

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG121418\
 Data File : BG038579.D
 Acq On : 14 Dec 2018 22:25
 Operator : JU/SJ
 Sample : PB115428BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB115428BL

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-BG121218.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	5.139	345	349	356	rVB	16474	26568	1.64%	0.366%
2	5.609	423	429	443	rVB	340693	575309	35.55%	7.919%
3	7.213	694	702	716	rBV	342553	611803	37.81%	8.421%
4	7.589	758	766	776	rBV	320268	579093	35.79%	7.971%
5	8.059	838	846	854	rBV	68892	121504	7.51%	1.672%
6	8.382	893	901	910	rVB	248977	452538	27.96%	6.229%
7	9.234	1039	1046	1058	rBV	197977	377958	23.36%	5.202%
8	10.879	1318	1326	1334	rBV	78948	154319	9.54%	2.124%
9	13.317	1734	1741	1756	rBV	534275	871368	53.85%	11.994%
10	14.698	1969	1976	1984	rBV2	128203	221428	13.68%	3.048%
11	16.185	2223	2229	2249	rBV3	381559	640708	39.59%	8.819%
12	17.442	2435	2443	2452	rBV	188981	295093	18.24%	4.062%
13	20.045	2880	2886	2895	rVB	1171497	1618254	100.00%	22.274%
14	21.719	3164	3171	3178	rBV2	223457	371005	22.93%	5.107%
15	25.015	3721	3732	3745	rVB2	113215	348306	21.52%	4.794%

