

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG100618\
 Data File : BG037193.D
 Acq On : 5 Oct 2018 17:05
 Operator : JU/SJ
 Sample : PB113546BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB113546BL

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	5.138	310	315	321	rBV	16364	27220	1.46%	0.279%
2	5.603	388	394	415	rBV	328701	638682	34.30%	6.546%
3	7.189	658	664	686	rBV	320799	702688	37.74%	7.202%
4	7.559	721	727	746	rBV	326414	665631	35.75%	6.822%
5	8.029	801	807	820	rBV	95895	205706	11.05%	2.108%
6	8.352	855	862	876	rBV	264964	515319	27.67%	5.282%
7	9.192	997	1005	1022	rBV	169242	420707	22.59%	4.312%
8	10.838	1278	1285	1299	rBV	126393	281883	15.14%	2.889%
9	13.276	1693	1700	1718	rBV	569931	1030138	55.32%	10.558%
10	14.657	1928	1935	1951	rBV2	251113	446164	23.96%	4.573%
11	16.143	2182	2188	2205	rBV	459912	837148	44.96%	8.580%
12	17.400	2395	2402	2422	rBV2	356488	633618	34.03%	6.494%
13	20.009	2840	2846	2860	rBV	1261318	1862144	100.00%	19.085%
14	21.666	3121	3128	3142	rBV	393119	753302	40.45%	7.721%
15	23.928	3508	3513	3518	rBV8	12310	25399	1.36%	0.260%
16	24.862	3661	3672	3688	rBV2	227344	711198	38.19%	7.289%

