

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091018\
 Data File : BG036635.D
 Acq On : 10 Sep 2018 21:37
 Operator : JU/SJ
 Sample : J4776-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EO-02-090718-A

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-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	5.160	308	318	328	rBV	120260	193112	4.38%	0.915%
2	5.624	389	397	407	rBV	963950	1506874	34.18%	7.137%
3	7.210	658	667	678	rBV	903898	1550511	35.17%	7.344%
4	7.586	722	731	743	rBV	990665	1712846	38.85%	8.113%
5	8.057	802	811	820	rVB	209017	350150	7.94%	1.658%
6	8.380	858	866	878	rBV	732900	1296311	29.40%	6.140%
7	9.214	1000	1008	1016	rBV	653000	1162057	26.36%	5.504%
8	10.859	1279	1288	1295	rBV	296654	526230	11.94%	2.492%
9	13.303	1697	1704	1717	rBV	1787935	2833986	64.28%	13.423%
10	14.678	1930	1938	1945	rBV2	490150	799996	18.15%	3.789%
11	16.165	2184	2191	2212	rBV	1096923	1853337	42.04%	8.778%
12	17.422	2398	2405	2414	rBV	657147	1010636	22.92%	4.787%
13	20.031	2842	2849	2861	rBV	3136115	4408680	100.00%	20.881%
14	21.688	3124	3131	3145	rBV	609643	1126759	25.56%	5.337%
15	24.896	3667	3677	3692	rBV3	214094	781872	17.73%	3.703%

