

Data Path : S:\HPCHEM1\BNA_G\DATA\BG120817\
 Data File : BG031419.D
 Acq On : 8 Dec 2017 12:19
 Operator : SJ/JU
 Sample : I6731-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 113017-TIDO-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	5.203	321	326	335	rVB	39906	63155	1.06%	0.333%
2	5.685	402	408	422	rBV	226634	415707	7.00%	2.190%
3	7.295	675	682	696	rBV	151217	314022	5.28%	1.655%
4	7.659	737	744	760	rBV	464619	837497	14.09%	4.413%
5	8.123	816	823	832	rBV	146610	261171	4.40%	1.376%
6	8.453	871	879	891	rBV	615456	1147451	19.31%	6.046%
7	9.293	1015	1022	1036	rBV	457414	942279	15.86%	4.965%
8	10.944	1295	1303	1321	rBV	170261	385260	6.48%	2.030%
9	13.370	1709	1716	1737	rBV	1664075	2758210	46.42%	14.533%
10	14.751	1944	1951	1959	rBV2	377838	667969	11.24%	3.520%
11	16.238	2197	2204	2216	rBV	979432	1707739	28.74%	8.998%
12	17.489	2411	2417	2431	rBV2	627540	1040919	17.52%	5.485%
13	20.080	2852	2858	2873	rBV	4270841	5942357	100.00%	31.311%
14	21.766	3137	3145	3157	rBV	713402	1300022	21.88%	6.850%
15	25.062	3694	3706	3721	rBV2	388069	1194739	20.11%	6.295%