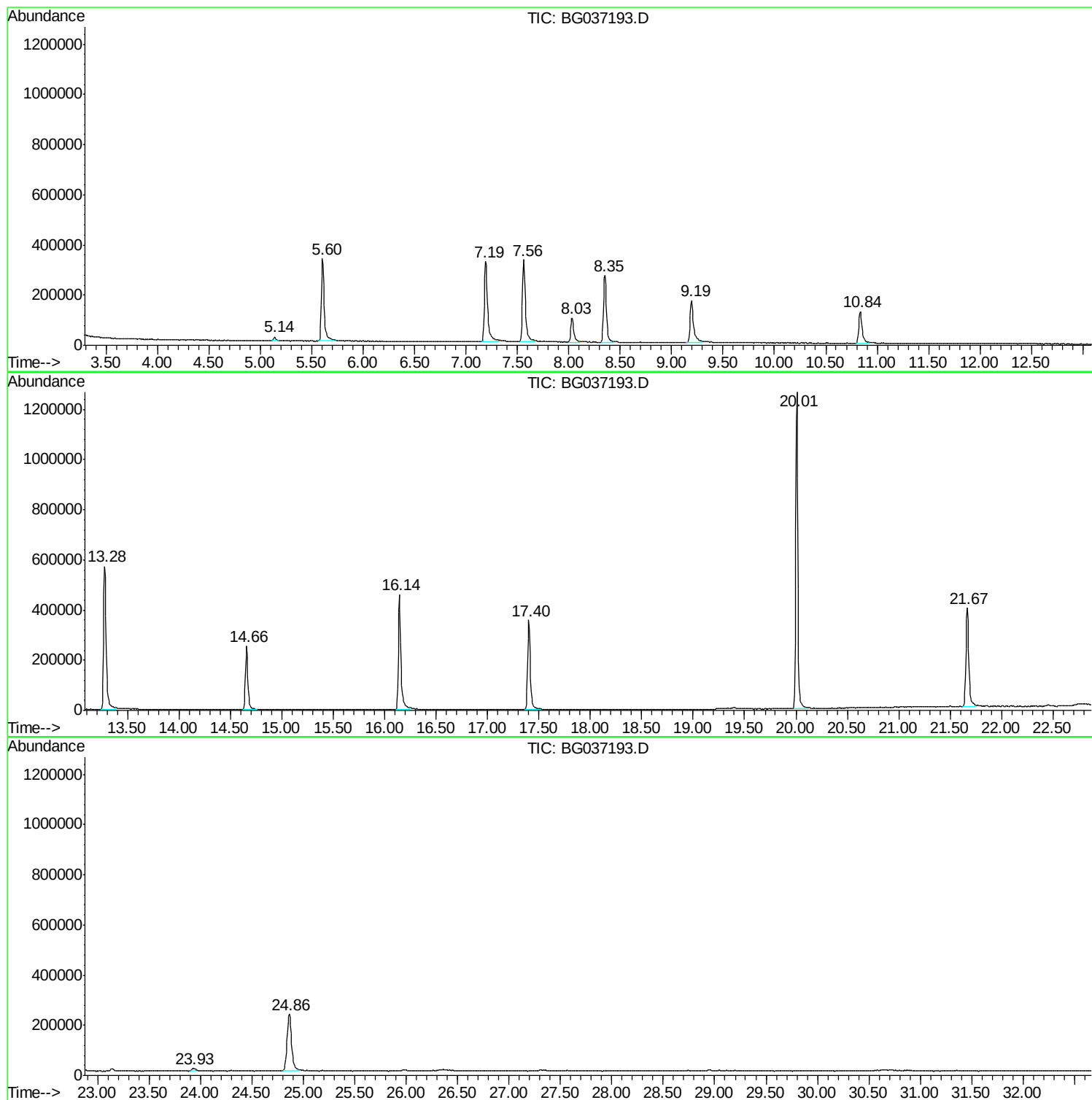
Sum of corrected areas: 9756947

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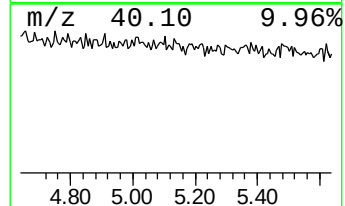
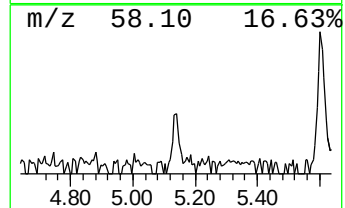
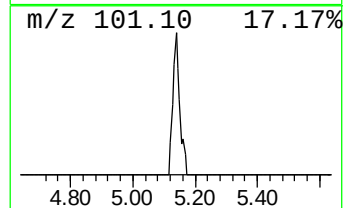
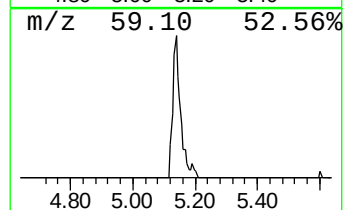
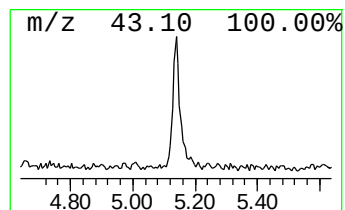
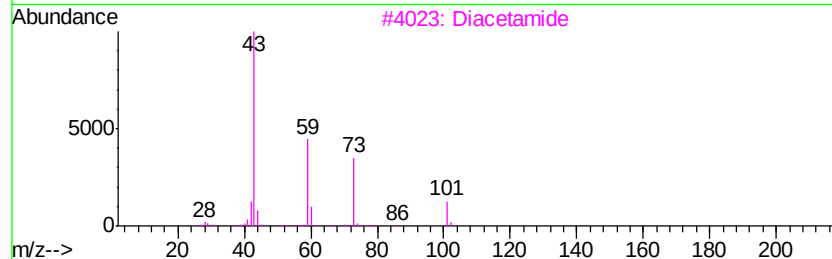
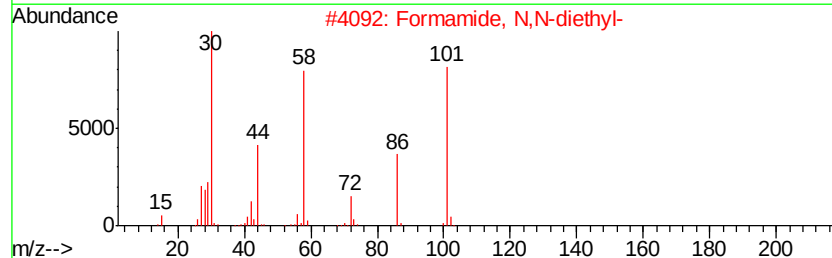
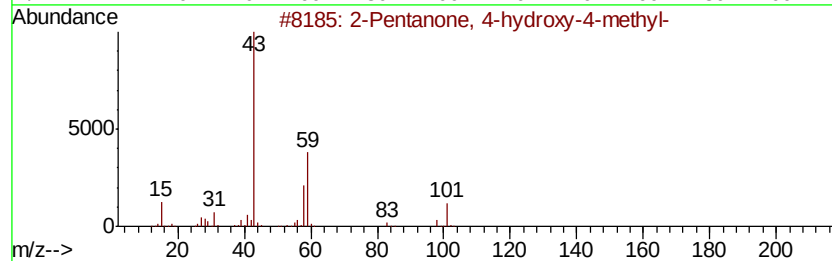
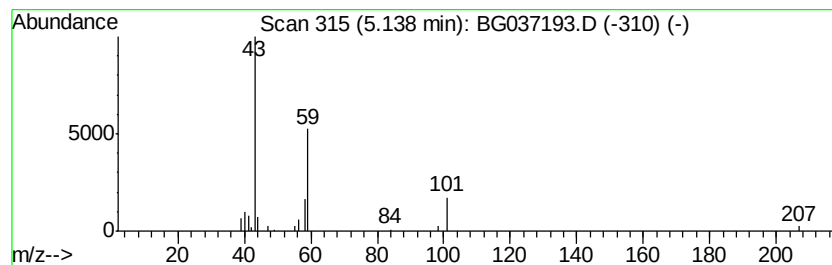
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	2.65 ng	27220	1,4-Dichlorobenzene-d4	8.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	33
2		Formamide, N,N-diethyl-	101	C5H11NO	000617-84-5	9
3		Diacetamide	101	C4H7NO2	000625-77-4	9
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		3-Octanol	130	C8H18O	000589-98-0	9



