

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG090718\
 Data File : BG036596.D
 Acq On : 7 Sep 2018 14:25
 Operator : JU/SJ
 Sample : PB112677BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB112677BL

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG082318.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	5.156	312	318	325	rBV	41364	62855	1.41%	0.287%
2	5.626	390	398	410	rBV	999315	1599279	35.93%	7.310%
3	7.212	659	668	683	rBV	1082680	1820946	40.91%	8.323%
4	7.588	723	732	743	rBV	966631	1668818	37.49%	7.627%
5	8.058	804	812	820	rVB	194526	333985	7.50%	1.526%
6	8.376	859	866	875	rBV	685633	1205205	27.08%	5.508%
7	9.216	1000	1009	1021	rBV	618845	1094718	24.59%	5.003%
8	10.861	1282	1289	1297	rBV	291781	513030	11.53%	2.345%
9	13.305	1697	1705	1719	rBV	1763430	2798362	62.87%	12.790%
10	14.680	1932	1939	1947	rBV2	515828	813403	18.27%	3.718%
11	16.166	2185	2192	2210	rBV	1237582	2030463	45.62%	9.280%
12	17.424	2399	2406	2417	rBV2	685336	1070110	24.04%	4.891%
13	19.721	2793	2797	2803	rBV	58592	64994	1.46%	0.297%
14	20.027	2843	2849	2860	rBV	3354785	4451182	100.00%	20.344%
15	20.761	2969	2974	2979	rBV	51201	61987	1.39%	0.283%
16	21.683	3125	3131	3141	rBV2	691221	1209876	27.18%	5.530%
17	24.897	3668	3678	3701	rBV2	304956	1080114	24.27%	4.937%