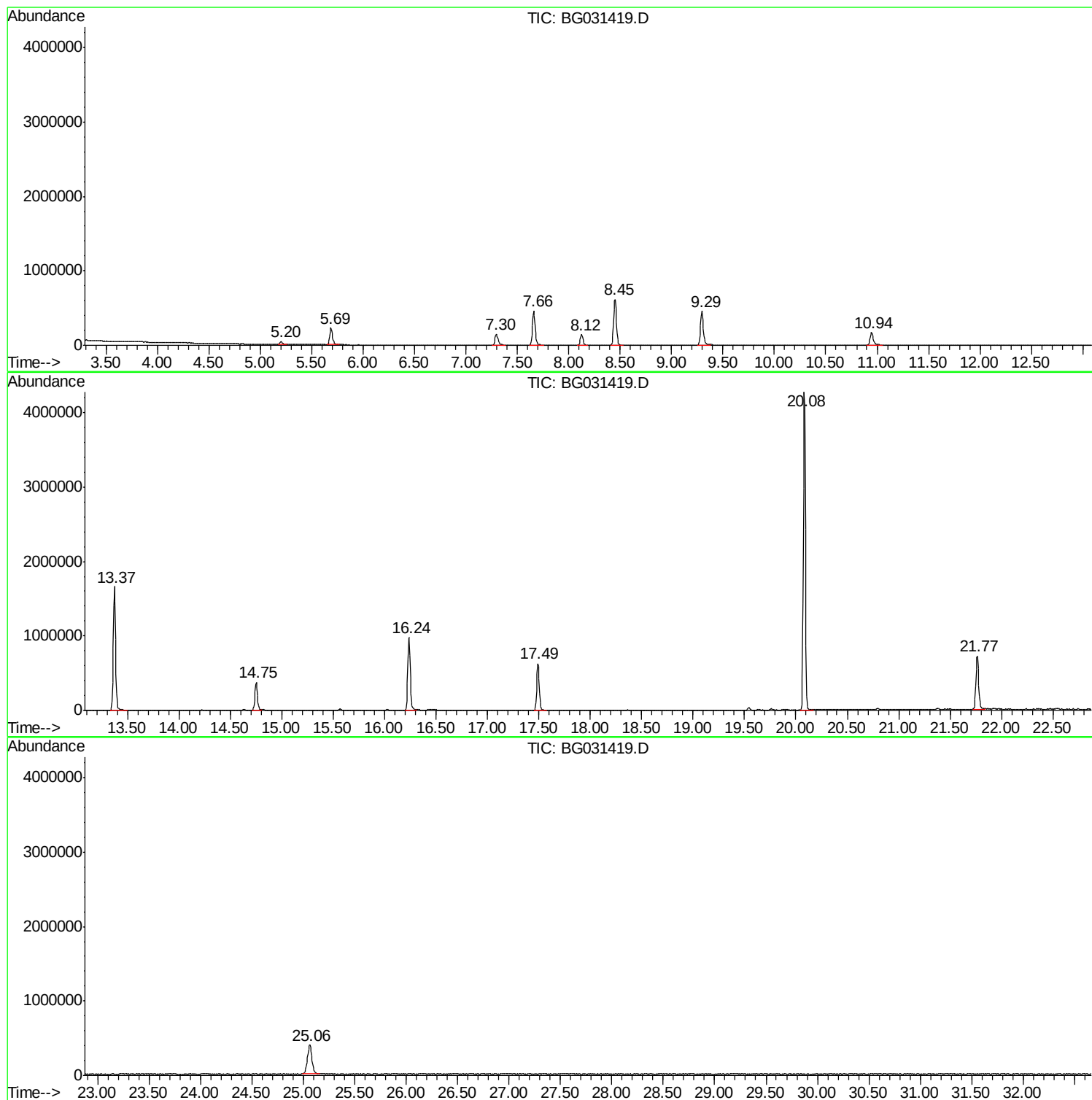
Sum of corrected areas: 18978497

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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG111017.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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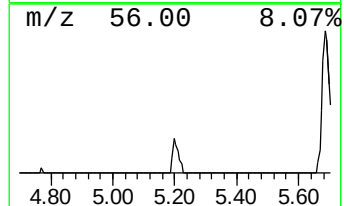
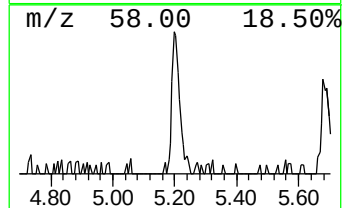
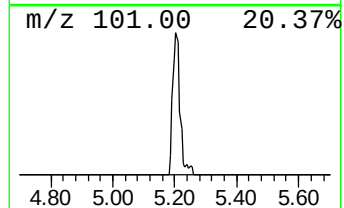
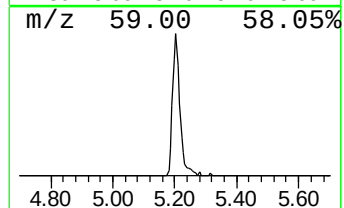
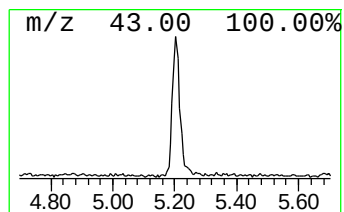
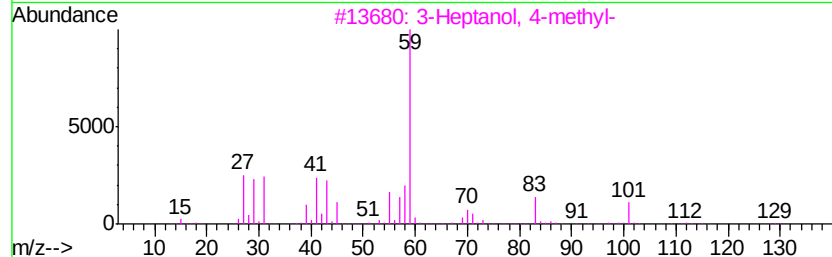
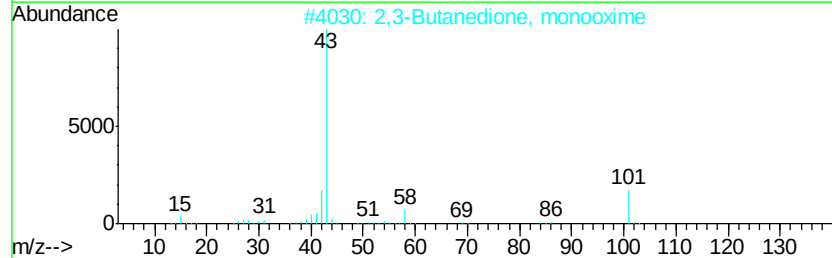
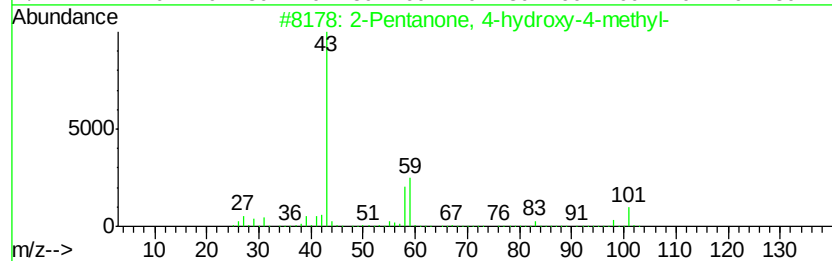
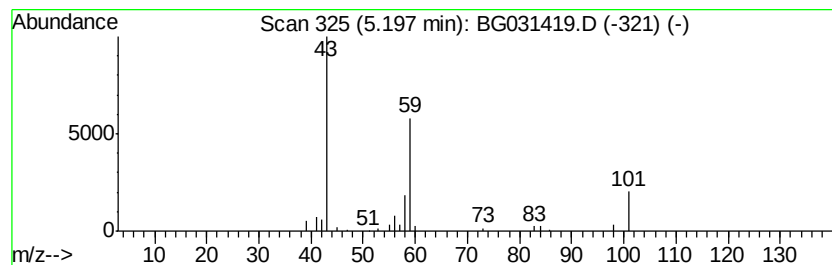
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	4.84 ng	63155	1,4-Dichlorobenzene-d4	8.12

