

Data Path : Z:\HPCHEM1\BNA G\DATA\BG113017\
 Data File : BG031293.D
 Acq On : 30 Nov 2017 18:30
 Operator : SJ/JU
 Sample : I6634-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 112717-JETT-E-D-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	5.208	321	327	336	rBV	56631	93582	1.46%	0.421%
2	5.690	402	409	428	rBV	308089	563422	8.79%	2.536%
3	7.294	675	682	704	rBV	187435	387854	6.05%	1.746%
4	7.664	737	745	756	rBV	601578	1084231	16.92%	4.881%
5	8.128	817	824	833	rBV	201081	358157	5.59%	1.612%
6	8.457	872	880	891	rBV	840696	1553945	24.25%	6.995%
7	9.298	1015	1023	1036	rBV	626871	1211581	18.91%	5.454%
8	10.949	1296	1304	1325	rBV	242070	532110	8.30%	2.395%
9	13.375	1710	1717	1730	rBV	2123841	3487953	54.44%	15.701%
10	14.750	1944	1951	1966	rVB2	466901	844715	13.18%	3.802%
11	16.237	2198	2204	2221	rBV	1020211	1752885	27.36%	7.890%
12	17.494	2410	2418	2433	rBV2	710494	1138305	17.77%	5.124%
13	20.085	2853	2859	2876	rBV	4734779	6407270	100.00%	28.842%
14	21.765	3137	3145	3156	rBV2	801098	1443582	22.53%	6.498%
15	25.062	3695	3706	3724	rBV2	450790	1355794	21.16%	6.103%

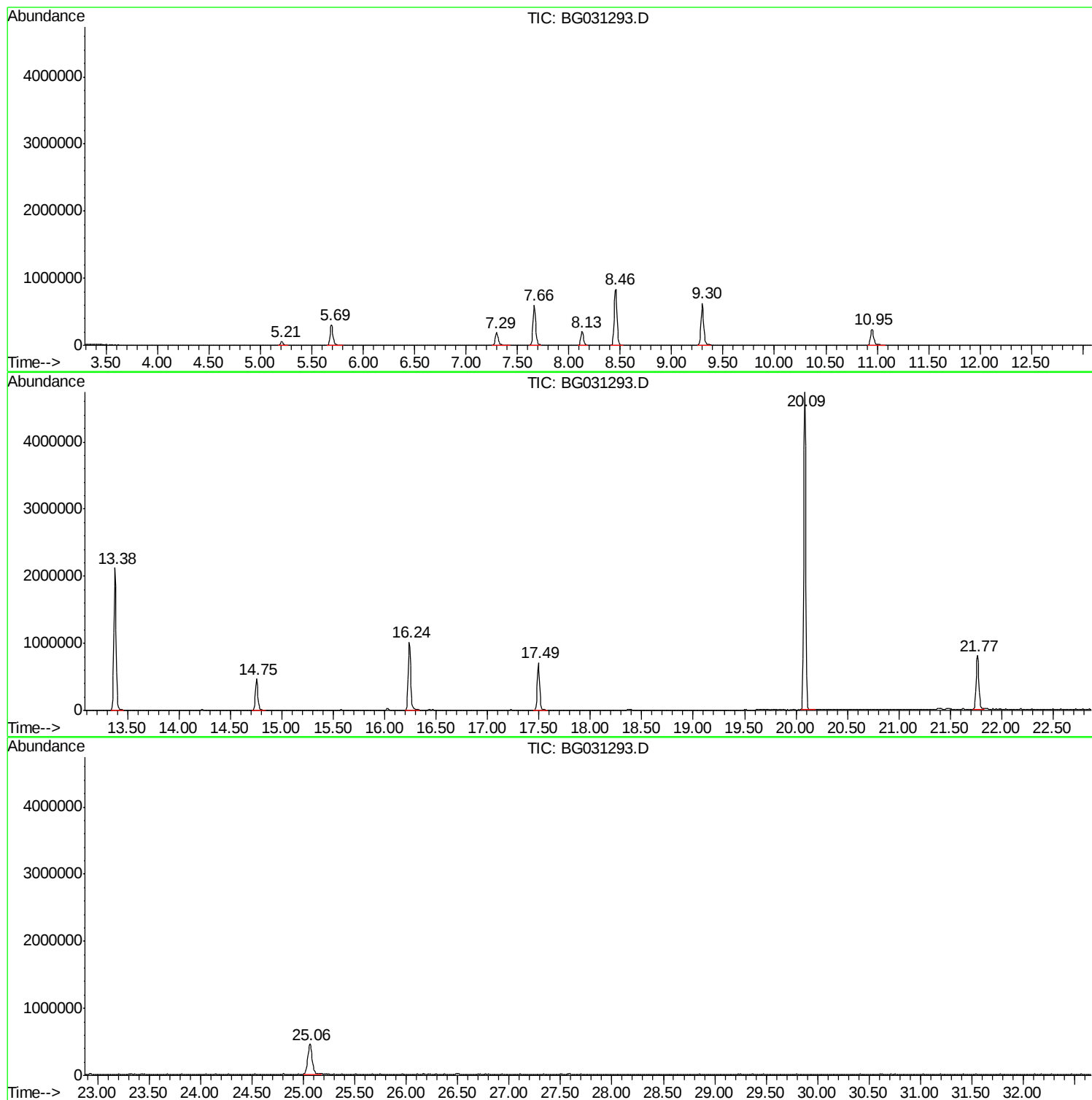
Sum of corrected areas: 22215386

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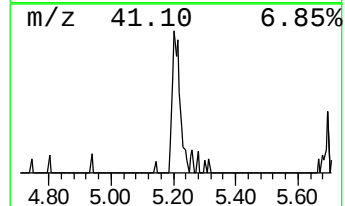
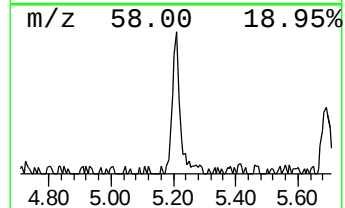
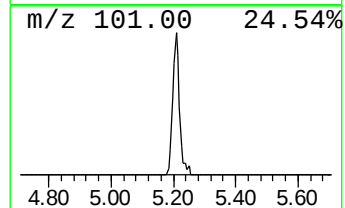
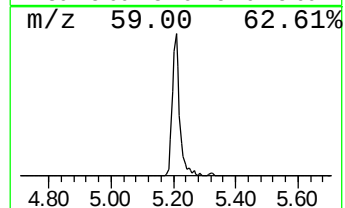
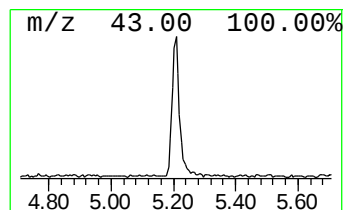
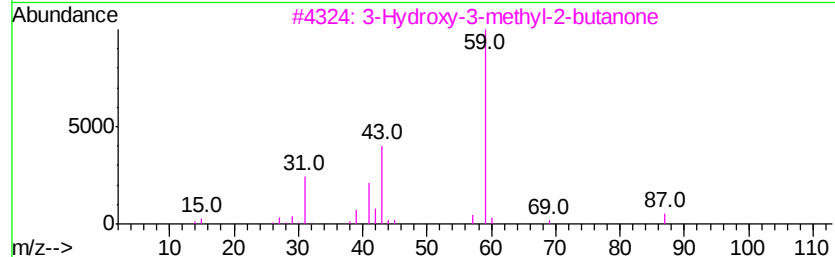
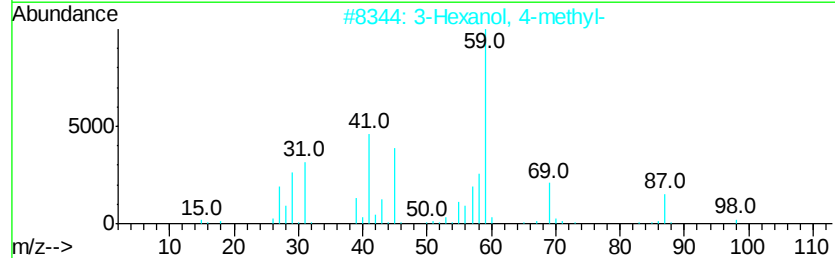
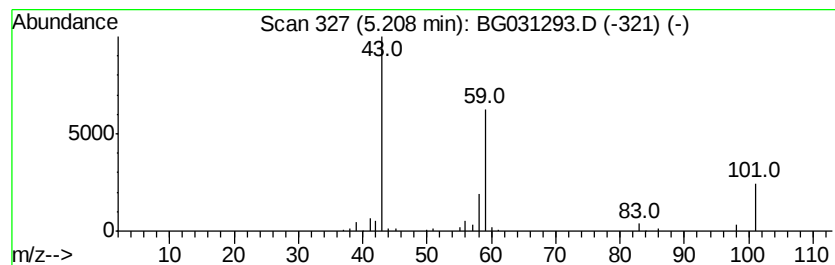
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.21	5.23 ng	93582	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	40
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