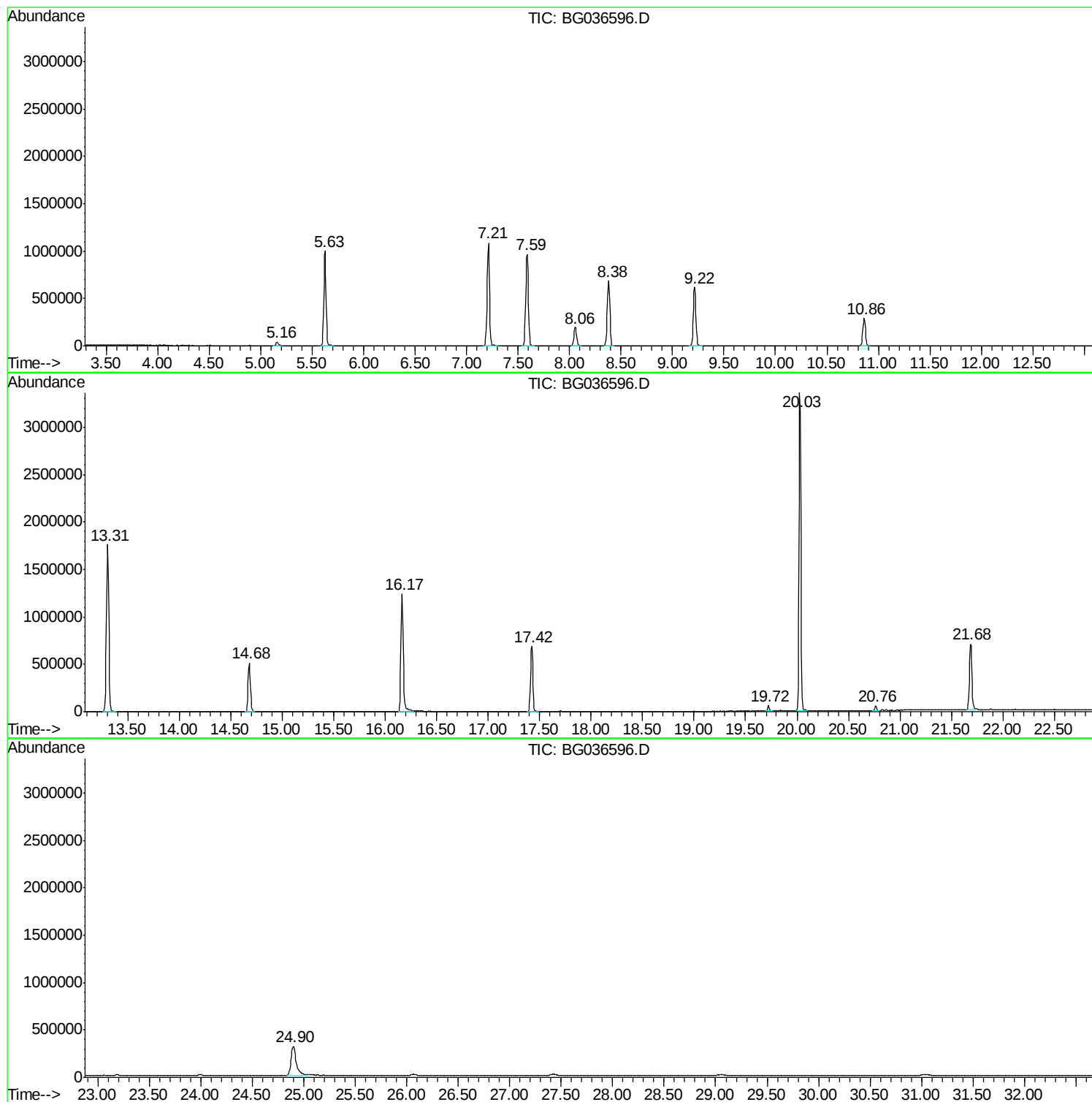
Sum of corrected areas: 21879327

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082318.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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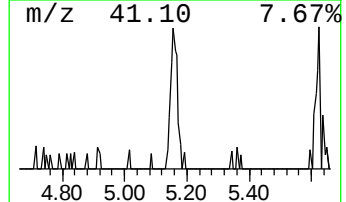
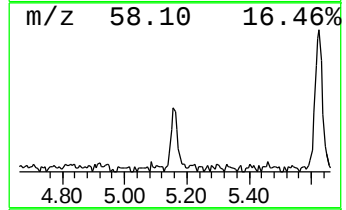
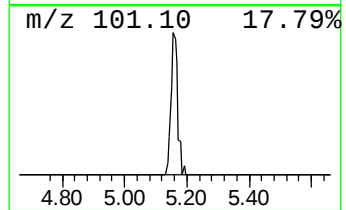
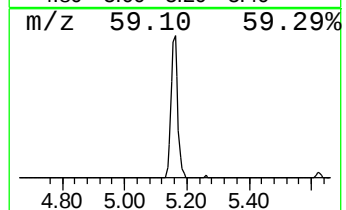
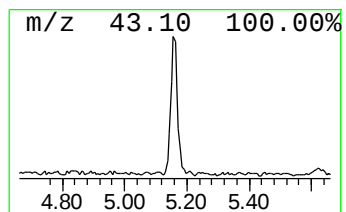
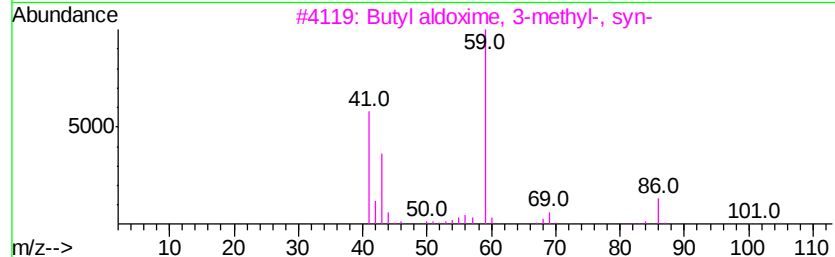
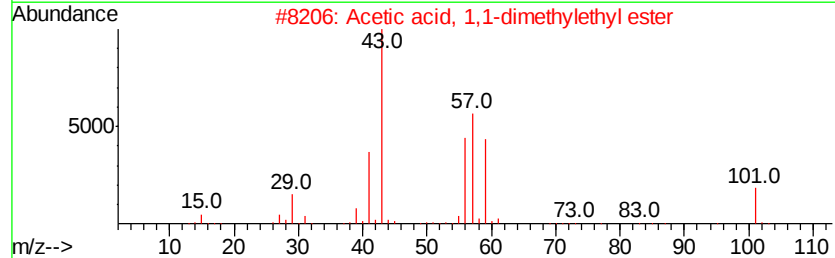
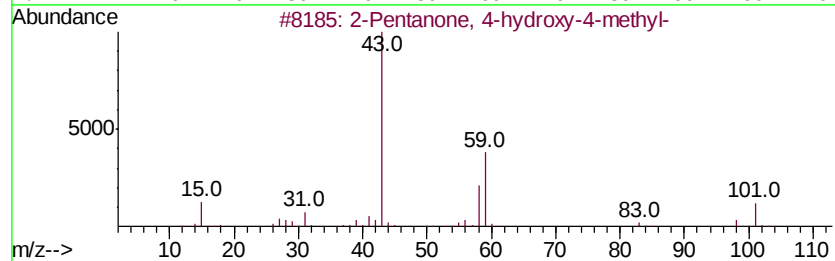
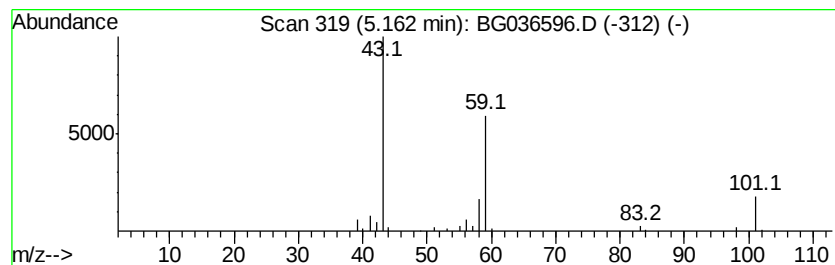
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	3.76 ng	62855	1,4-Dichlorobenzene-d4	8.06

