

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120118\
 Data File : BG038293.D
 Acq On : 1 Dec 2018 18:46
 Operator : JU/SJ
 Sample : PB115213BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB115213BL

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-BG111318.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	5.159	313	318	326	rVB	22153	36958	1.68%	0.369%
2	5.623	390	397	412	rVB	443693	742553	33.75%	7.404%
3	7.227	662	670	679	rBV	451511	817744	37.17%	8.154%
4	7.603	726	734	746	rBV	440447	775278	35.24%	7.731%
5	8.079	807	815	822	rBV	88210	157343	7.15%	1.569%
6	8.396	862	869	880	rVB	324471	592368	26.92%	5.907%
7	9.254	1007	1015	1025	rBV	269752	508117	23.09%	5.067%
8	10.899	1287	1295	1306	rBV	110493	219595	9.98%	2.190%
9	13.332	1702	1709	1724	rBV	737487	1183029	53.77%	11.797%
10	14.712	1937	1944	1954	rBV2	195302	324743	14.76%	3.238%
11	16.199	2191	2197	2211	rBV2	603523	957306	43.51%	9.546%
12	17.462	2405	2412	2421	rBV	289188	465031	21.14%	4.637%
13	20.059	2848	2854	2867	rBV	1680844	2200158	100.00%	21.939%
14	21.739	3133	3140	3151	rBV2	310447	536038	24.36%	5.345%
15	25.059	3687	3705	3717	rBV3	164286	512231	23.28%	5.108%

