

Data Path : Z:\HPCHEM1\BNA G\DATA\BG113017\
 Data File : BG031294.D
 Acq On : 30 Nov 2017 19:07
 Operator : SJ/JU
 Sample : I6634-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 112717-JETT-E-D-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-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	5.210	320	327	340	rBV	70059	114029	1.68%	0.464%
2	5.691	402	409	421	rBV	361505	639195	9.45%	2.604%
3	7.295	674	682	694	rBV	211594	427968	6.32%	1.743%
4	7.665	737	745	756	rBV	683236	1256755	18.57%	5.119%
5	8.130	817	824	832	rBV2	239343	418605	6.19%	1.705%
6	8.453	871	879	895	rVB	1041426	1879534	27.77%	7.656%
7	9.299	1015	1023	1034	rBV	742738	1416302	20.93%	5.769%
8	10.950	1295	1304	1314	rBV	271258	579042	8.56%	2.359%
9	13.376	1709	1717	1737	rBV	2500134	4048728	59.83%	16.492%
10	14.751	1942	1951	1967	rBV2	520330	934116	13.80%	3.805%
11	16.238	2197	2204	2223	rBV	1178915	1971150	29.13%	8.029%
12	17.495	2410	2418	2432	rBV2	747005	1223882	18.08%	4.985%
13	20.086	2852	2859	2871	rBV	4837338	6767389	100.00%	27.567%
14	21.767	3138	3145	3155	rBV	865913	1481432	21.89%	6.035%
15	25.063	3695	3706	3724	rBV3	454461	1391076	20.56%	5.666%

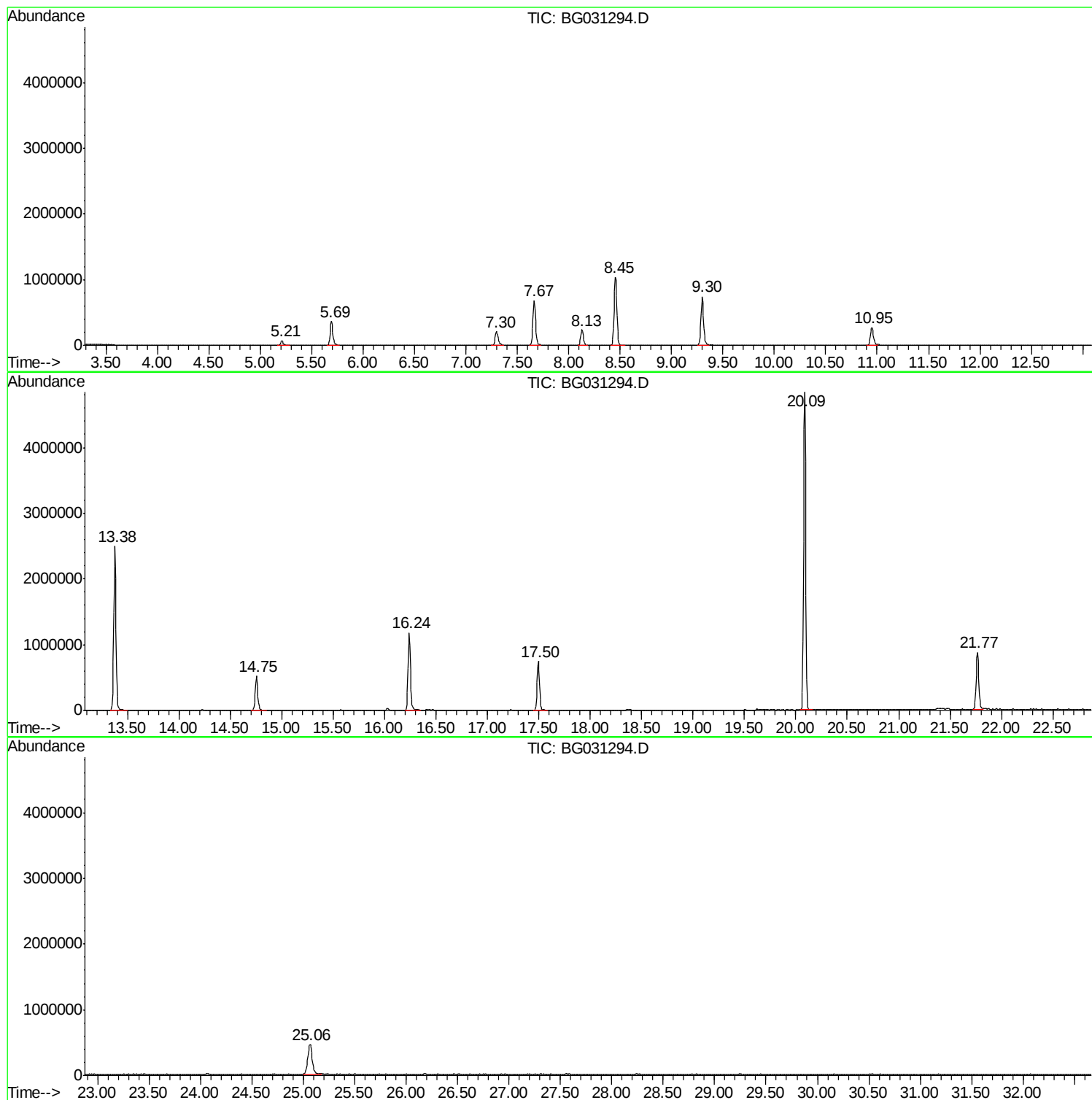
Sum of corrected areas: 24549203

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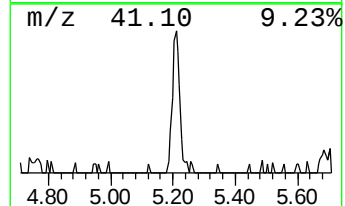
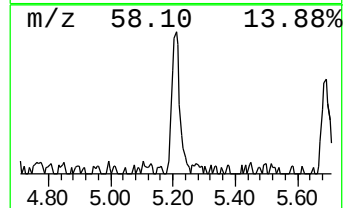
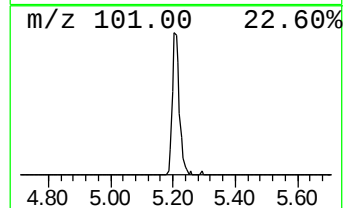
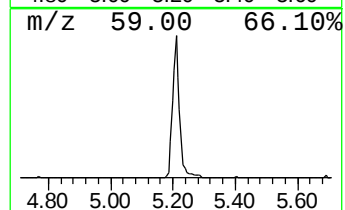
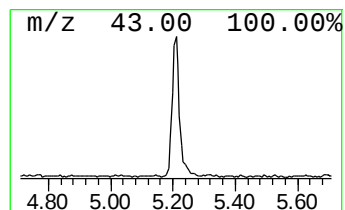
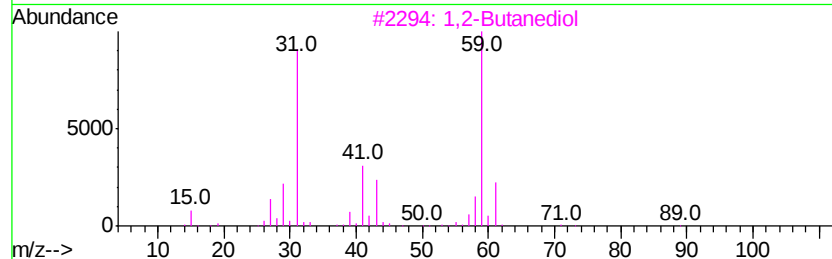
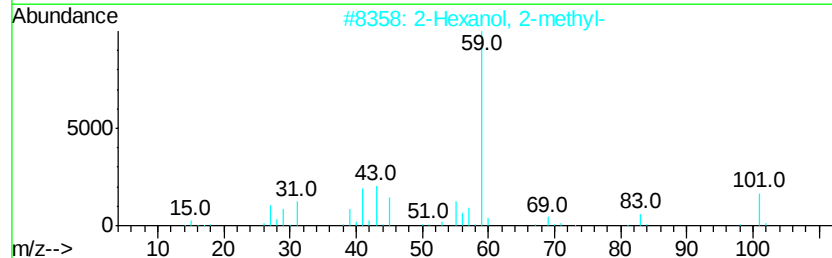
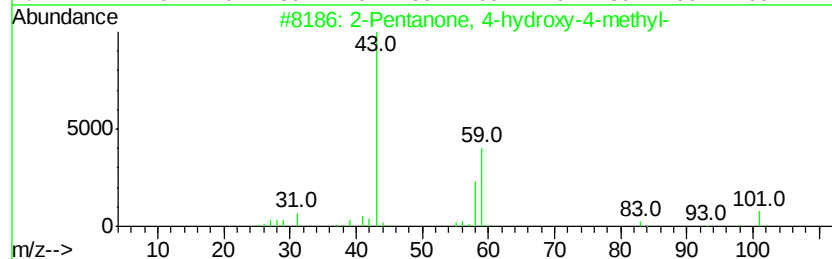
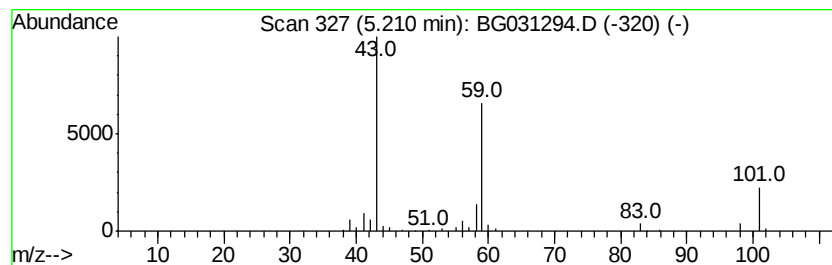
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.21	5.45 ng	114029	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	50
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
