

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG092618\
 Data File : BG036995.D
 Acq On : 26 Sep 2018 23:43
 Operator : JU/SJ
 Sample : PB113238BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB113238BL

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:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG092018.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	5.146	311	316	325	rBV	30705	52723	1.58%	0.333%
2	5.616	390	396	415	rBV	692474	1261288	37.84%	7.959%
3	7.203	659	666	692	rBV	694589	1401946	42.06%	8.847%
4	7.573	722	729	746	rBV	675687	1296511	38.90%	8.181%
5	8.043	803	809	821	rBV	123689	240493	7.22%	1.518%
6	8.360	856	863	880	rBV	511539	1011701	30.35%	6.384%
7	9.200	999	1006	1029	rBV	394868	868352	26.05%	5.480%
8	10.846	1279	1286	1306	rBV	158338	337825	10.14%	2.132%
9	13.290	1695	1702	1718	rBV	1145348	1959078	58.78%	12.362%
10	14.665	1929	1936	1951	rBV2	269597	481033	14.43%	3.035%
11	16.151	2183	2189	2217	rBV	774607	1445258	43.36%	9.120%
12	17.414	2397	2404	2418	rBV	370482	648103	19.45%	4.090%
13	20.017	2841	2847	2868	rBV	2358618	3332993	100.00%	21.032%
14	21.680	3123	3130	3147	rBV	427067	777172	23.32%	4.904%
15	24.882	3664	3675	3691	rBV2	227874	732541	21.98%	4.623%