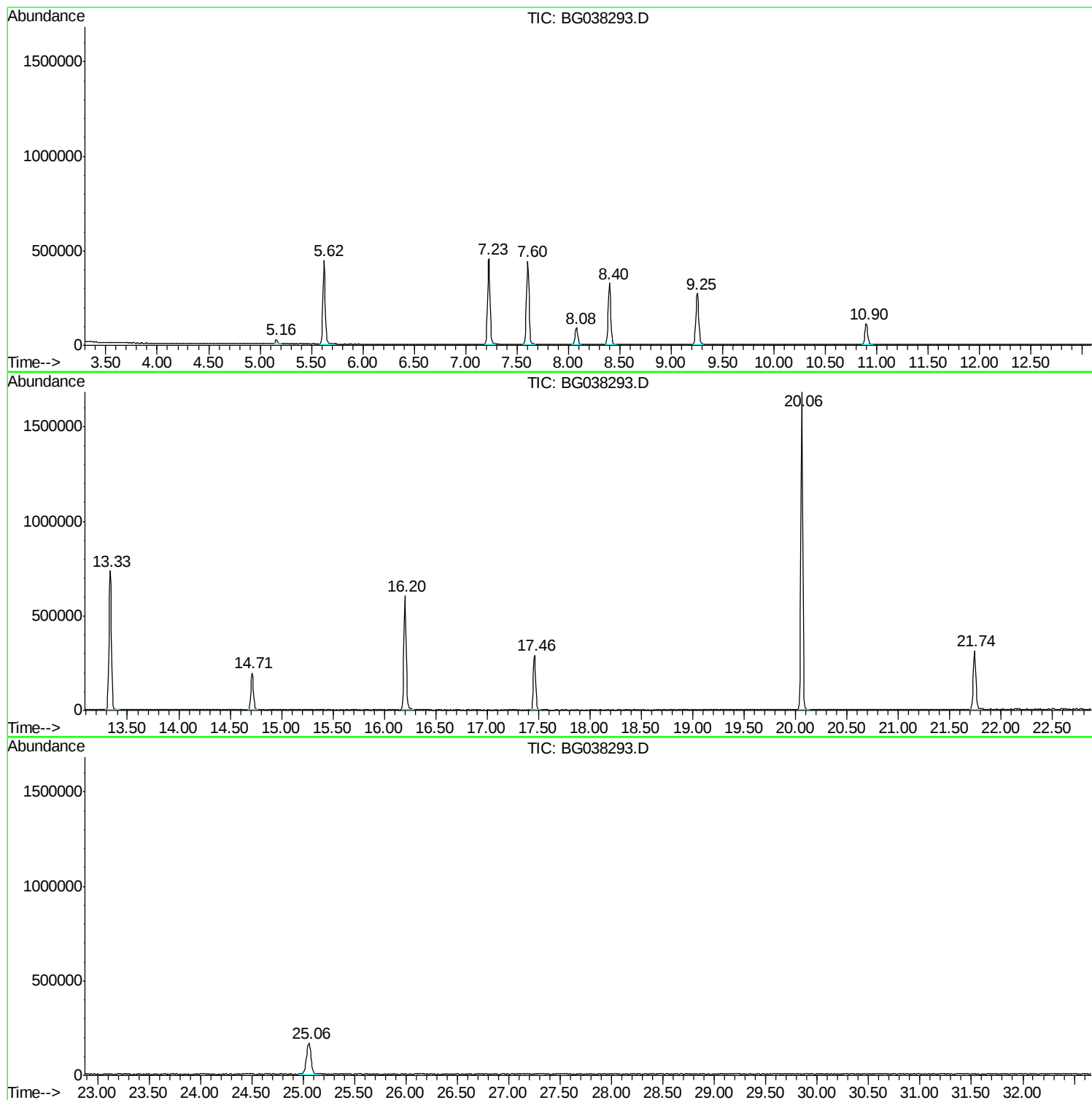
Sum of corrected areas: 10028492

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



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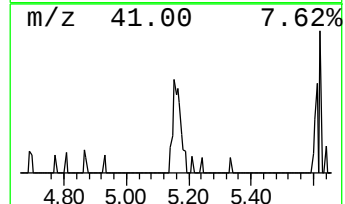
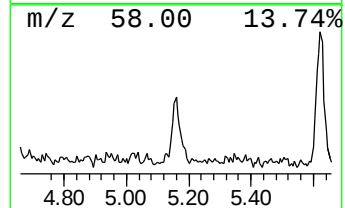
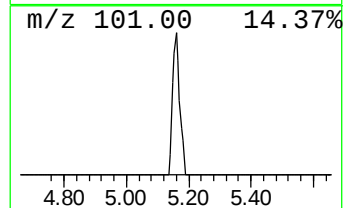
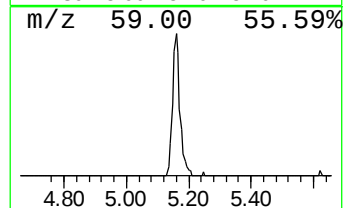
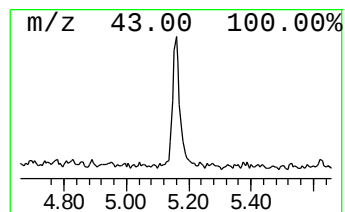
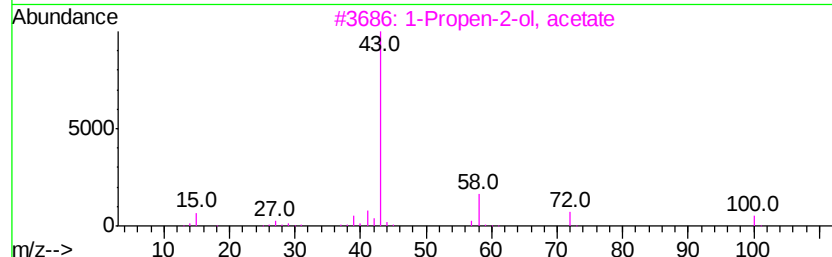
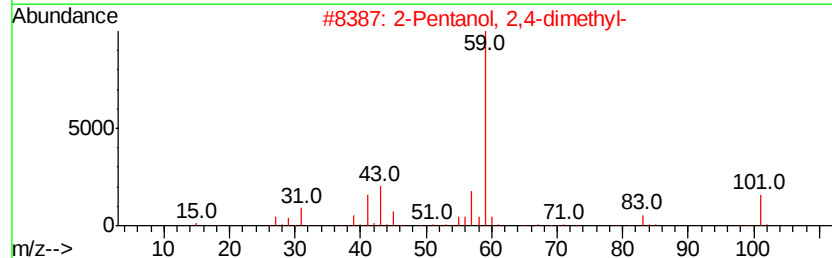
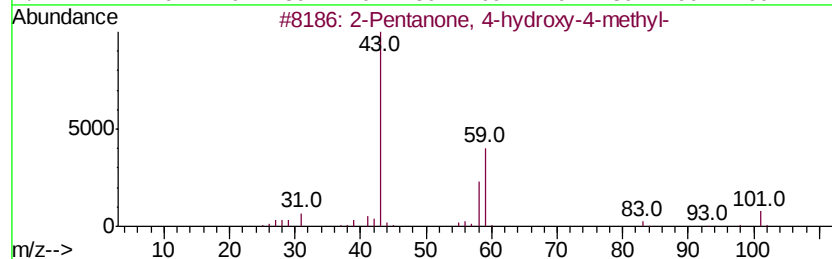
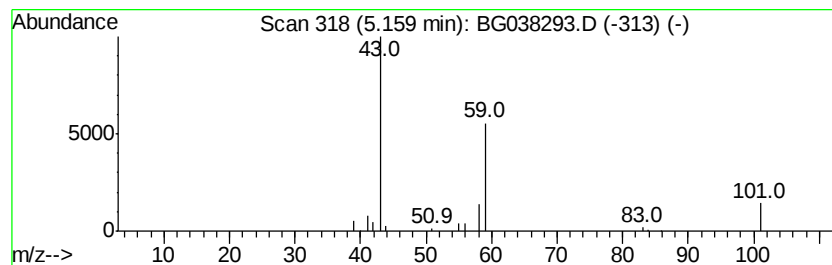
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TIC Library : C:\Database\NIST11.L
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Peak Number 1 unknown5.16 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	4.70 ng	36958	1,4-Dichlorobenzene-d4	8.08