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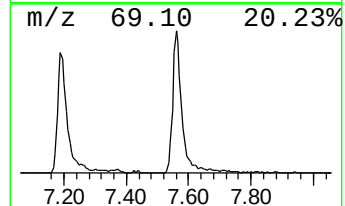
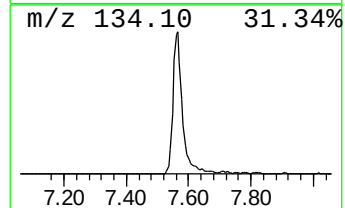
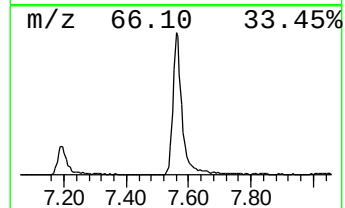
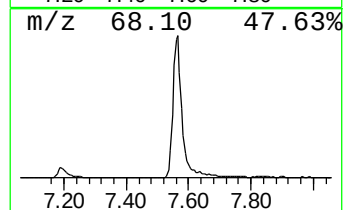
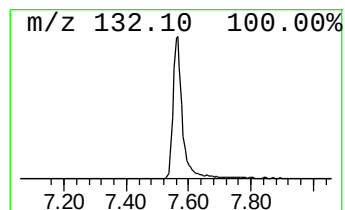
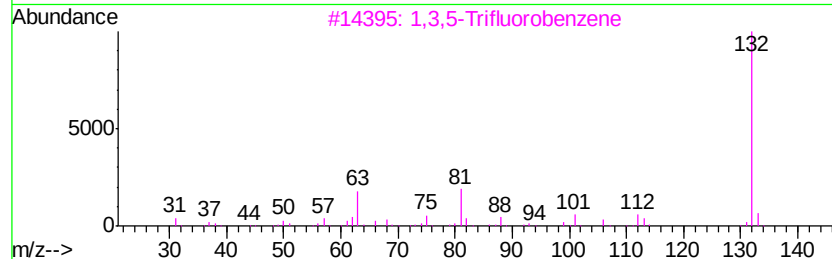
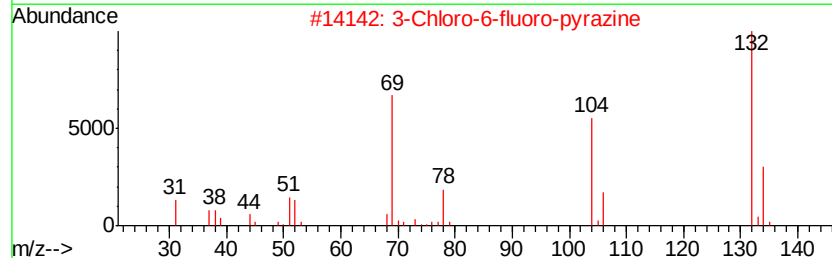
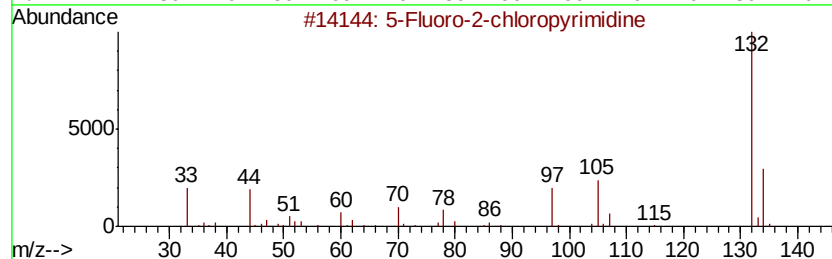
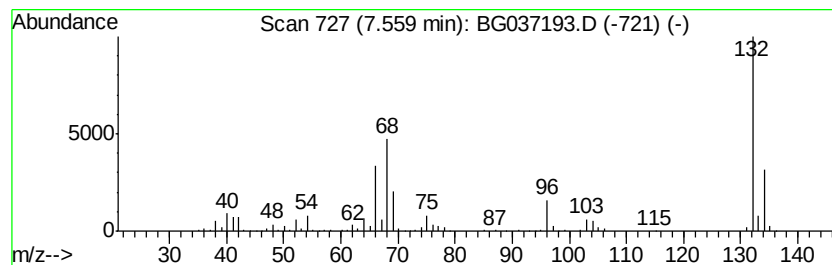
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Peak Number 2 unknown7.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	64.72 ng	665631	1,4-Dichlorobenzene-d4	8.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
3		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	10
4		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.14	2.6	ng	27220	1	8.03	205706	20.0
unknown7.56	7.56	64.7	ng	665631	1	8.03	205706	20.0