

# Calculation of Natural Gas Calorific Value with the Agilent 990 PRO Micro GC System

#### Author

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#### **Abstract**

This Application Note gives a functionality introduction of the Agilent 990 process (PRO) Micro GC on how to autonomously analyze and calculate the natural gas energy content.

## Introduction

Natural gas is an important energy resource. It is widely used in different industries to melt, dry, bake, and glaze a product. General households use it for cooking, heating, and lighting. It can also be used as fuel for vehicles. The chemical energy of natural gas is released by the reaction of the natural gas with oxygen. The trade of natural gas is mainly determined by its energy content. Usually, gas chromatography (GC) analysis of natural gas is used to estimate its energy content.

Several organizations, such as the Gas Processors Association (GPA), which collaborates with the American Petroleum Institute, ASTM international formerly American Society of Testing and Materials—and the International Organization for Standardization (ISO) have developed different standards for calculation of natural gas energy based on the individual compound energy values and other physical constants.

The 990 process (PRO) Micro GC can work as an intelligent process GC to quickly analyze the composition of a natural gas stream, then automatically make onboard calculations of its energy content according to the above mentioned standards.

With the assistance of Agilent PROstation software, a user can load predefined energy calculation methods (following GPA/ATSM/ ISO/GOST standards) to the 990 PRO Micro GC. With the completion of each chromatography analysis, the 990 PRO Micro GC generates the normalized mole concentration of each targeted component, then autonomously feeds them to an internal energy content calculation process. Finally, a report for the analyzed gas stream can be generated with information on energy content, including total superior/inferior calorific value (depending on water existing in either the liquid or gaseous state), density, relative density, and Wobbe index.

This Application Note demonstrates the composition analysis and calorific value calculation of natural gas by the 990 PRO Micro GC.

### **Experimental**

Simulated natural gas was analyzed on a two-channel Agilent 990 PRO Micro GC. Channel 1 is a 10 m Agilent J&W CP-PoraPLOT U backflush channel for nitrogen, methane, carbon dioxide, and ethane analysis. Channel 2 is a 6 m Agilent J&W CP-Sil 5CB straight channel for propane, isobutane, butane, 2,2-dimethylpropane, isopentane, pentane, and hexane analysis.

The 990 Micro GC used has PRO and energy meter licenses. The PRO license enables the autonomous running of the instrument at prescheduled time and onboard data processing, including integration, identification, and quantitation based on the preloaded method. The energy meter license enables the automatic onboard calculation of fuel gas energy content based on the PRO GC quantitation result. Table 1 lists the analytical methods used for natural gas composition analysis. The composition of the gas sample is tabulated in Table 2. The analytical parameters were prewritten into the 990 PRO Micro GC mainboard with assistance of Agilent PROstation software. The external standard calibration curve for each targeted component was developed using PROstation.

**Table 1.** Configuration and analytical conditions ofthe Agilent 990 PRO Micro GC.

The Agilent 990 PRO Micro GC Parameters										
Channel Type	10 m Agilent J&W CP-PoraPLOT U, backflush	6 m Agilent J&W CP-Sil 5 CB, straight								
Sampling Time	30 seconds	30 seconds								
Injector Temperature	110 °C	110 °C								
Column Pressure	200 kPa	175 kPa								
Column Temperature	80 °C	70 °C								
Backflush Time	11.3 seconds	NA								

Table 2. Composition of simulated natural gas.

Compound	Concentration (mol%)				
Nitrogen	2.04%				
Carbon Dioxide	3.12%				
Ethane	0.575%				
Propane	0.084%				
Isobutane	0.011%				
Butane	0.011%				
2,2-Dimethylpropane	0.0106%				
Isopentane	0.0097%				
Pentane	0.011%				
Hexane	0.0102%				
Methane	Balance				

The normalization method was defined and written into the mainboard together with the calibration and energy calculation methods prior to the real sample analysis. When the analysis was initiated, the PRO GC deploys these methods for onboard data collection and calculation to generate energy information about the sample. In this application, the calculation method was developed based on ISO standard 6976-2016 as shown in the method setting (Figure 1).

