

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG083118\
 Data File : BG036514.D
 Acq On : 31 Aug 2018 20:26
 Operator : JU/SJ
 Sample : PB112535BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB112535BL

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	5.162	314	319	325	rVB	34223	48873	1.28%	0.265%
2	5.626	391	398	413	rBV	841335	1331233	34.98%	7.218%
3	7.212	661	668	680	rBV	871231	1474615	38.74%	7.995%
4	7.588	725	732	745	rBV	802609	1366082	35.89%	7.407%
5	8.058	806	812	821	rBV	193380	329057	8.65%	1.784%
6	8.382	858	867	875	rBV	594078	1023788	26.90%	5.551%
7	9.216	1002	1009	1019	rBV	483102	866210	22.76%	4.696%
8	10.867	1282	1290	1300	rVB	266276	476922	12.53%	2.586%
9	13.305	1698	1705	1714	rBV	1387189	2180872	57.30%	11.824%
10	14.686	1933	1940	1957	rVB2	434759	710947	18.68%	3.855%
11	16.167	2185	2192	2205	rBV	1042684	1624938	42.69%	8.810%
12	17.424	2399	2406	2415	rBV2	668087	1000045	26.28%	5.422%
13	20.033	2844	2850	2862	rBV	2934856	3806024	100.00%	20.636%
14	21.689	3124	3132	3147	rBV	724160	1202622	31.60%	6.520%
15	24.903	3668	3679	3698	rBV2	303152	1001605	26.32%	5.431%