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	23
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	10
5			5-Hexen-2-one	98	C6H10O	000109-49-9	10



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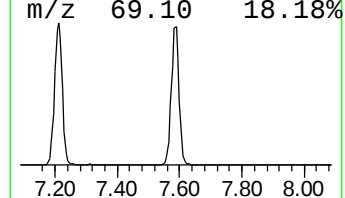
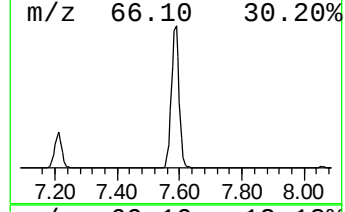
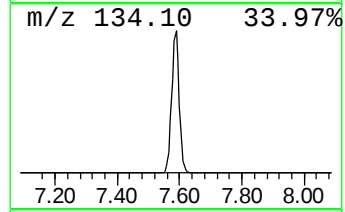
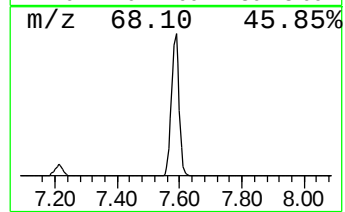
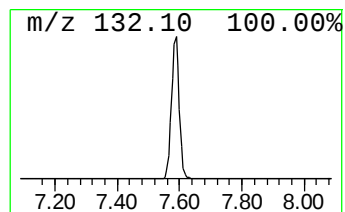
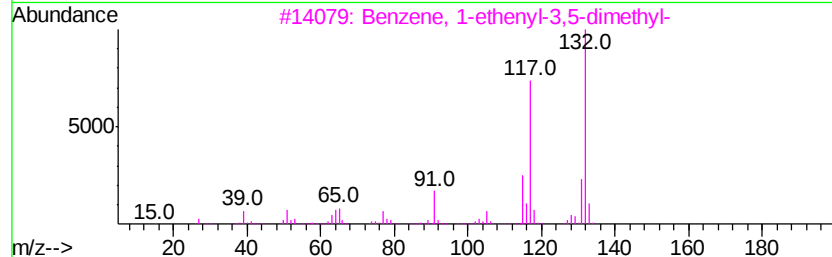
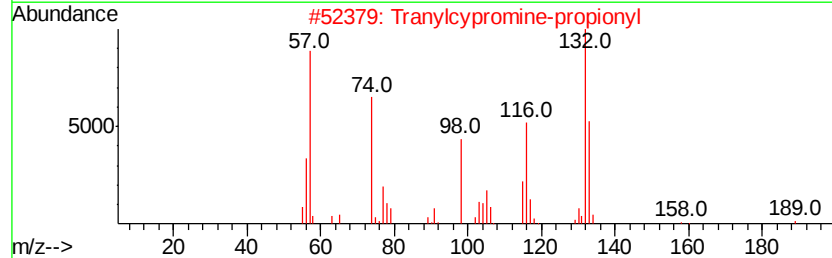
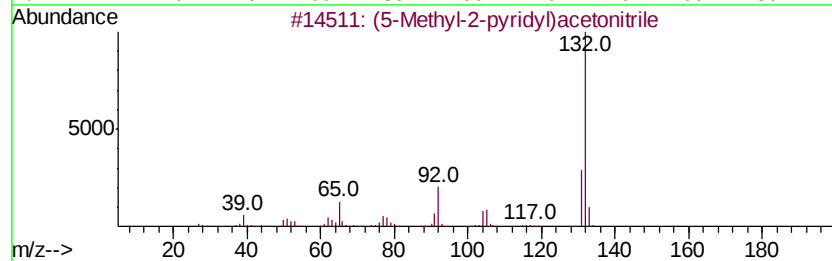
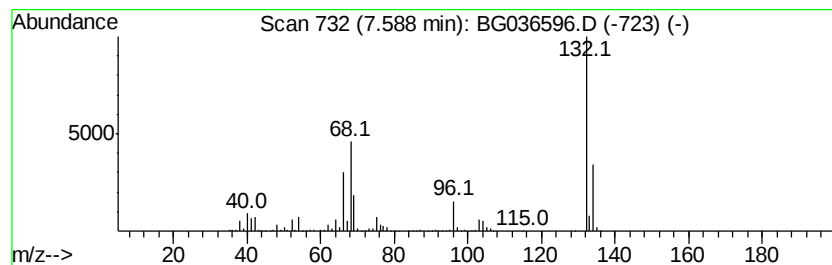
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Peak Number 2 unknown7.59 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	99.93 ng	1668820	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
2		Tranlycypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.16	3.8	ng	62855	1	8.06	333985	20.0
unknown7.59	7.59	99.9	ng	1668820	1	8.06	333985	20.0