

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG092218\
 Data File : BG036911.D
 Acq On : 23 Sep 2018 3:18
 Operator : JU/SJ
 Sample : J5019-18
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 091918-TIPC-F-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:\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	5.153	311	317	330	rVB	24651	42755	1.04%	0.291%
2	5.617	389	396	415	rBV	182053	358291	8.75%	2.441%
3	7.203	660	666	687	rBV	108657	265657	6.49%	1.810%
4	7.579	722	730	749	rBV	361054	722800	17.65%	4.925%
5	8.043	803	809	825	rBV	106451	215823	5.27%	1.471%
6	8.366	855	864	883	rBV	547574	1045280	25.53%	7.122%
7	9.207	999	1007	1024	rBV	456085	950459	23.21%	6.476%
8	10.852	1279	1287	1299	rBV	154040	314629	7.68%	2.144%
9	13.296	1695	1703	1723	rBV	1348895	2319093	56.64%	15.802%
10	14.118	1837	1843	1856	rVB	91760	147566	3.60%	1.005%
11	14.671	1930	1937	1952	rBV3	282260	500643	12.23%	3.411%
12	16.157	2183	2190	2208	rBV	739693	1223919	29.89%	8.339%
13	17.415	2397	2404	2417	rBV	401190	669069	16.34%	4.559%
14	20.023	2842	2848	2859	rBV	2966668	4094465	100.00%	27.898%
15	21.322	3064	3069	3073	rBV3	33715	58433	1.43%	0.398%
16	21.680	3123	3130	3140	rBV2	464755	798839	19.51%	5.443%
17	23.161	3377	3382	3390	rVB3	25700	51968	1.27%	0.354%
18	23.966	3512	3519	3526	rVB4	28054	62551	1.53%	0.426%
19	24.888	3666	3676	3688	rBV3	272254	834069	20.37%	5.683%