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			Propane, 2-methoxy-	74	C4H10O	000598-53-8	12



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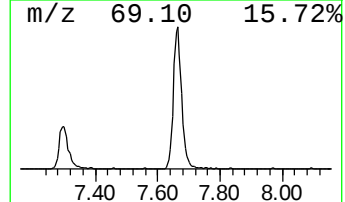
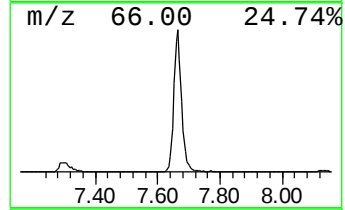
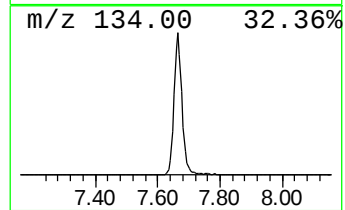
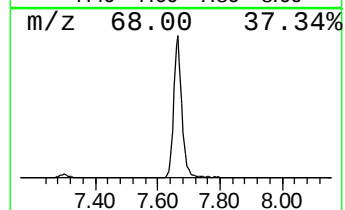
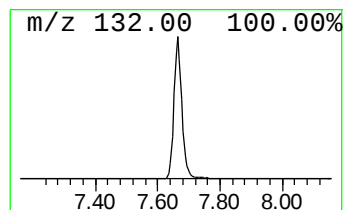
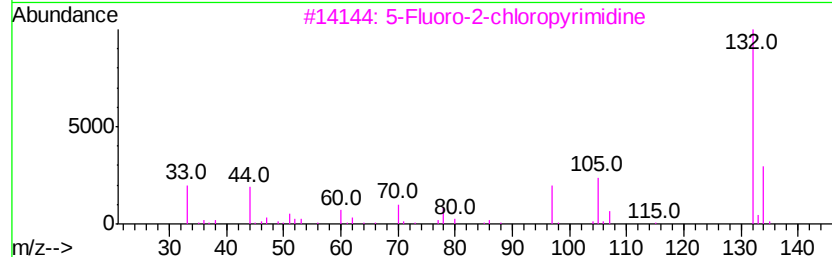
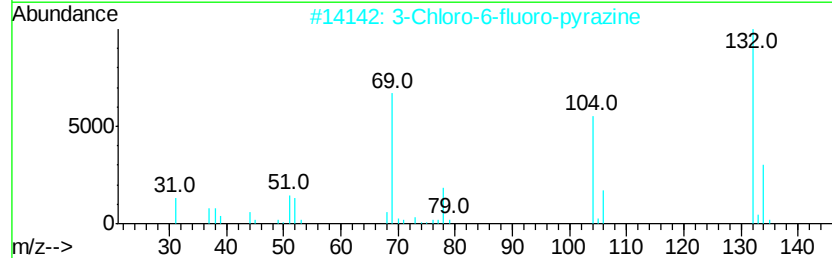
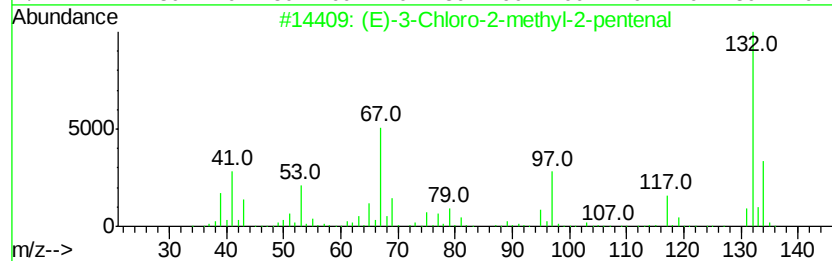
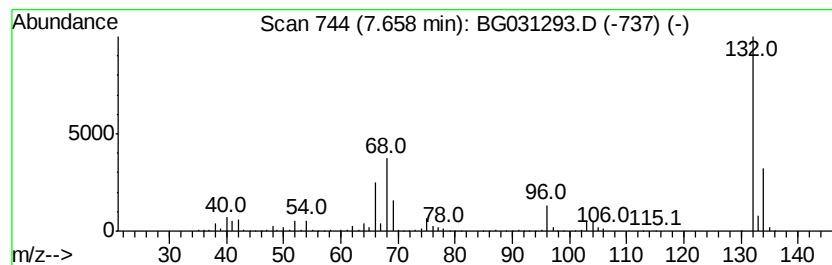
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Peak Number 2 unknown7.66 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.66	60.55 ng	1084230	1,4-Dichlorobenzene-d4	8.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	25
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.21	5.2	ng	93582	1	8.13	358157	20.0
unknown7.66	7.66	60.5	ng	1084230	1	8.13	358157	20.0