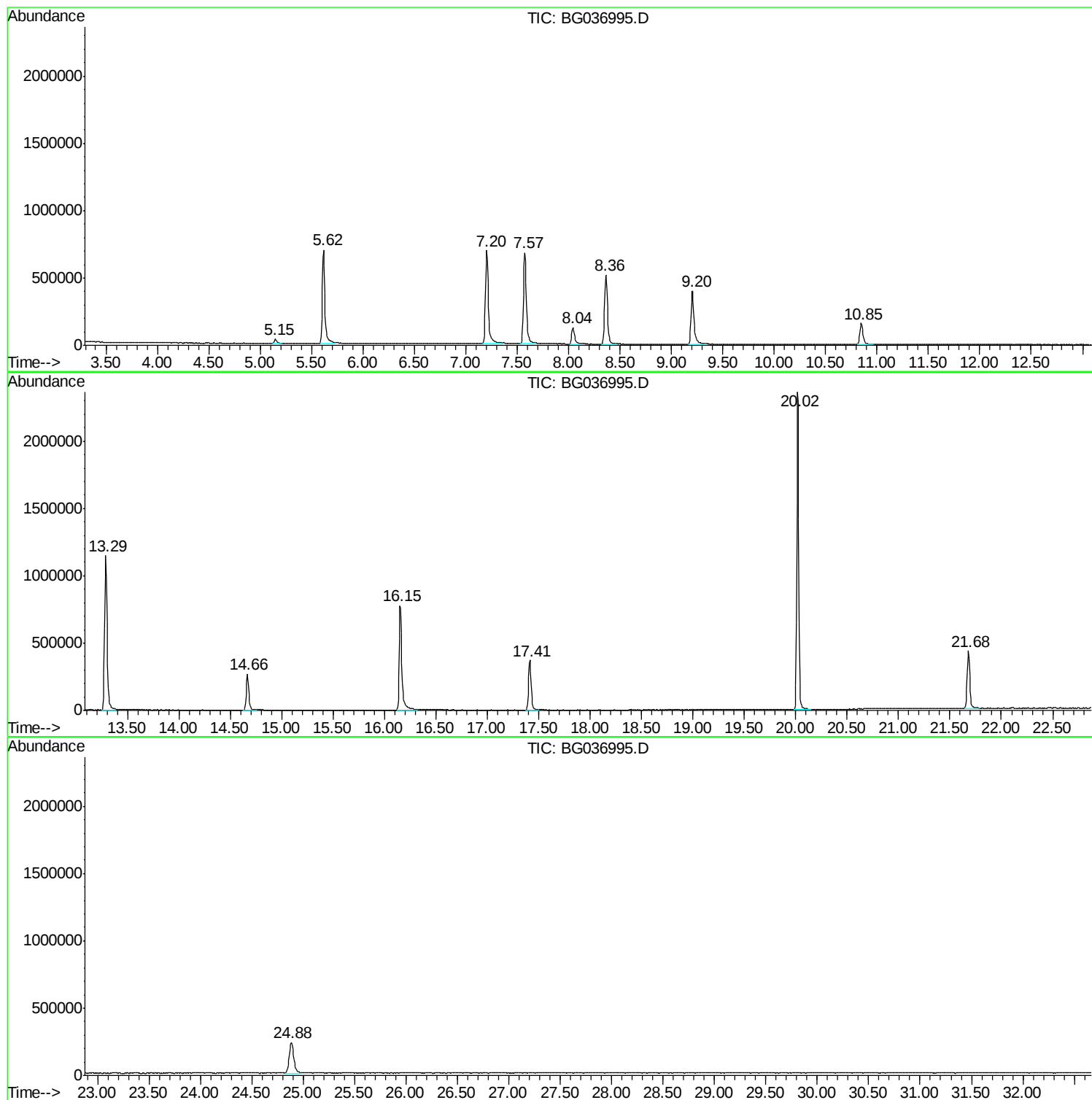
Sum of corrected areas: 15847017

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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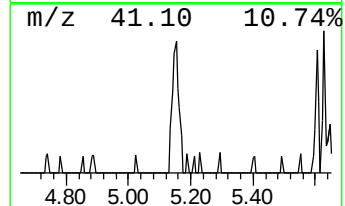
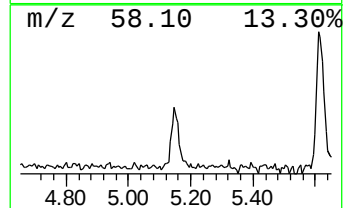
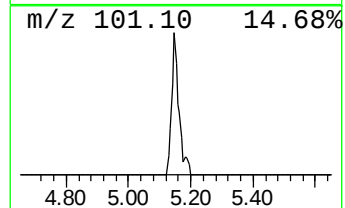
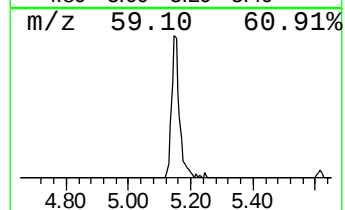
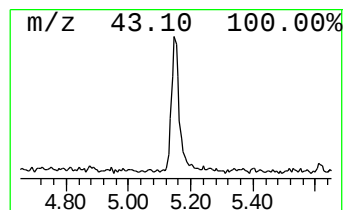
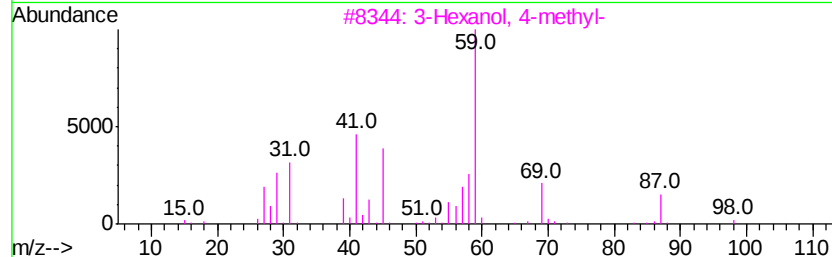
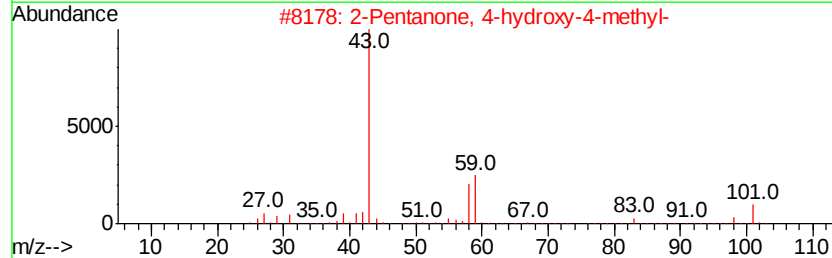
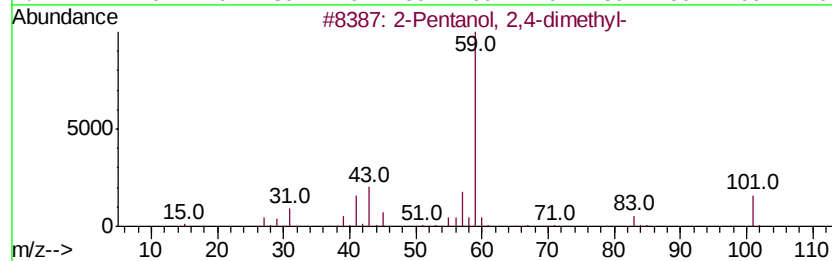
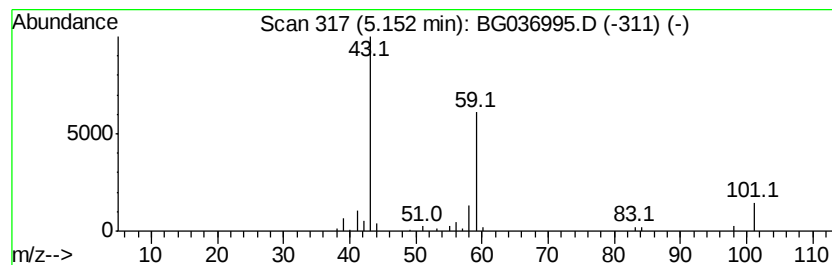
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Peak Number 1 2-Pentanol, 2,4-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.15	4.38 ng	52723	1,4-Dichlorobenzene-d4	8.04