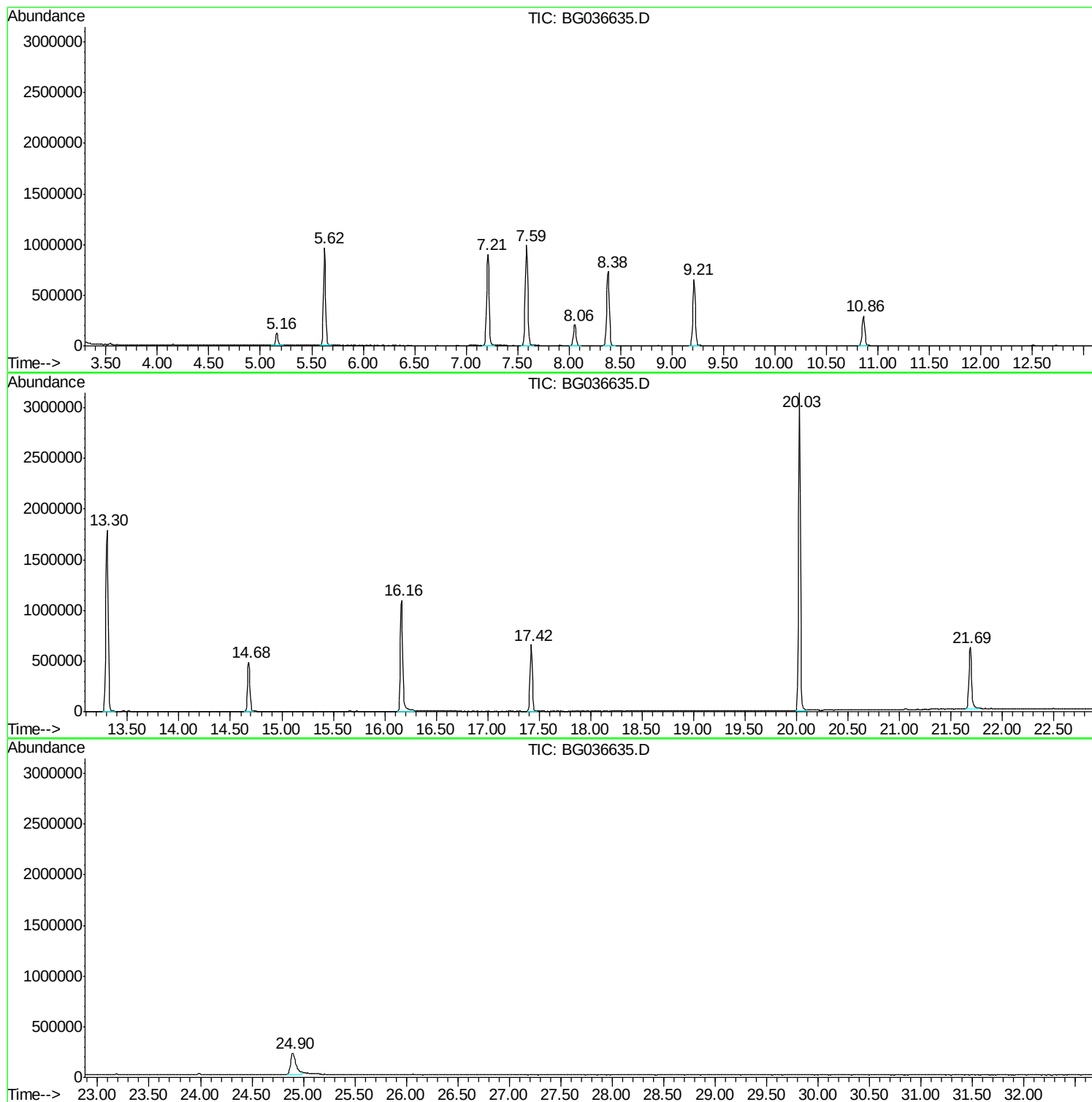
Sum of corrected areas: 21113357

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



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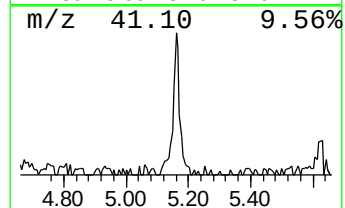
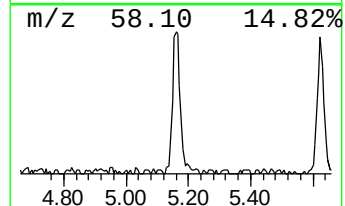
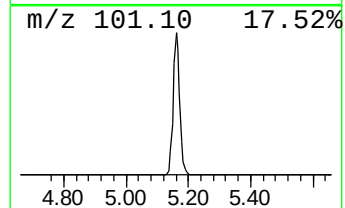
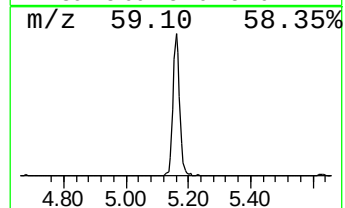
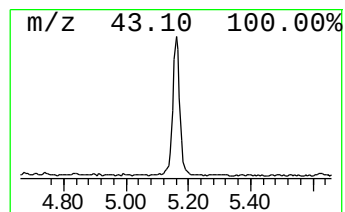
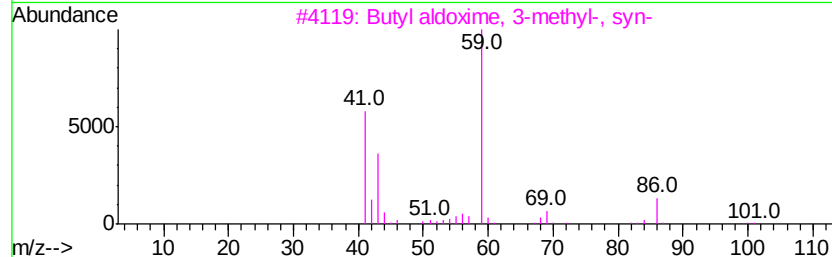
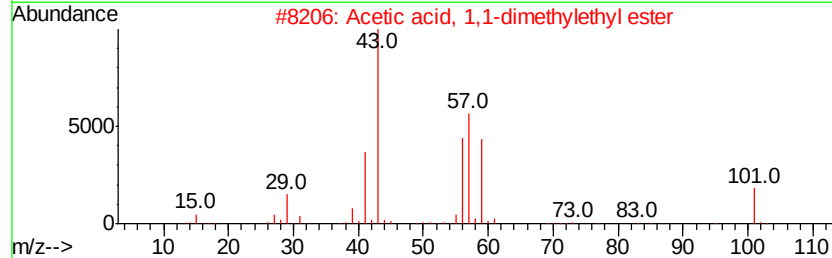
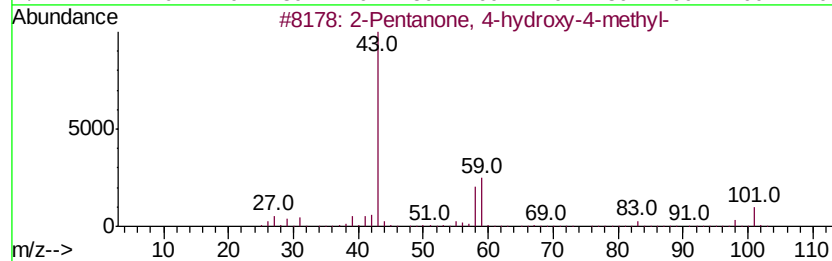
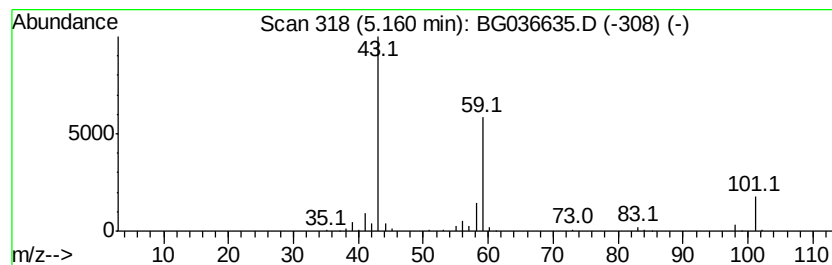
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	11.03 ng	193112	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	50
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	50
3			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	37
4			3-Hexanol	102	C6H14O	000623-37-0	36
5			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	33



