

Data Path : Z:\HPCHEM1\BNA G\DATA\BG112117\
 Data File : BG031176.D
 Acq On : 22 Nov 2017 6:36
 Operator : SJ/JU
 Sample : I6539-04
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 111517-TIDO-F-U-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	5.215	322	328	338	rVB	94693	151311	2.55%	0.739%
2	5.697	403	410	422	rBV	278595	498481	8.41%	2.434%
3	7.307	677	684	696	rBV	169429	344453	5.81%	1.682%
4	7.671	739	746	756	rBV	554417	975232	16.45%	4.761%
5	8.135	819	825	835	rVB	180910	315187	5.32%	1.539%
6	8.464	873	881	892	rBV	777878	1383662	23.33%	6.755%
7	9.304	1016	1024	1038	rBV	592247	1114767	18.80%	5.443%
8	10.955	1296	1305	1320	rBV	229985	471628	7.95%	2.303%
9	13.382	1710	1718	1731	rBV	1937219	3180839	53.64%	15.530%
10	14.757	1944	1952	1971	rVB2	431187	765399	12.91%	3.737%
11	16.243	2195	2205	2224	rBV	1012347	1668003	28.13%	8.144%
12	17.495	2410	2418	2433	rBV2	614107	1029294	17.36%	5.025%
13	20.092	2853	2860	2869	rBV	4168013	5930132	100.00%	28.953%
14	21.766	3138	3145	3155	rBV2	783782	1369568	23.10%	6.687%
15	25.068	3696	3707	3724	rVB2	428883	1284112	21.65%	6.269%

