

Data Path : Z:\HPCHEM1\BNA G\DATA\BG113017\
 Data File : BG031302.D
 Acq On : 1 Dec 2017 00:08
 Operator : SJ/JU
 Sample : I6635-02
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 112817-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	5.209	322	327	338	rBV	62353	102312	1.62%	0.467%
2	5.690	400	409	423	rBV	300939	533422	8.46%	2.435%
3	7.294	675	682	694	rBV	172704	359705	5.70%	1.642%
4	7.665	738	745	756	rBV	573707	1035633	16.42%	4.727%
5	8.129	816	824	834	rBV	211899	375821	5.96%	1.715%
6	8.452	871	879	895	rVB	859455	1571069	24.91%	7.171%
7	9.298	1014	1023	1037	rBV	621974	1203436	19.08%	5.493%
8	10.949	1297	1304	1315	rBV	241237	514668	8.16%	2.349%
9	13.375	1709	1717	1729	rBV	2108889	3406225	54.01%	15.548%
10	14.750	1944	1951	1964	rBV2	463500	840451	13.33%	3.836%
11	16.237	2197	2204	2221	rBV	990555	1727264	27.39%	7.884%
12	17.494	2411	2418	2433	rBV2	705617	1136639	18.02%	5.188%
13	20.085	2852	2859	2873	rBV	4698952	6306916	100.00%	28.789%
14	21.766	3138	3145	3159	rBV	850693	1442095	22.87%	6.583%
15	25.062	3695	3706	3723	rBV2	450621	1352083	21.44%	6.172%

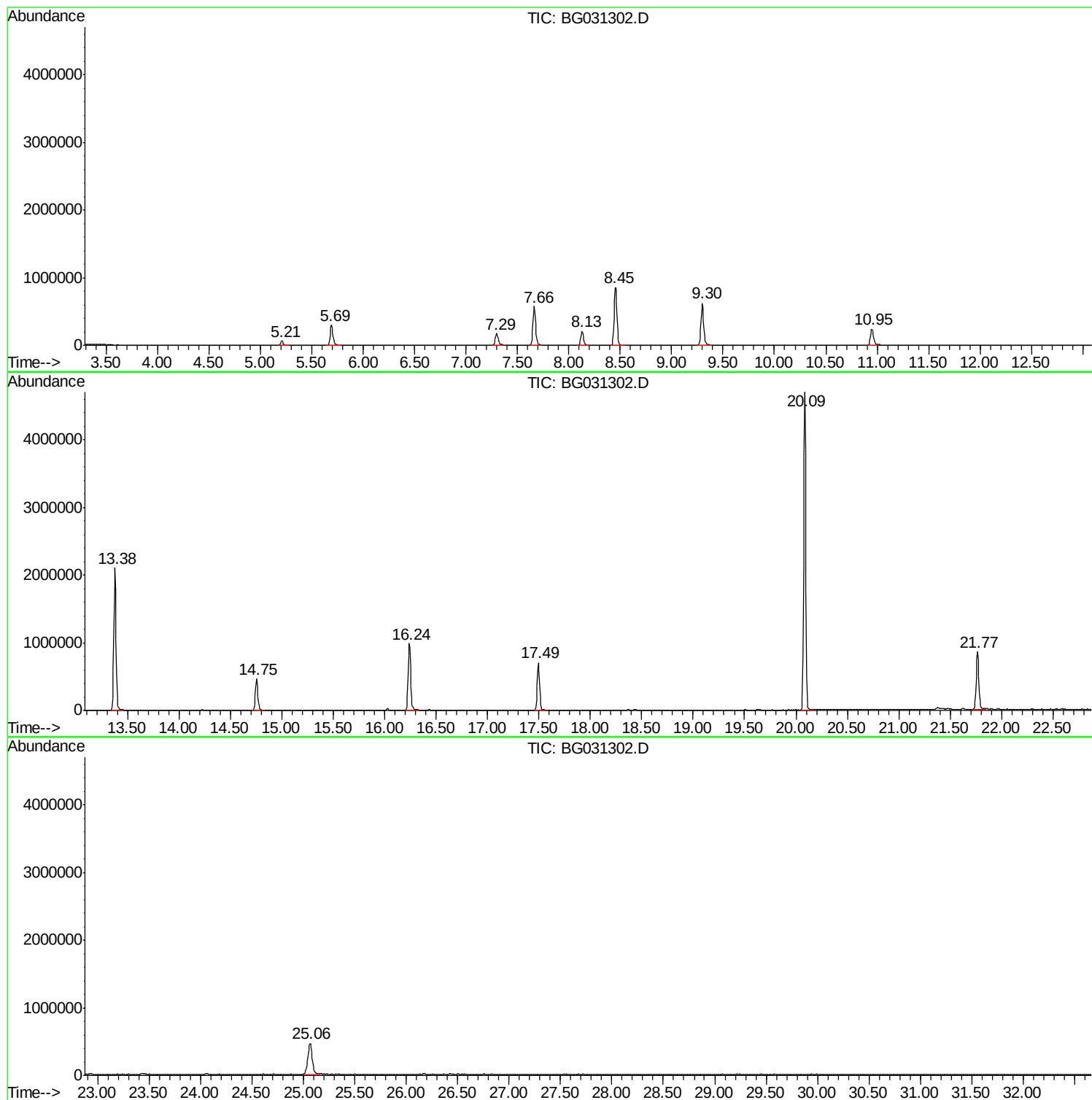
Sum of corrected areas: 21907739

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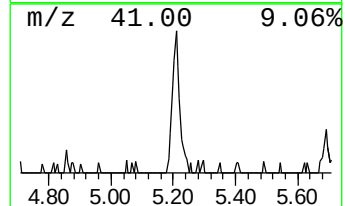
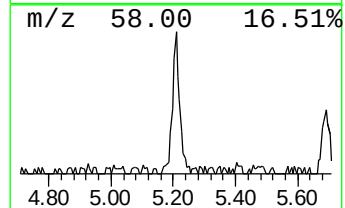
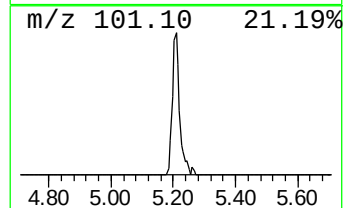
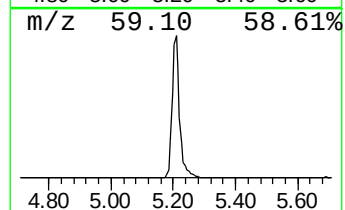
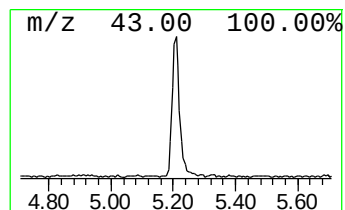
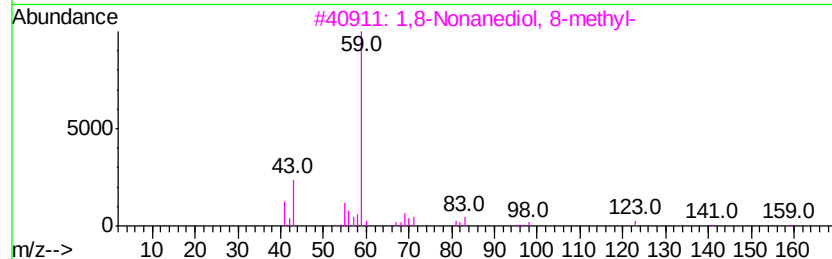
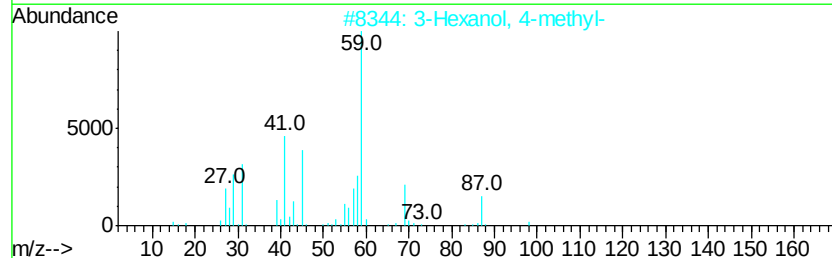
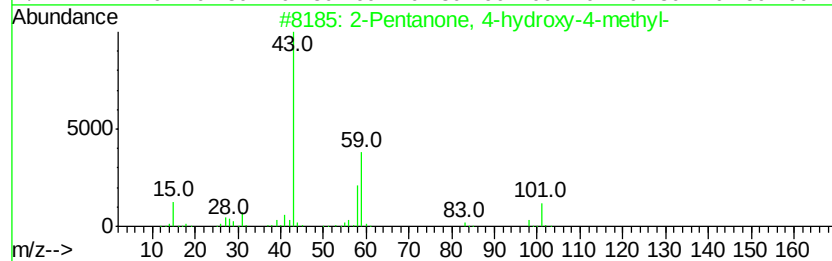
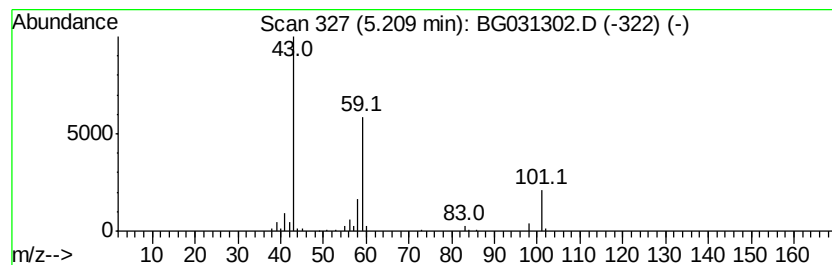
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.21	5.44 ng	102312	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	25
3	1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	23