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	32
3			3-Heptanol, 4-methyl-	130	C8H18O	014979-39-6	25
4			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	23
5			1,2-Butanediol	90	C4H10O2	000584-03-2	23



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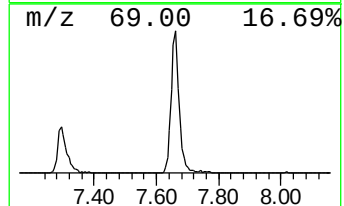
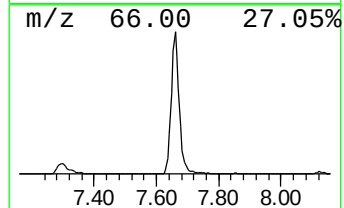
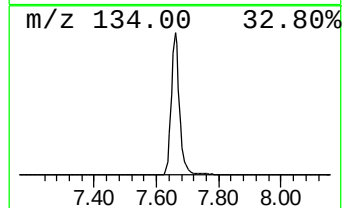
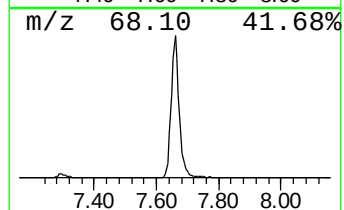
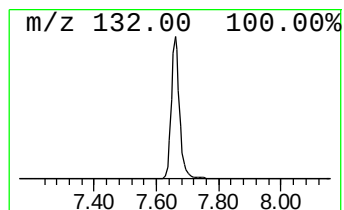
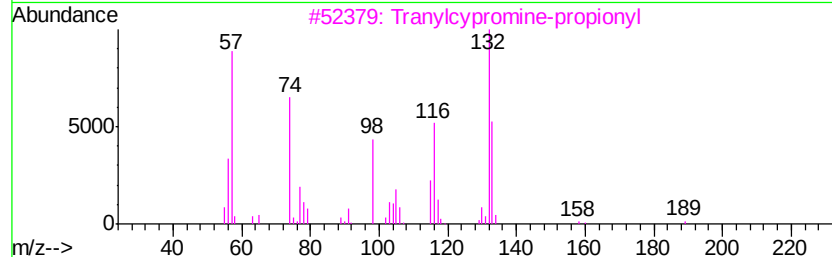
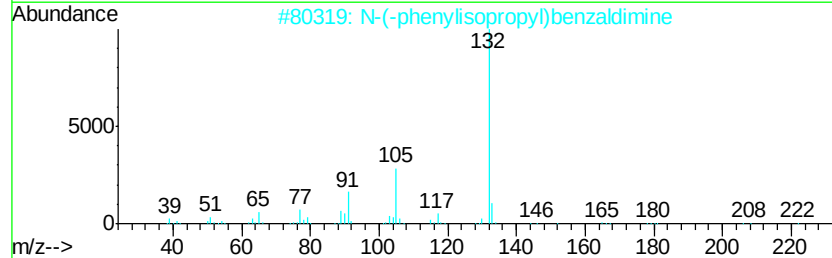
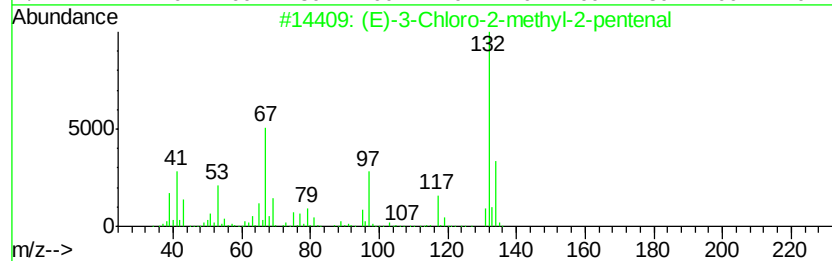
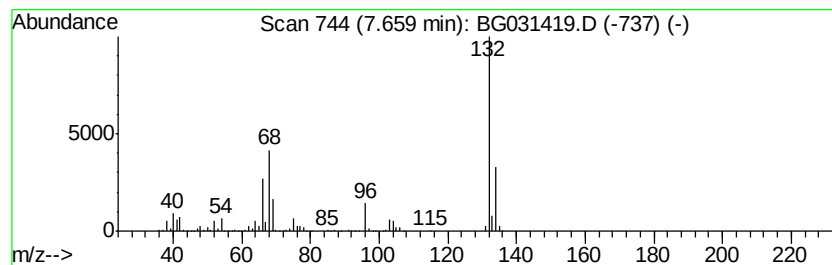
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Peak Number 2 unknown7.66 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.66	64.13 ng	837497	1,4-Dichlorobenzene-d4	8.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	27
2		N-(-phenylisopropyl)benzalimine	223	C16H17N	1000379-01-3	10
3		Tranlycypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		5-Aminoindole	132	C8H8N2	005192-03-0	10
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.20	4.8	ng	63155	1	8.12	261171	20.0
unknown7.66	7.66	64.1	ng	837497	1	8.12	261171	20.0