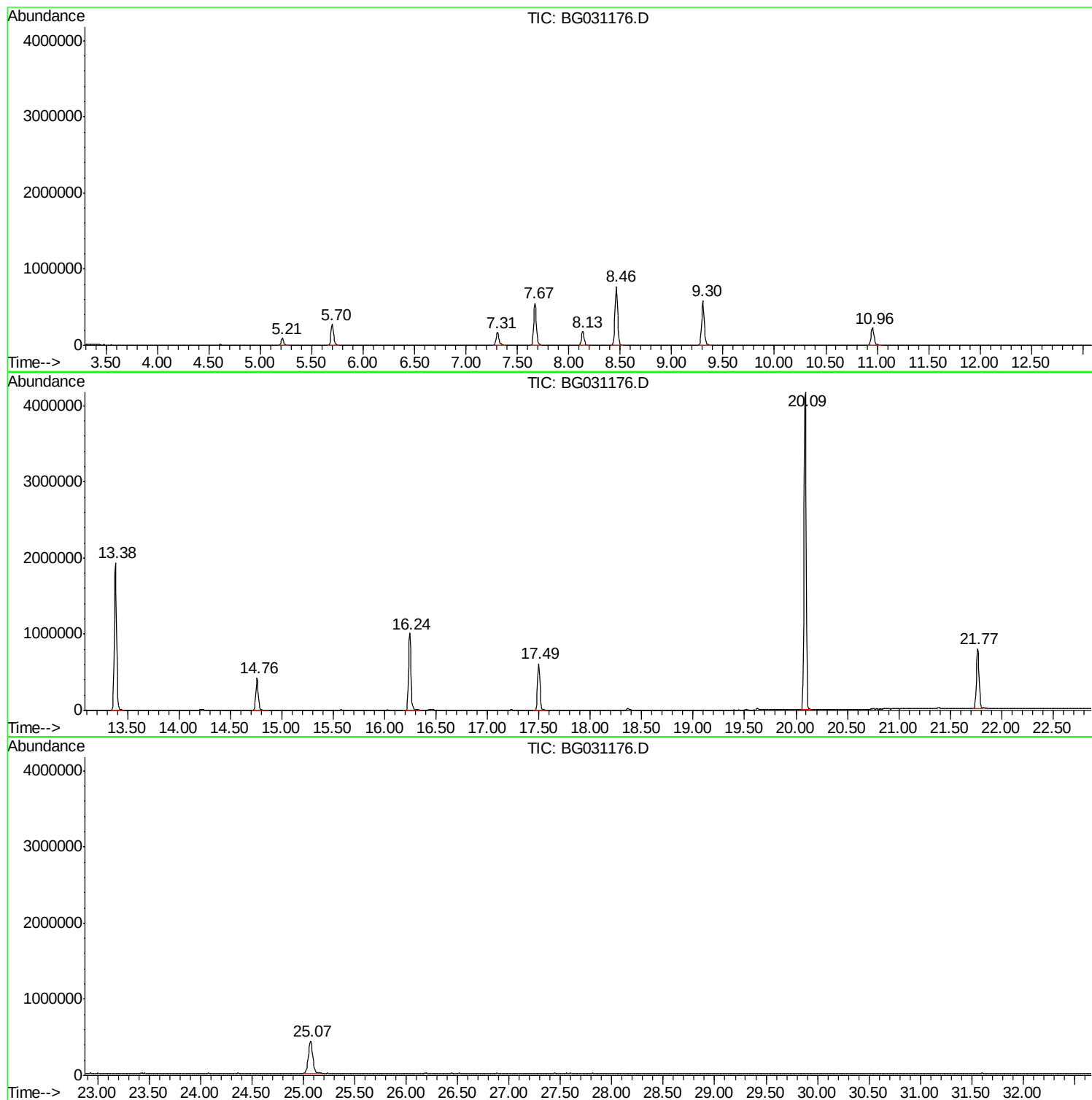
Sum of corrected areas: 20482068

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111517-TIDO-F-U-S

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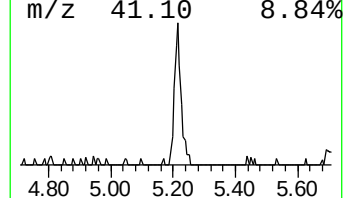
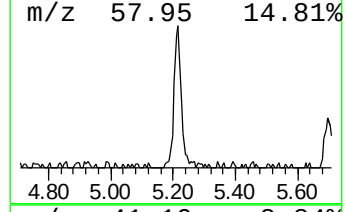
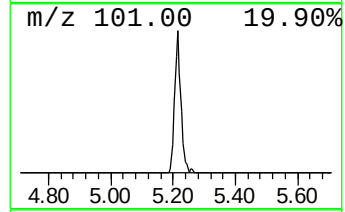
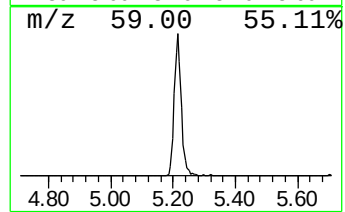
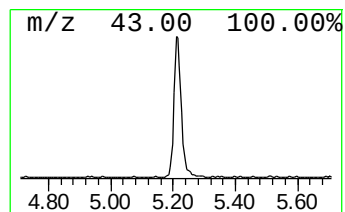
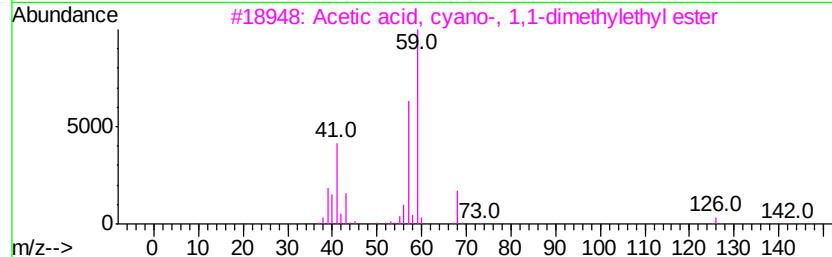
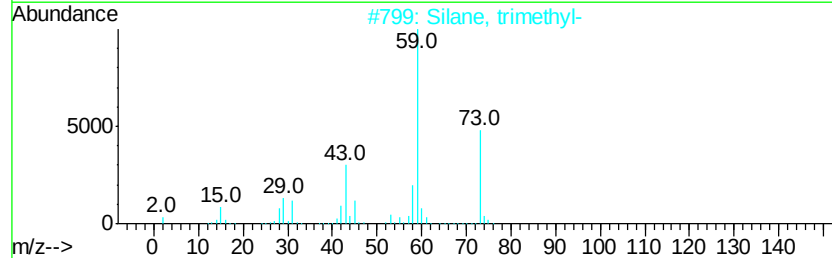
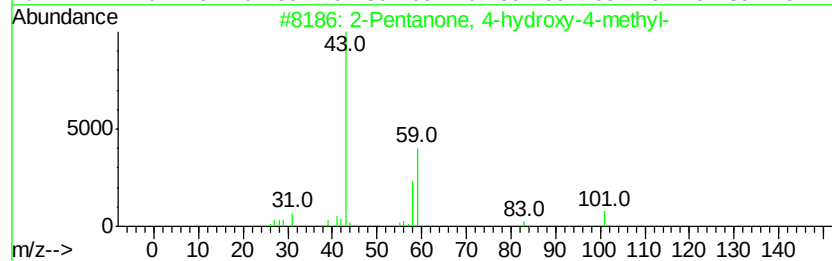
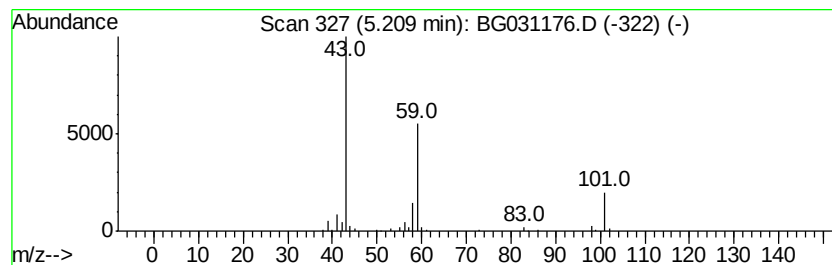
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.21	9.60 ng	151311	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			Silane, trimethyl-	74	C3H10Si	000993-07-7	23
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17
5			Acetone	58	C3H6O	000067-64-1	10