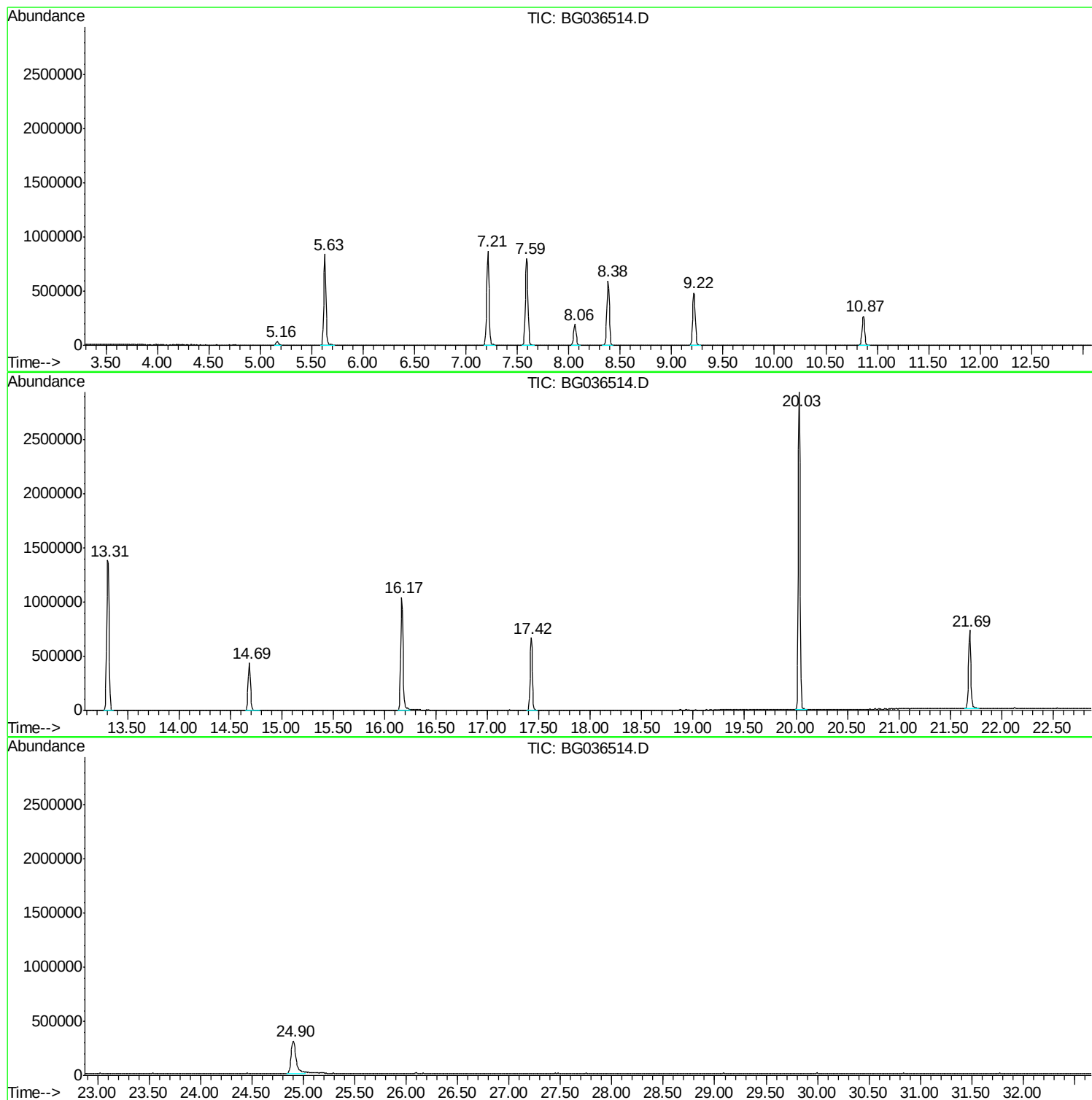
Sum of corrected areas: 18443833

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



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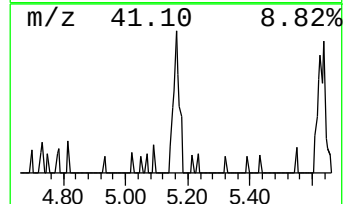
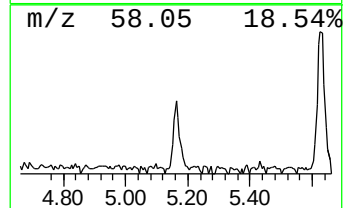
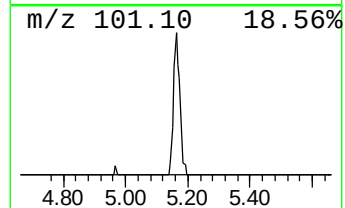
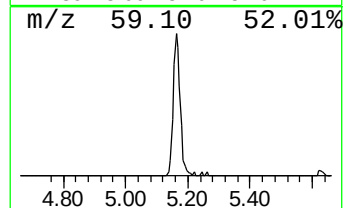
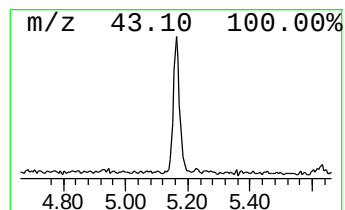
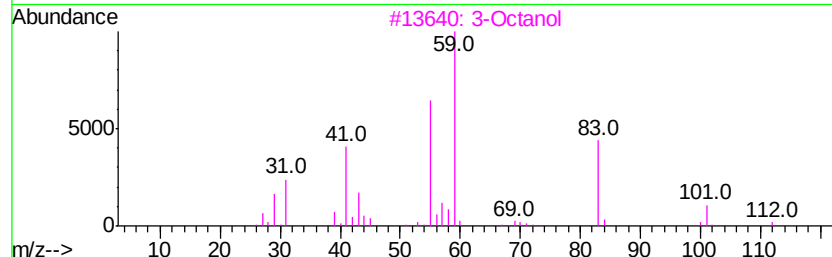
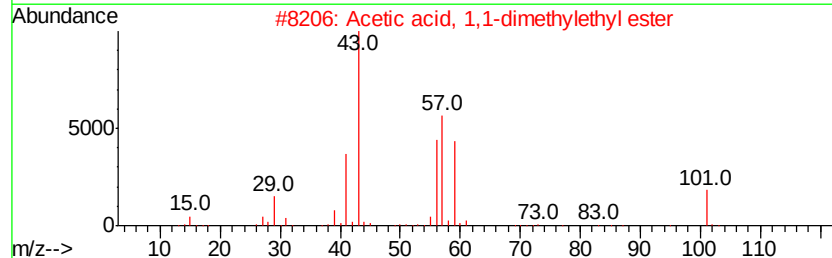
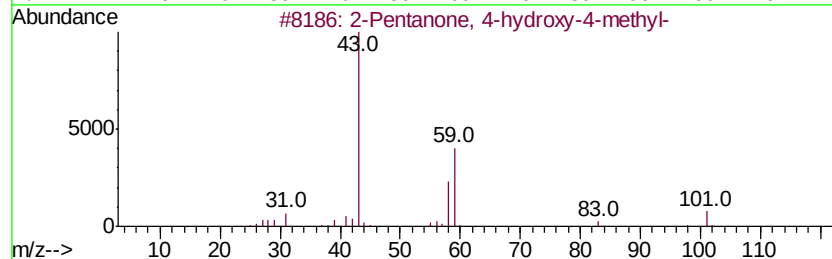
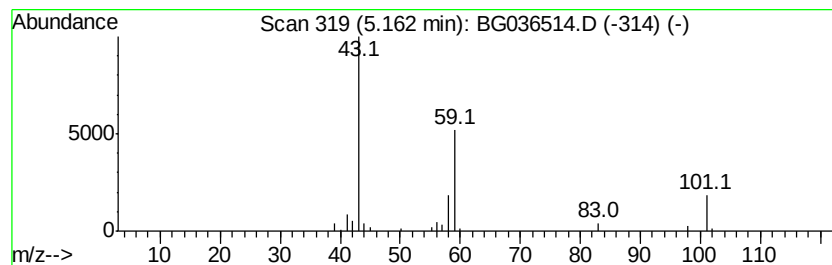
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	2.97 ng	48873	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	56
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3			3-Octanol	130	C8H18O	000589-98-0	28
4			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	23