17	Colouistion Mathe		γ	Corre	onent Constar	te			
	Calculation Method			Lomp	urient Lonstar	115			
ISO 6976-1	995 Calorific v	alue unit c	onversion	No Conversio	n	•			
ISO 6976-21	016								
GPA 2172	🗔 Sum C	C6+ uniden	tified compor	nents					
C ASTM D358									
C GOST-2266	57								
C GOST-3136	69								
	🗖 Back	flush to de	tector C6+ Sj	plit					
Combustion T	moorahura 🕞								
	mperature 273.15 K 💌		Caburtada	atas Carrotau t			where the second second	T	
Reference Tem			Saturated W	ater Constant:	2.31	Mole %	at the Selected Reference	e Temperature	
Compressibility /	Air (Zair): 0.99963								
									_
alorific Power	,							==	
alorific Power	Calculation Method		)	Compo	nent Consta	ints			, ) [ [
alorific Power	Calculation Method		)	Compo	nent Consta	ints			, .
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Component	Calculation Method	Index 1	SF 0.0173	Compo MW 28.0135	nent Consta Hs.Mol 0	nts Bj O	Comp. Type 0. Component	_	, (
Component Active C	Calculation Method Sort	_	_	MW	Hs.Mol	Bj			
Component Active C 1 2 2 3 3	Calculation Method Sort Component Name , nitrogen	1	0.0173	MW 28.0135	Hs.Mol 0	Bi O	0. Component		,
Component Active C I I 2 3 4 4 4 4	Calculation Method Sort Component Name . nitrogen 2. methane	1 2 3 4	0.0173 0.0436 0.0728 0.0894	MW 28.0135 16.043 44.01 30.07	Hs.Mol 0 890.63 0 1560.69	Bj 0 0 0 0	0. Component 0. Component 0. Component 0. Component		
Component Active C 2 2 3 3 4 4 5	Calculation Method Sort Component Name . nitrogen 2. methane 3. CO2 . exthane 3. propane	1 2 3 4 5	0.0173 0.0436 0.0728 0.0894 0.1288	MW/ 28.0135 16.043 44.01 30.07 44.097	Hs.Mol 0 890.63 0 1560.69 2219.17	Bi 0 0 0 0 0 0	0. Component 0. Component 0. Component 0. Component 0. Component		
Component Active C 1 2 3 4 4 5 5 5 6 6 6	Calculation Method Sort Introgen Introgen Introgen Introgen I cO2 I. ethane I. ethane I. ethane I. propane I. ibutane	1 2 3 4 5 6	0.0173 0.0436 0.0728 0.0894 0.1288 0.1703	MW 28.0135 16.043 44.01 30.07 44.097 58.123	Hs.Mol 0 890.63 0 1560.69 2219.17 2868.2	Bj 0 0 0 0 0 0 0 0	0. Component 0. Component 0. Component 0. Component 0. Component 0. Component		
Component Active C 1 2 4 4 3 4 4 5 5 5 6 7	Calculation Method Sort Component Name . nitrogen 2. methane 3. CO2 4. ethane 5. propane 3. ibutane 7. mbutane	1 2 3 4 5 6 7	0.0173 0.0436 0.0728 0.0894 0.1288 0.1703 0.1703	MW 28.0135 16.043 44.01 30.07 44.097 58.123 58.123	Hs.Mol 0 890.63 0 1560.69 2219.17 2868.2 2877.4	Bj           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0	0. Component 0. Component 0. Component 0. Component 0. Component 0. Component 0. Component		
Component	Calculation Method Sort Component Name . nitrogen 2. methane 3. CO2 4. ethane 5. propane 5. propane 5. ptotane 7. rubutane 8. neo-pentane	1 2 3 4 5 6 7 8	0.0173 0.0436 0.0728 0.0894 0.1288 0.1703 0.1783 0.2025	MW 28.0135 16.043 44.01 30.07 44.097 58.123 58.123 58.123 72.15	Hs.Mol 0 890.63 0 1560.69 2219.17 2868.2 2877.4 3514.61	Bi           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0	0. Component 0. Component 0. Component 0. Component 0. Component 0. Component 0. Component 0. Component		
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Component Active C Active C Activ	Calculation Method Sort Component Name . nitrogen 2. methane 8. CO2 9. echane 9. copane 9. robutane 9. novpentane 9. neo-pentane 10. jepentane 0. n.pentane	1 2 3 4 5 6 7 8 9 10	0.0173 0.0436 0.0728 0.0894 0.1288 0.1703 0.1783 0.2025 0.2168 0.2345	MW/ 28.0135 16.043 44.01 30.07 58.123 58.123 58.123 72.15 72.15 72.15	Hs.Mol           0           890.63           0           1560.69           2219.17           2868.2           2877.4           3514.61           3528.83           3535.77	Bi           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0	O. Component     O. Component		
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Component Active C Active C Activ	Calculation Method Sort Component Name . nitrogen 2. methane 8. CO2 9. echane 9. copane 9. robutane 9. novpentane 9. neo-pentane 10. jepentane 0. n.pentane	1 2 3 4 5 6 7 8 9 10	0.0173 0.0436 0.0728 0.0894 0.1288 0.1703 0.1783 0.2025 0.2168 0.2345	MW/ 28.0135 16.043 44.01 30.07 58.123 58.123 58.123 72.15 72.15 72.15	Hs.Mol           0           890.63           0           1560.69           2219.17           2868.2           2877.4           3514.61           3528.83           3535.77	Bi           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0	O. Component     O. Component		