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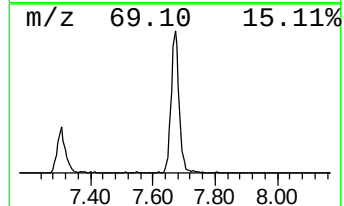
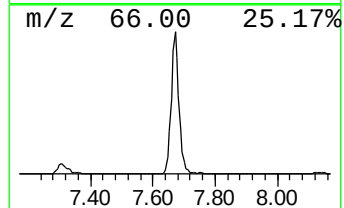
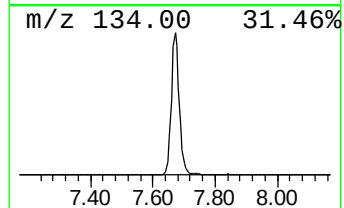
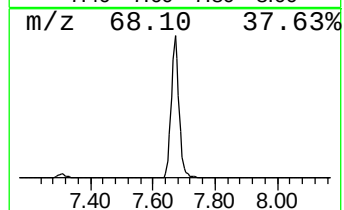
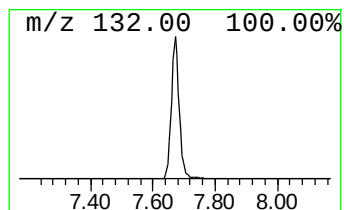
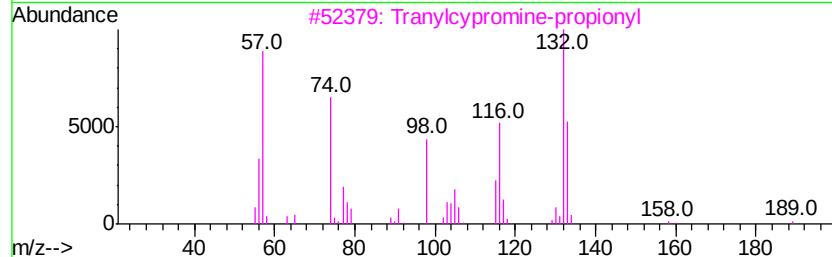
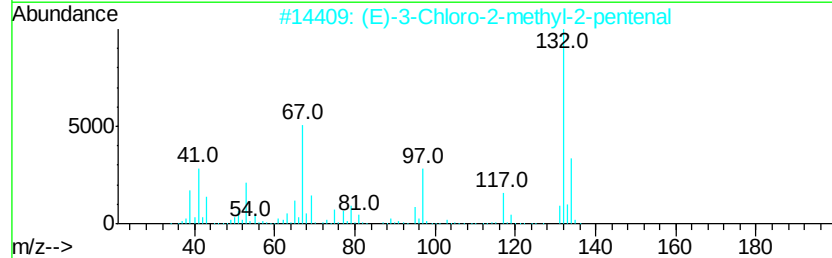
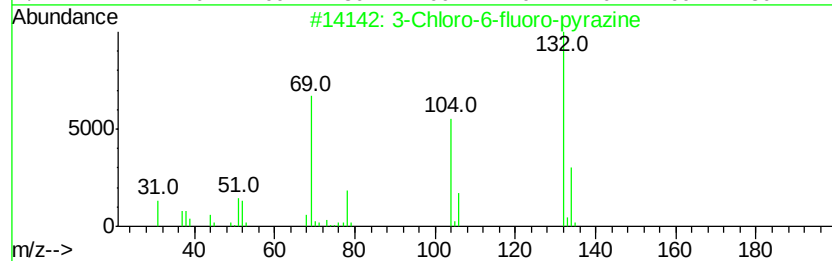
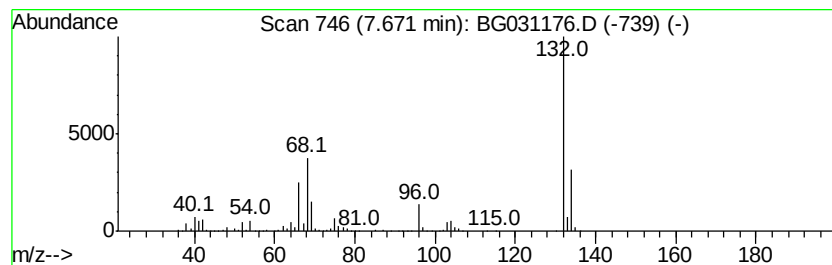
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Peak Number 2 unknown7.67 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.67	61.88 ng	975232	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-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	9.6	ng	151311	1	8.13	315187	20.0
unknown7.67	7.67	61.9	ng	975232	1	8.13	315187	20.0