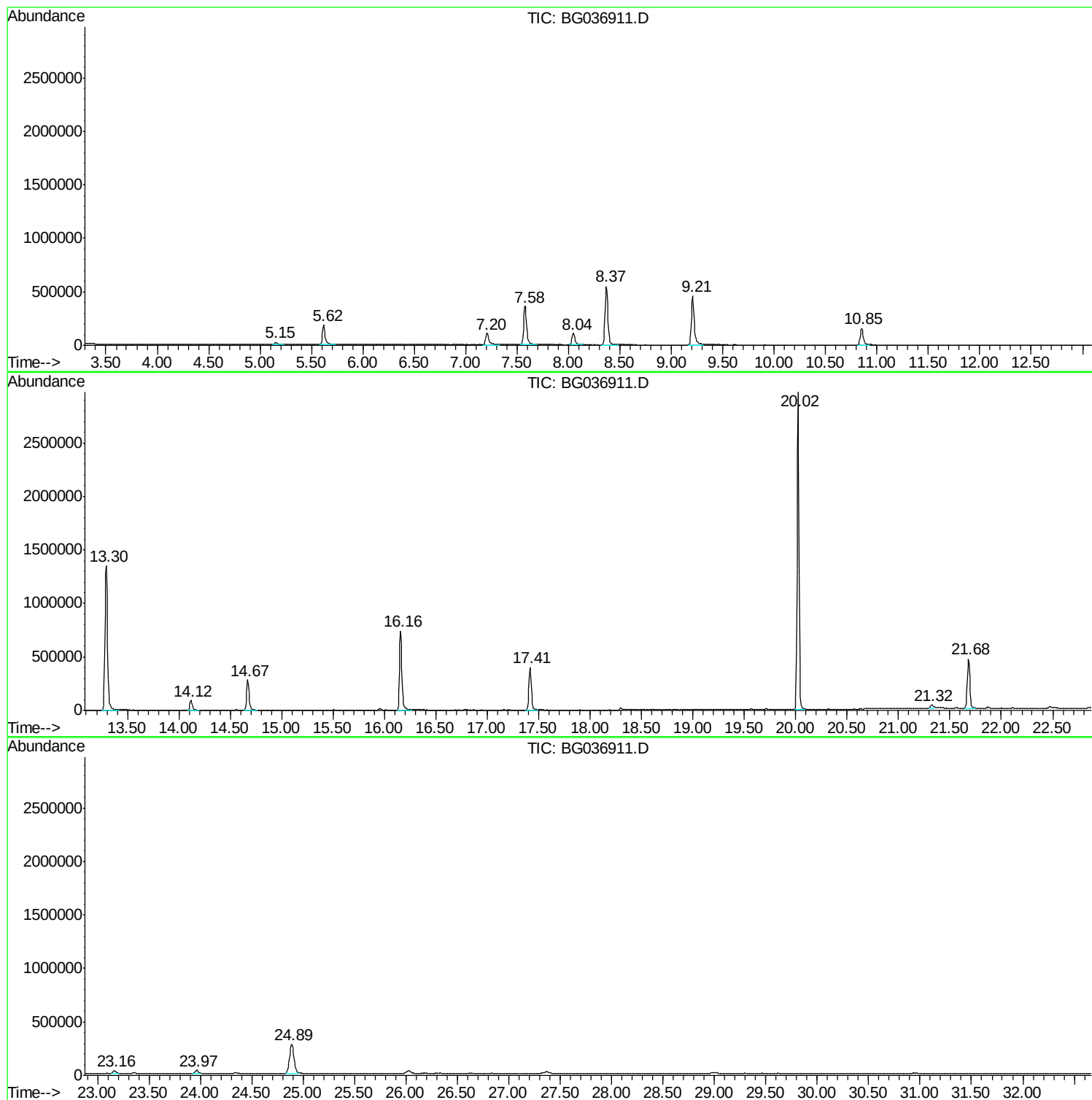
Sum of corrected areas: 14676309

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092018.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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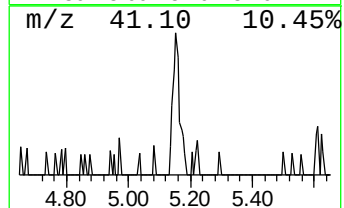
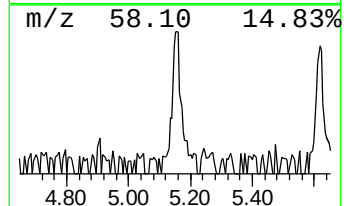
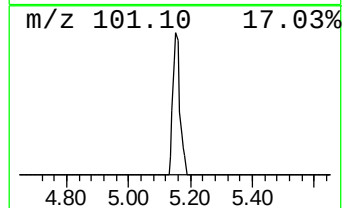
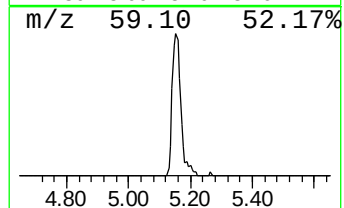
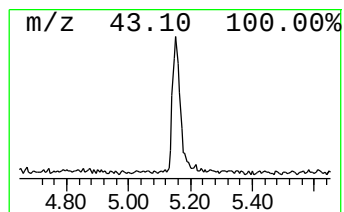
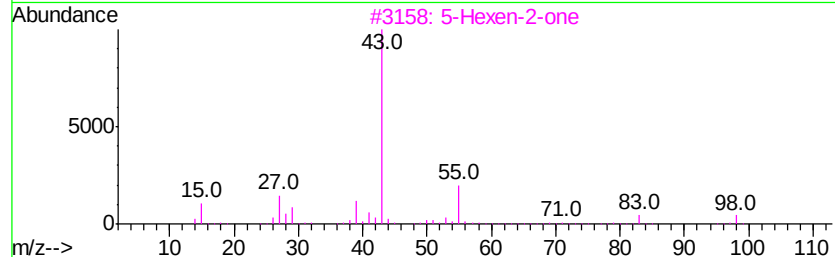
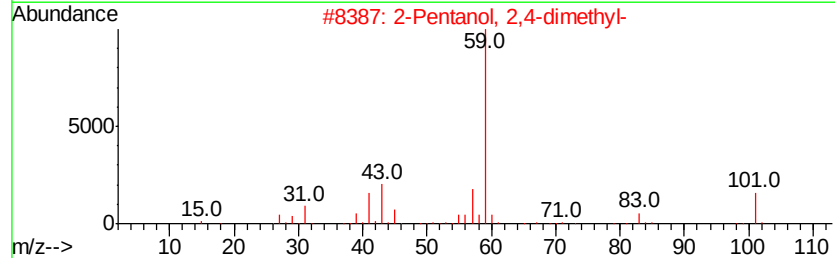
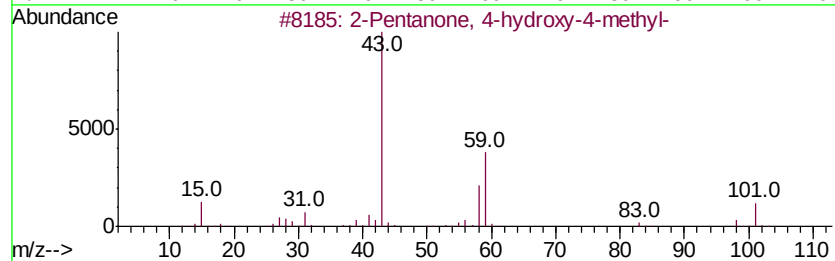
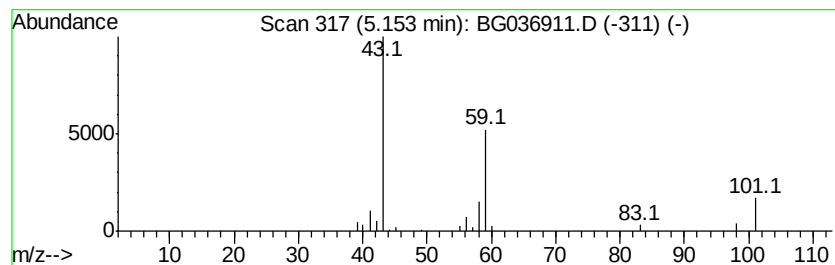
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.15	3.96 ng	42755	1,4-Dichlorobenzene-d4	8.04