3			1,2-Butanediol	90	C4H10O2	000584-03-2	17
4			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	17
5			Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	17



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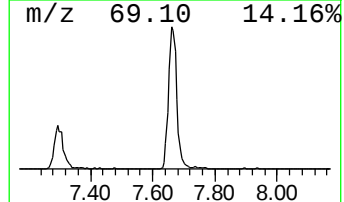
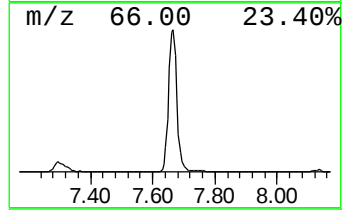
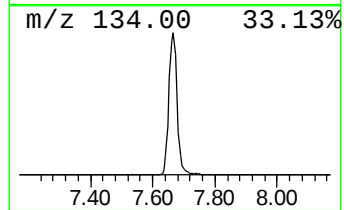
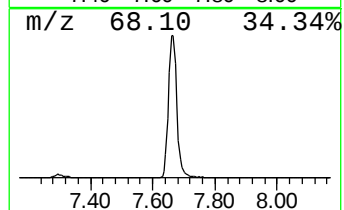
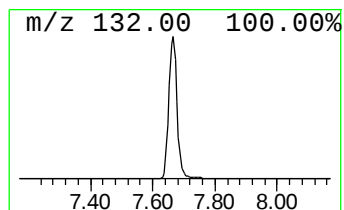
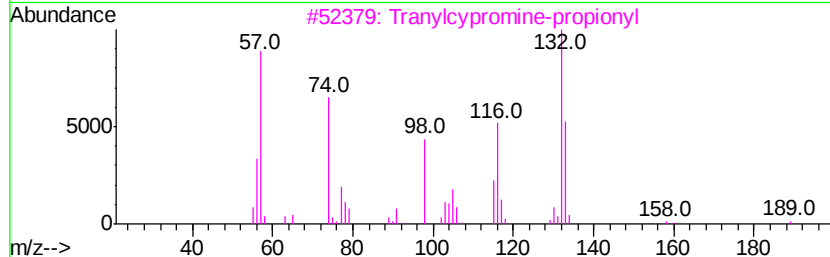
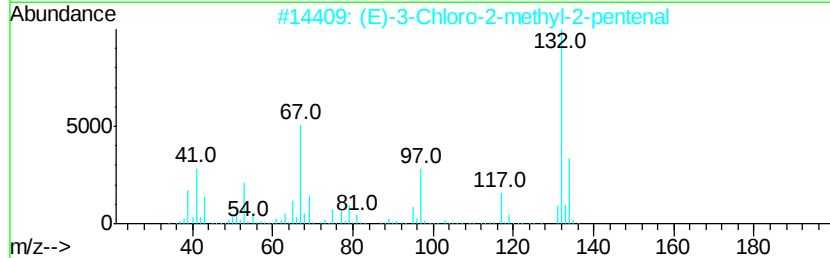
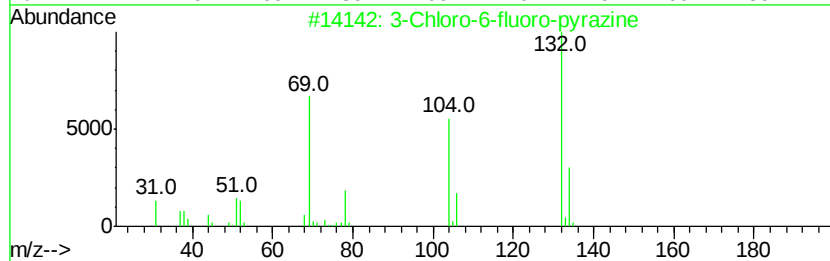
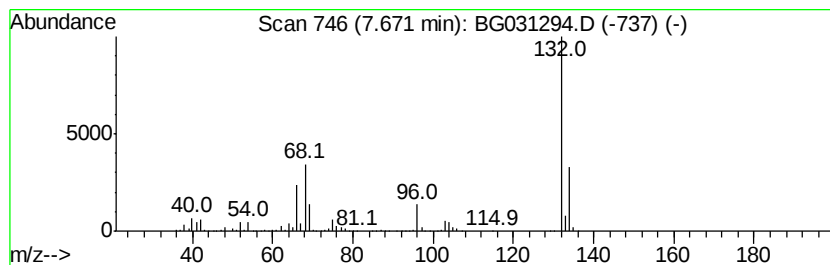
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Peak Number 2 unknown7.67 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.67	60.04 ng	1256760	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	17
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.21	5.5	ng	114029	1	8.13	418605	20.0
unknown7.67	7.67	60.0	ng	1256760	1	8.13	418605	20.0