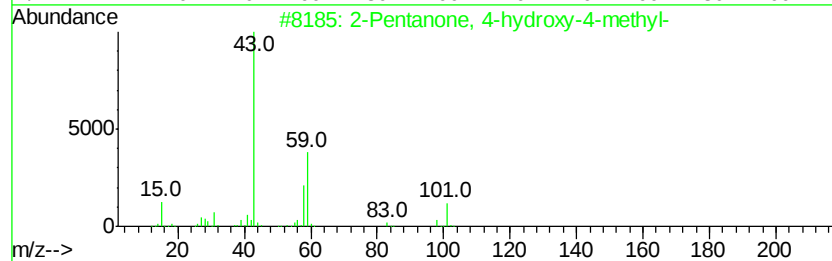
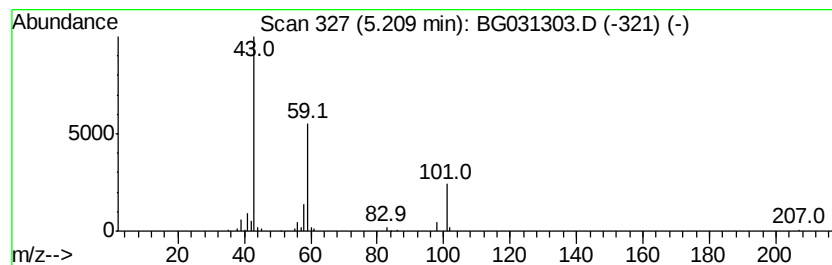
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG111017.M
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.21	4.29 ng	79444	1,4-Dichlorobenzene-d4	8.13

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	38
4			Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	28
5			2,2,3,3,5,8,11-Heptamethyl-4,7,1...	348	C18H40O4Si	1000367-07-2	28



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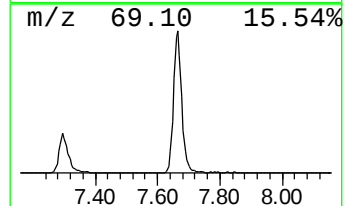
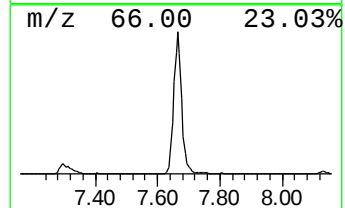
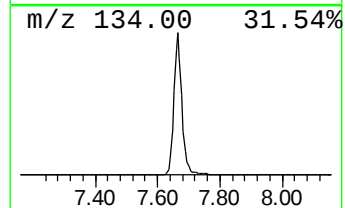
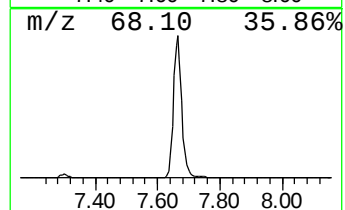
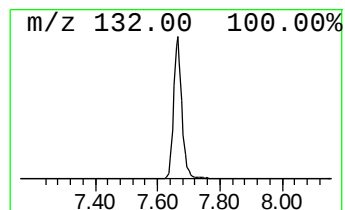
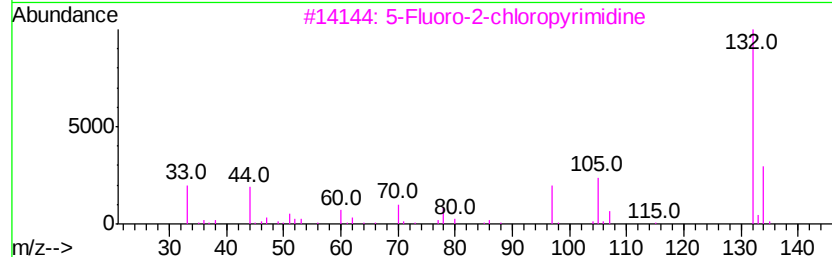
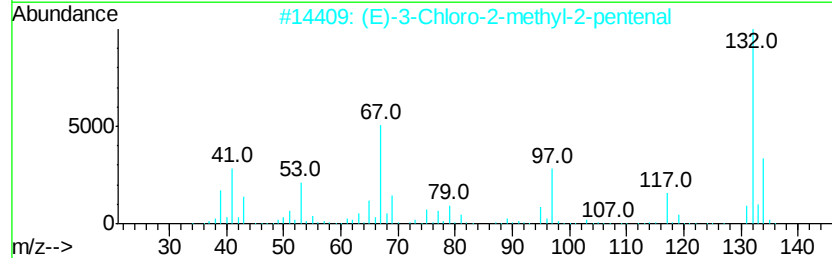
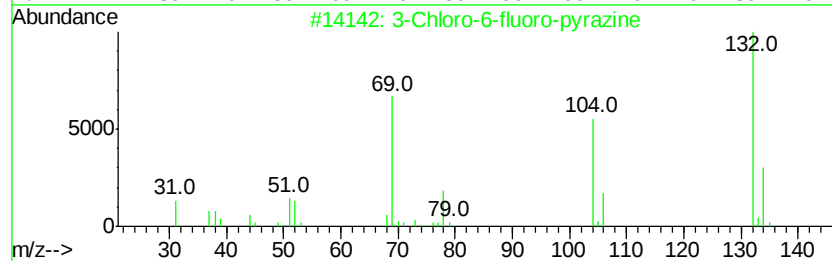
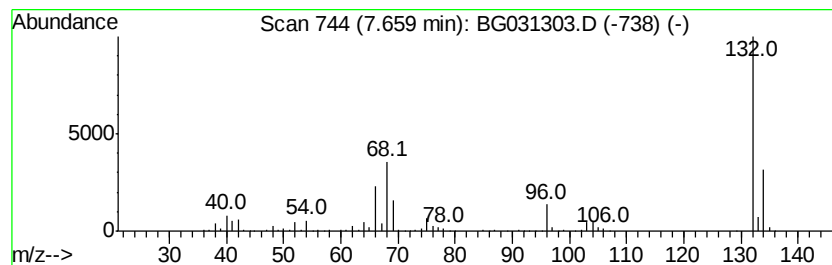
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Peak Number 2 unknown7.66 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.66	54.84 ng	1014940	1,4-Dichlorobenzene-d4	8.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	25
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.21	4.3	ng	79444	1	8.13	370179	20.0
unknown7.66	7.66	54.8	ng	1014940	1	8.13	370179	20.0