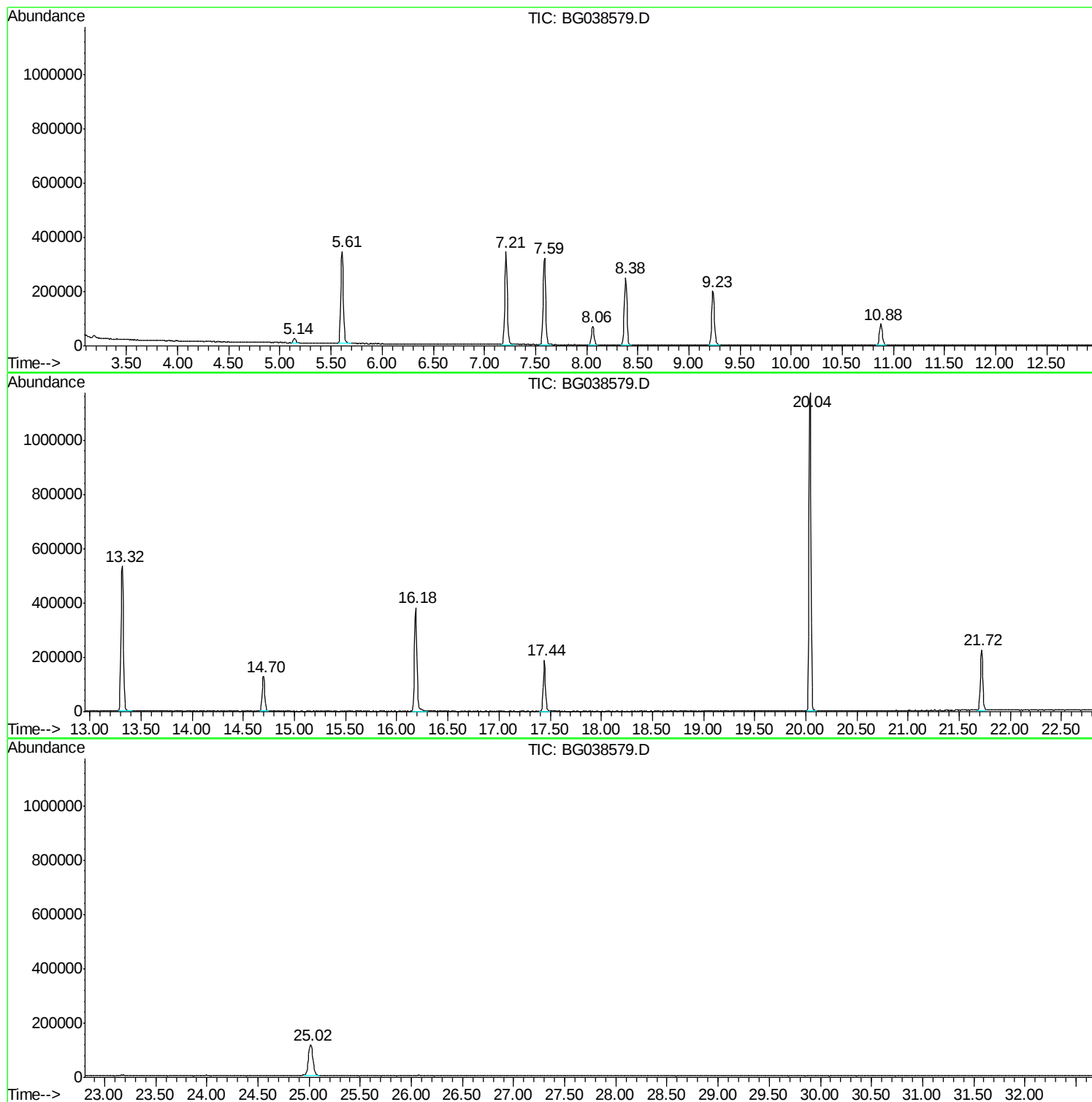
Sum of corrected areas: 7265254

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



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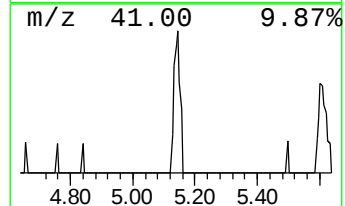
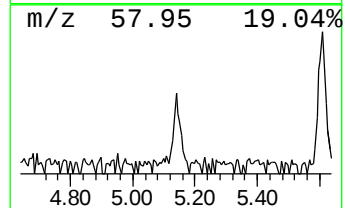
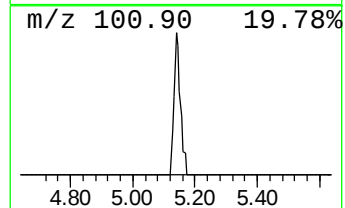
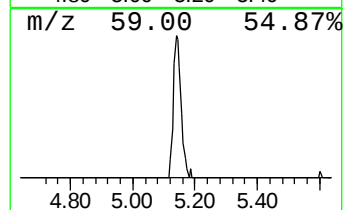
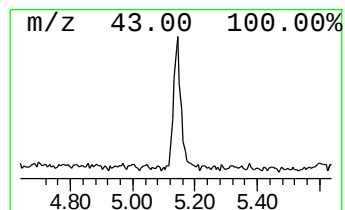
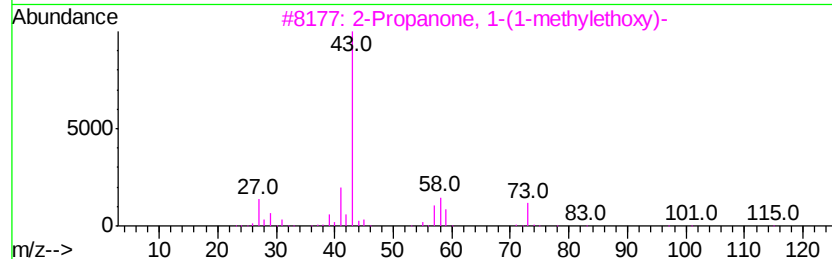
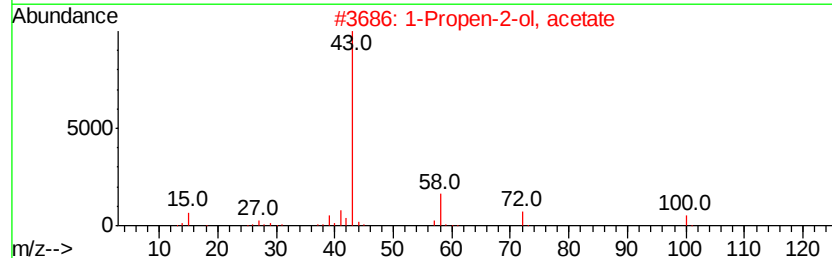
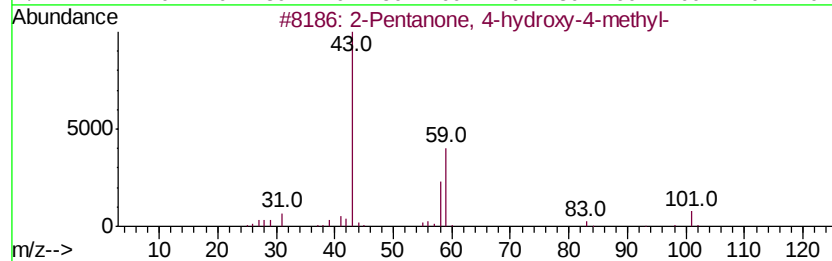
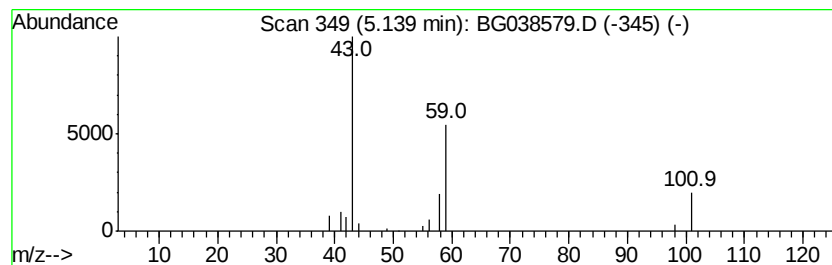
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	4.37 ng	26568	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	39
2			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
3			2-Propanone, 1-(1-methylethoxy)-	116	C6H12O2	042781-12-4	9
4			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	9
5			2-Nonanone, 9-hydroxy-	158	C9H18O2	025368-56-3	9