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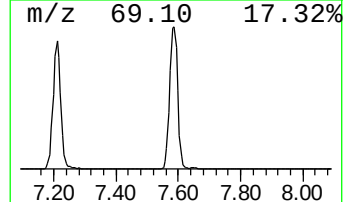
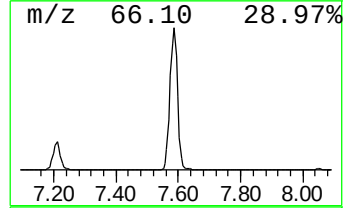
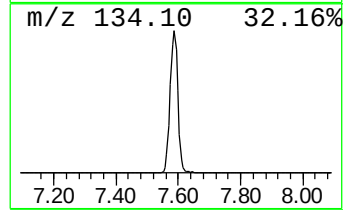
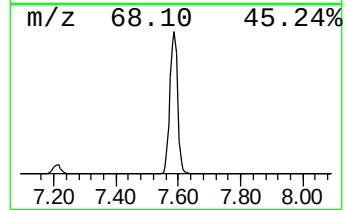
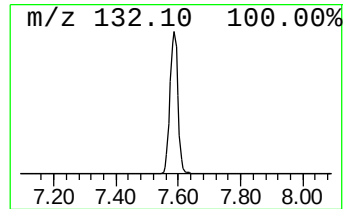
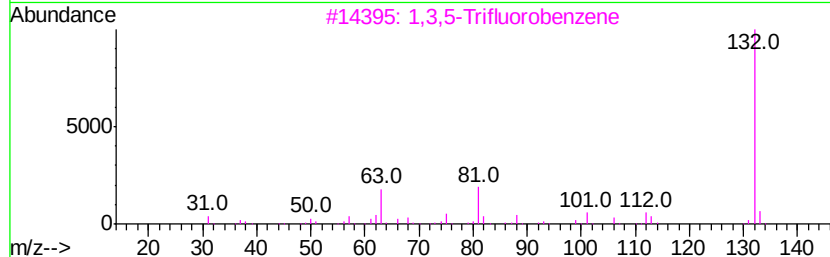
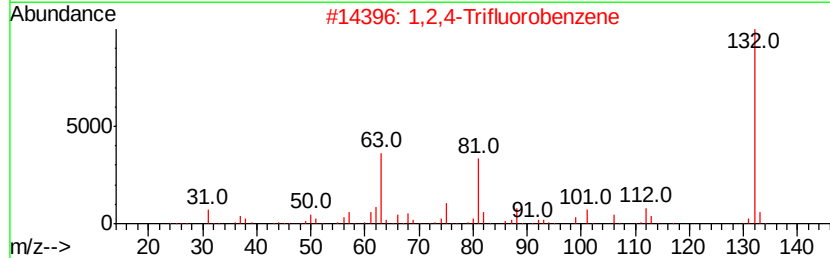
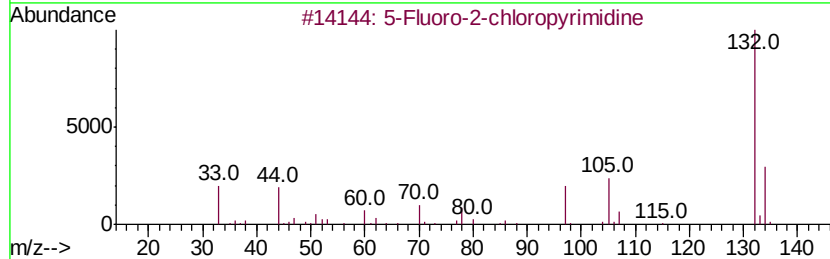
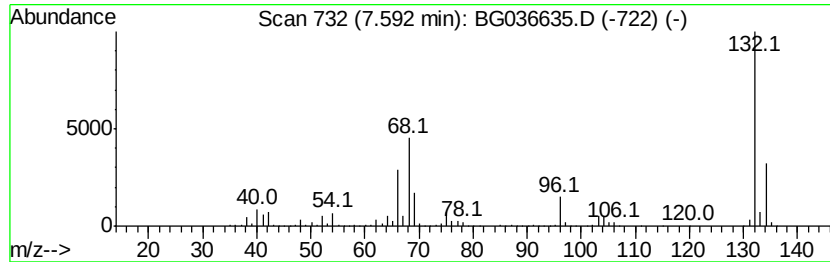
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Peak Number 2 unknown7.59 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	97.84 ng	1712850	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
2		1,2,4-Trifluorobenzene	132	C6H3F3	000367-23-7	9
3		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
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...	5.16	11.0	ng	193112	1	8.06	350150	20.0
unknown7.59	7.59	97.8	ng	1712850	1	8.06	350150	20.0