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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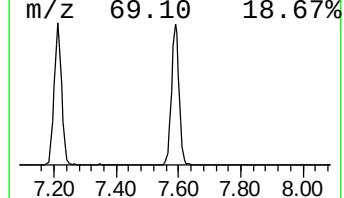
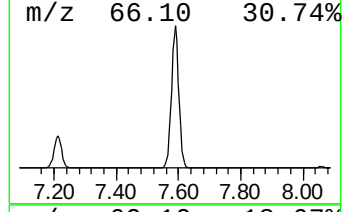
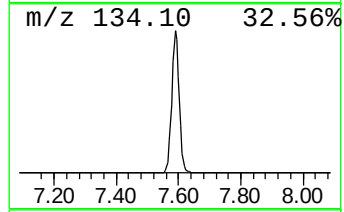
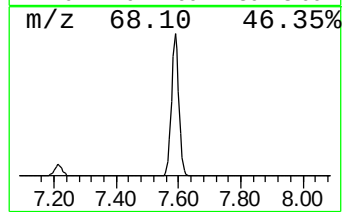
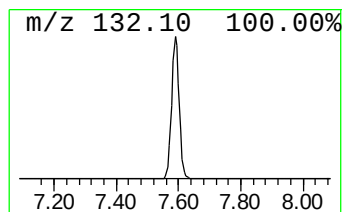
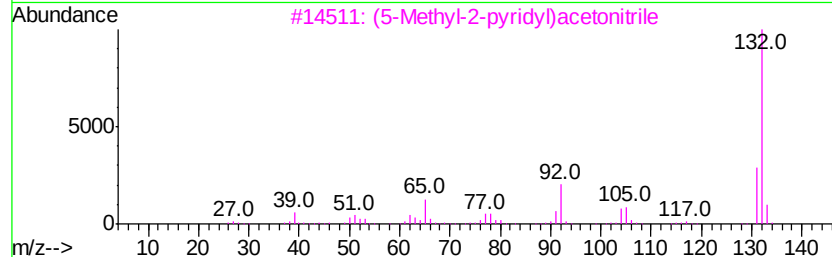
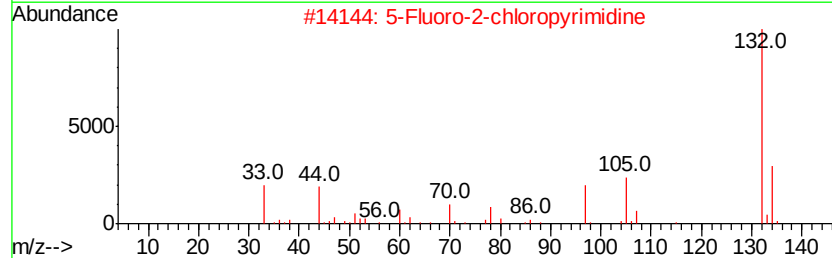
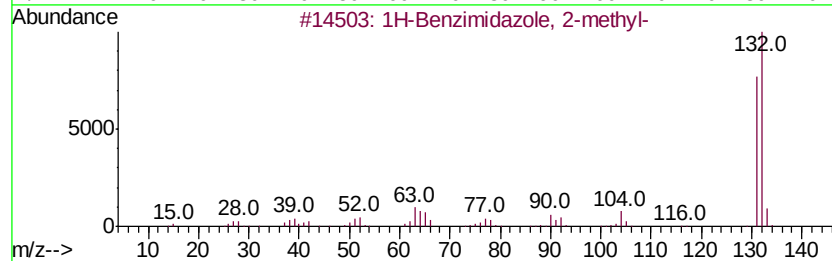
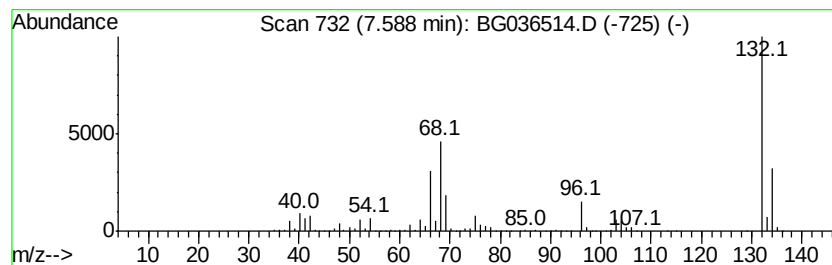
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Peak Number 2 unknown7.59 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	83.03 ng	1366080	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
4		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	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...	5.16	3.0	ng	48873	1	8.06	329057	20.0
unknown7.59	7.59	83.0	ng	1366080	1	8.06	329057	20.0