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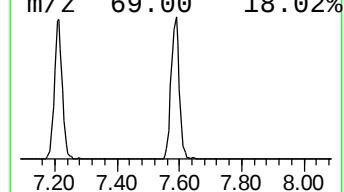
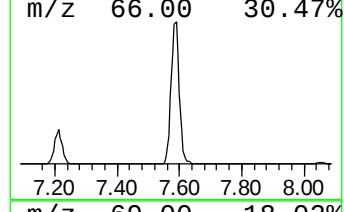
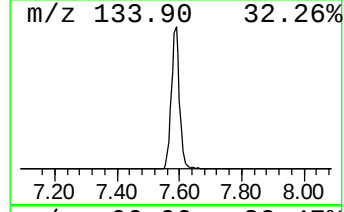
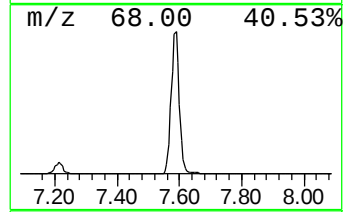
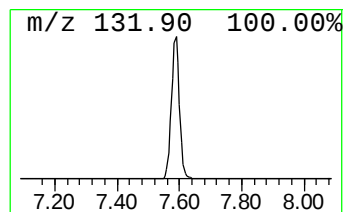
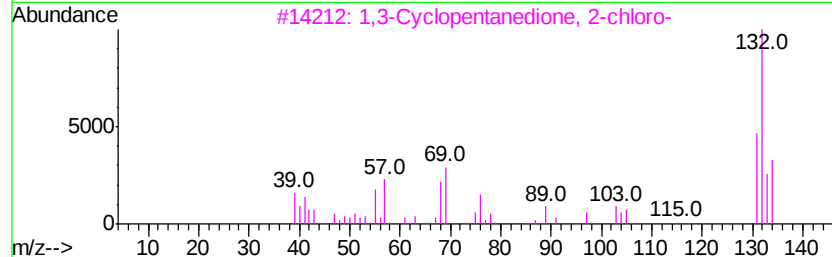
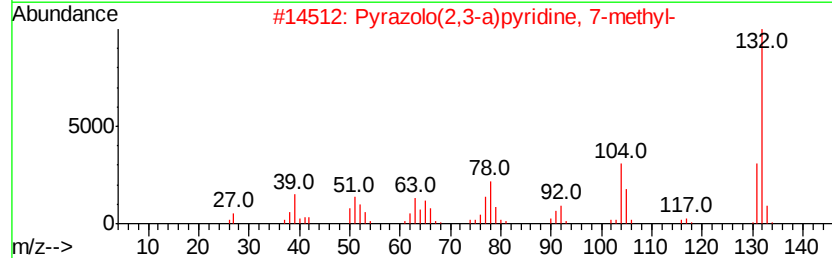
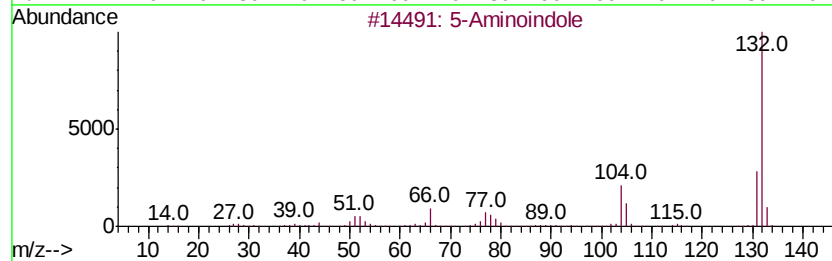
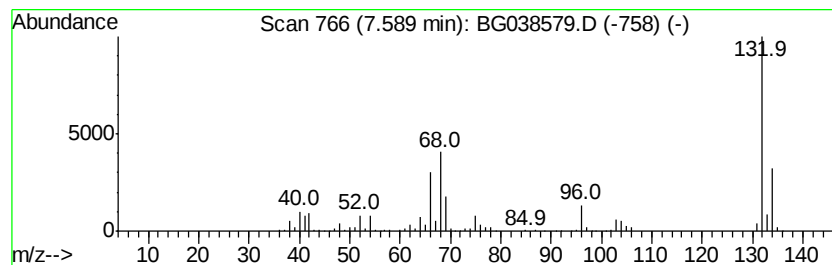
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Peak Number 2 unknown7.59 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	95.32 ng	579093	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Aminoindole	132	C8H8N2	005192-03-0	27
2		Pyrazolo(2,3-a)pyridine, 7-methyl-	132	C8H8N2	034760-58-2	22
3		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	13
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	12
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.14	4.4	ng	26568	1	8.06	121504	20.0
unknown7.59	7.59	95.3	ng	579093	1	8.06	121504	20.0