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	53
2			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	38
4			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
5			3-Heptanol	116	C7H16O	000589-82-2	25



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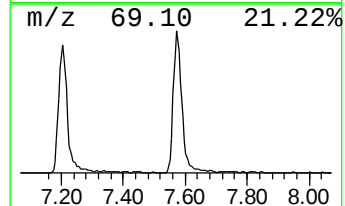
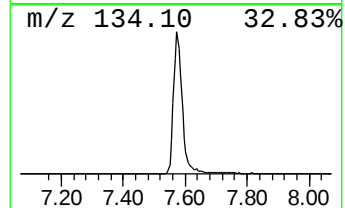
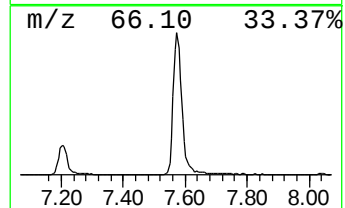
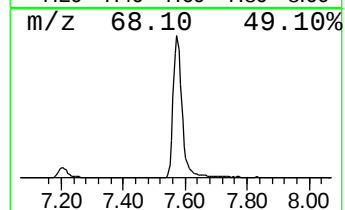
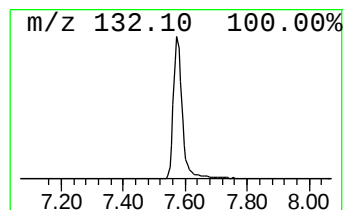
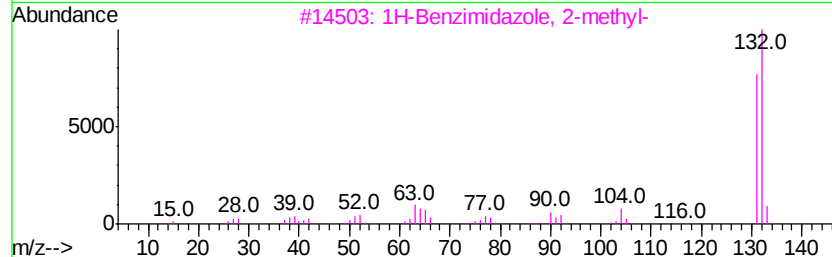
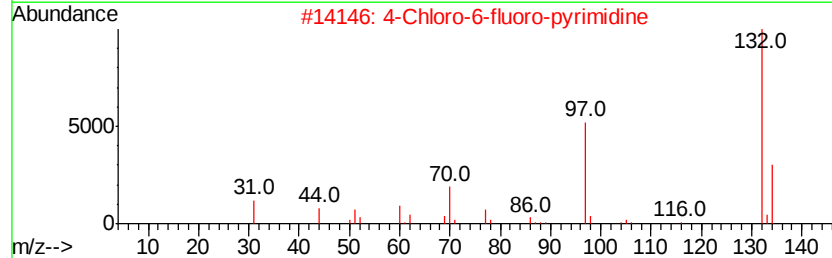
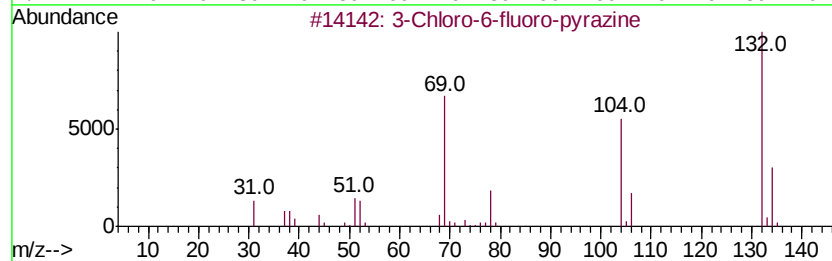
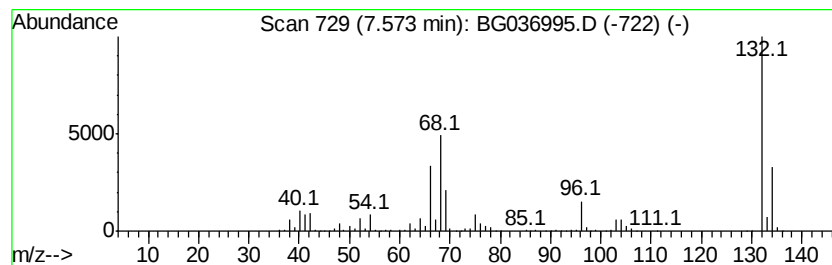
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Peak Number 2 unknown7.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.57	107.82 ng	1296510	1,4-Dichlorobenzene-d4	8.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	22
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanol, 2,4-d...	5.15	4.4	ng	52723	1	8.04	240493	20.0
unknown7.57	7.57	107.8	ng	1296510	1	8.04	240493	20.0