4	Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	17
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16



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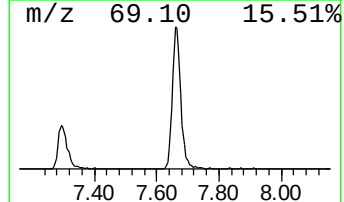
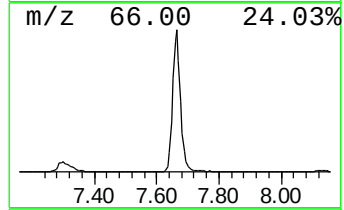
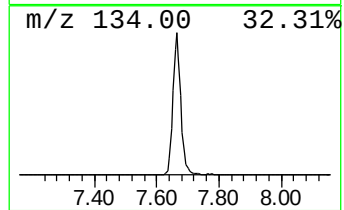
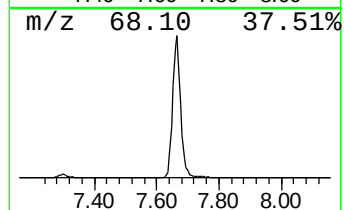
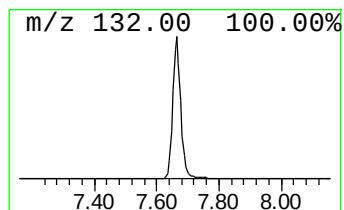
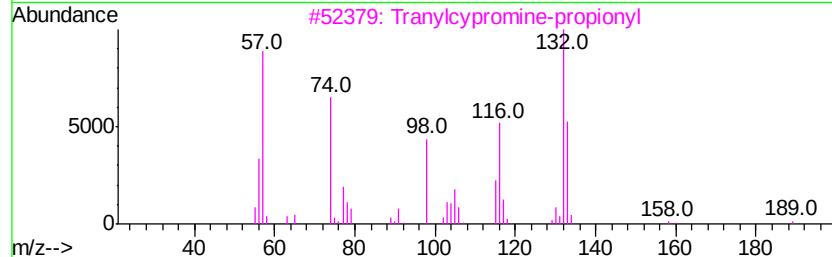
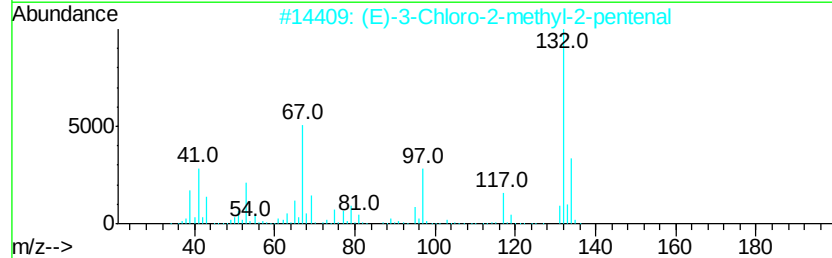
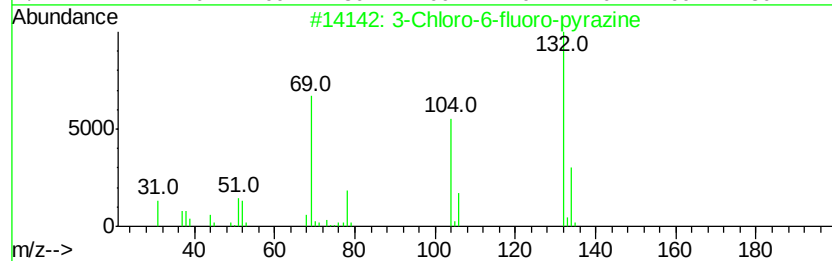
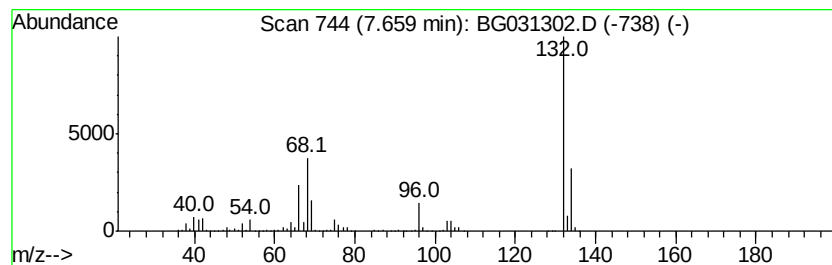
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Peak Number 2 unknown7.66 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.66	55.11 ng	1035630	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	12
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.21	5.4	ng	102312	1	8.13	375821	20.0
unknown7.66	7.66	55.1	ng	1035630	1	8.13	375821	20.0