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	28
3			5-Hexen-2-one	98	C6H10O	000109-49-9	9
4			Diacetamide	101	C4H7NO2	000625-77-4	9
5			Ethanone, 1-(2-methylcyclopropyl)-	98	C6H10O	000930-56-3	9



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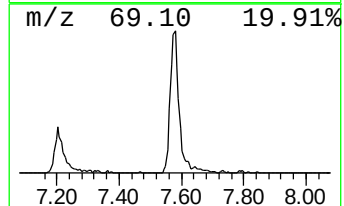
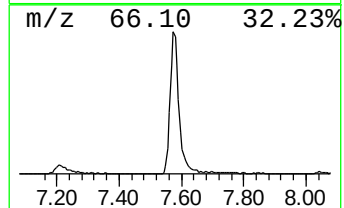
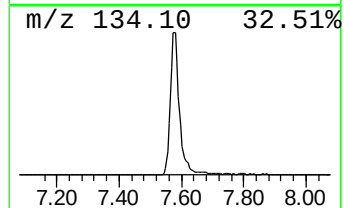
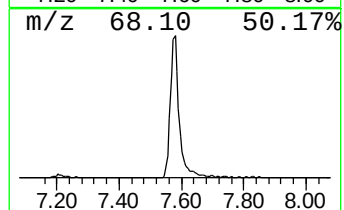
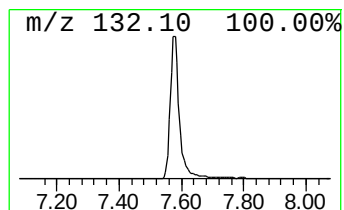
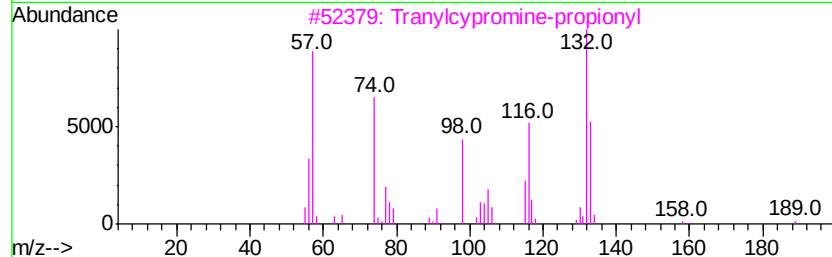
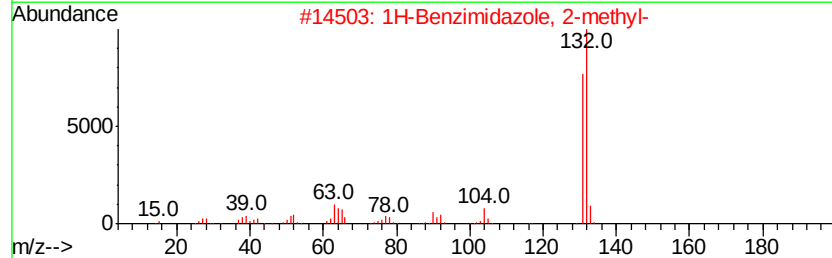
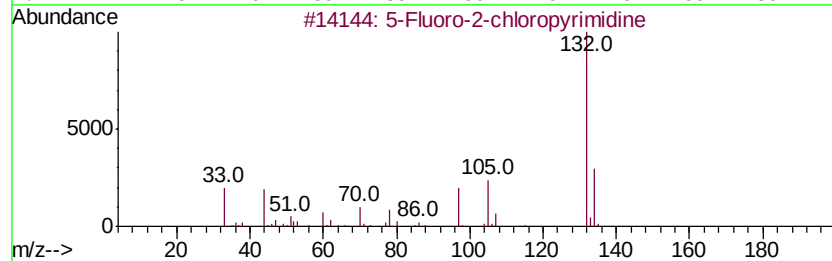
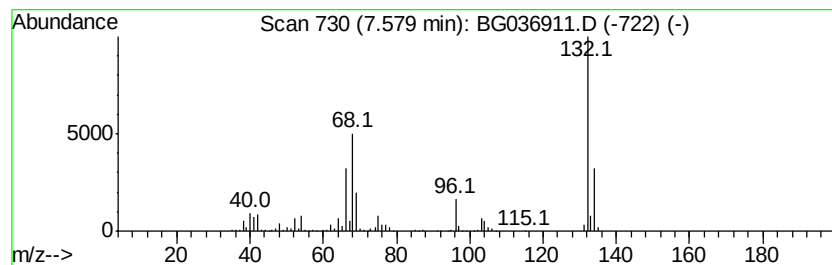
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Peak Number 2 unknown7.58 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	66.98 ng	722800	1,4-Dichlorobenzene-d4	8.04

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	27
2			1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3			Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4			(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
5			2-Cyclohexen-1-one	96	C6H8O	000930-68-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.15	4.0	ng	42755	1	8.04	215823	20.0
unknown7.58	7.58	67.0	ng	722800	1	8.04	215823	20.0