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40		
2	2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	9		
3	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9		
4	Diacetamide	101	C4H7NO2	000625-77-4	7		
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	7		



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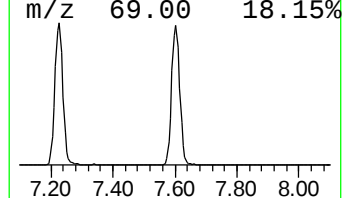
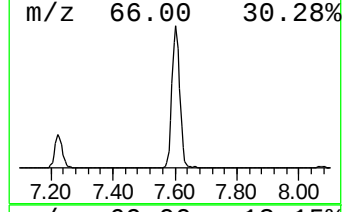
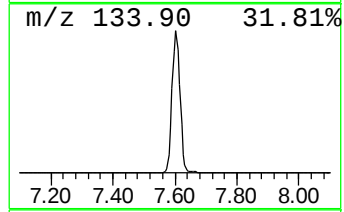
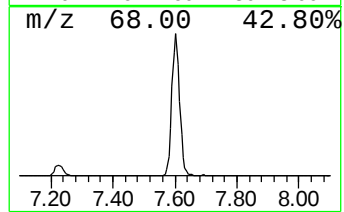
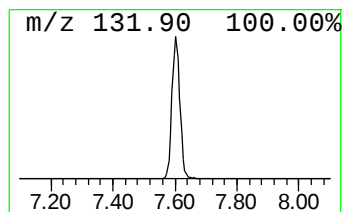
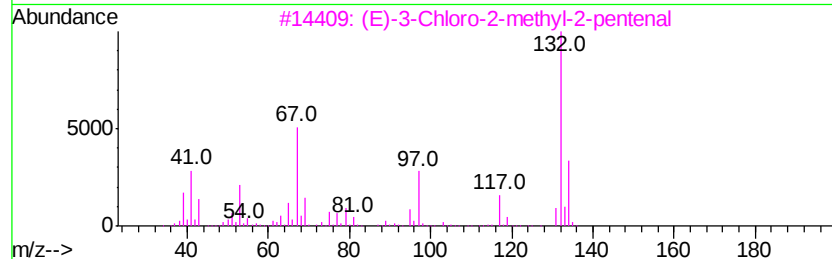
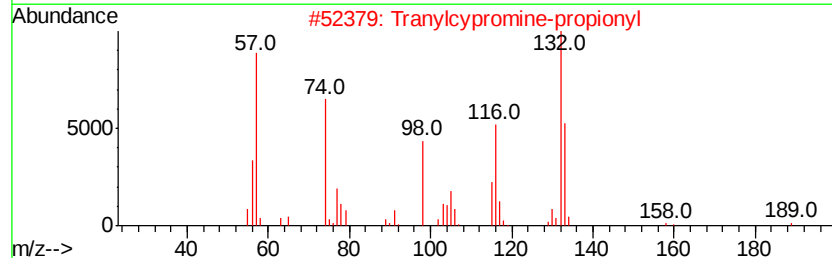
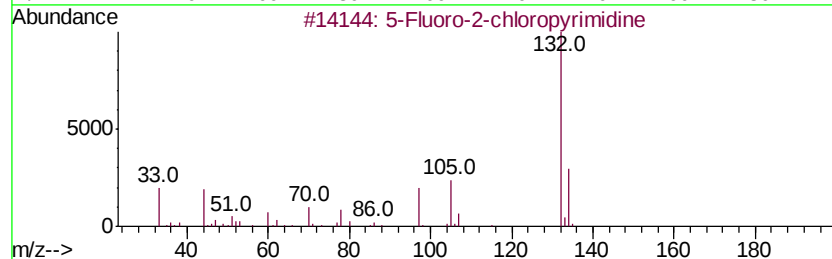
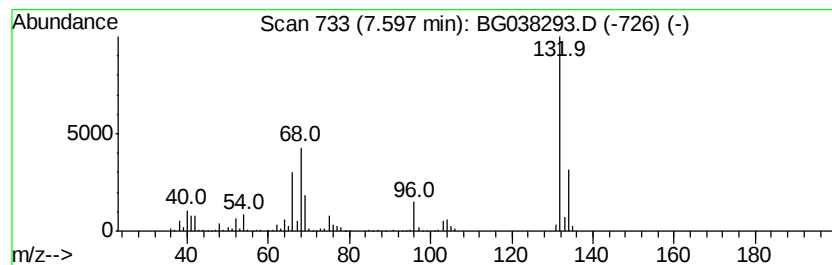
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Peak Number 2 unknown7.60 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.60	98.55 ng	775278	1,4-Dichlorobenzene-d4	8.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
2		Tranlycypramine-propionyl	189	C12H15NO	1000123-86-3	12
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		1,2,4-Trifluorobenzene	132	C6H3F3	000367-23-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown5.16	5.16	4.7	ng	36958	1	8.08	157343	20.0
unknown7.60	7.60	98.5	ng	775278	1	8.08	157343	20.0