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Component Active C Active C Activ	Calculation Method Sort Component Name . nitrogen 2. methane 8. CO2 9. echane 9. copane 9. robutane 9. novpentane 9. neo-pentane 10. jepentane 0. n.pentane	1 2 3 4 5 6 7 8 9 10	0.0173 0.0436 0.0728 0.0894 0.1288 0.1703 0.1783 0.2025 0.2168 0.2345	MW/ 28.0135 16.043 44.01 30.07 58.123 58.123 58.123 72.15 72.15 72.15	Hs.Mol           0           890.63           0           1560.69           2219.17           2868.2           2877.4           3514.61           3528.83           3535.77	Bi           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0	O. Component     O. Component		
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Component Active C Active C Activ	Calculation Method Sort Component Name . nitrogen 2. methane 8. CO2 9. echane 9. copane 9. robutane 9. novpentane 9. neo-pentane 10. jepentane 0. n.pentane	1 2 3 4 5 6 7 8 9 10	0.0173 0.0436 0.0728 0.0894 0.1288 0.1703 0.1783 0.2025 0.2168 0.2345	MW/ 28.0135 16.043 44.01 30.07 58.123 58.123 58.123 72.15 72.15 72.15	Hs.Mol           0           890.63           0           1560.69           2219.17           2868.2           2877.4           3514.61           3528.83           3535.77	Bi           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0	O. Component     O. Component		
Component Active C Active C Activ	Calculation Method Sort Component Name . nitrogen 2. methane 8. CO2 9. echane 9. copane 9. robutane 9. novpentane 9. neo-pentane 10. jepentane 0. n.pentane	1 2 3 4 5 6 7 8 9 10	0.0173 0.0436 0.0728 0.0894 0.1288 0.1703 0.1783 0.2025 0.2168 0.2345	MW/ 28.0135 16.043 44.01 30.07 58.123 58.123 58.123 72.15 72.15 72.15	Hs.Mol           0           890.63           0           1560.69           2219.17           2868.2           2877.4           3514.61           3528.83           3535.77	Bi           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0	O. Component     O. Component		
Component Active C Active C Activ	Calculation Method Sort Component Name . nitrogen 2. methane 8. CO2 9. echane 9. copane 9. robutane 9. novpentane 9. neo-pentane 10. jepentane 0. n.pentane	1 2 3 4 5 6 7 8 9 10	0.0173 0.0436 0.0728 0.0894 0.1288 0.1703 0.1783 0.2025 0.2168 0.2345	MW/ 28.0135 16.043 44.01 30.07 58.123 58.123 58.123 72.15 72.15 72.15	Hs.Mol           0           890.63           0           1560.69           2219.17           2868.2           2877.4           3514.61           3528.83           3535.77	Bi           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0           0	O. Component     O. Component		

Figure 1. Settings of the calorific value calculation method based on ISO standard in Agilent PROstation software.

## **Results and discussion**

Figures 2a and 2b show the chromatograms generated on channel 1 and channel 2 for simulated natural gas sample. The peaks were integrated onboard according to the integration parameters optimized for each analytical channel. The integration result was used to generate the quantitation result for each target component based on the prewritten ESTD curves. The concentration normalization was made over two analytical channels for all target components according to the setting in the normalization table (Figure 3). Then, the normalized concentrations were used for onboard calorific value calculation according to predefined energy calculation methods.

Figure 4 shows the quantitation and energy content calculation results for the simulated natural gas. The "energy" part shows the standard which is followed for calculation, and the key physical properties required to be calculated in the standard, such as compressibility, relative density/density, superior/inferior heating value in molar/weight/volume units, and Wobbe index. The types of properties shown in the report vary according to the standard requirement.

The quantitation results based on chromatography analysis are shown in the bottom of the report. Both the ESTD concentration and normalized concentration are reported. The report was generated in Agilent PROstation software as soon as each

chromatography separation was finished. The analysis cycle under the applied conditions in this work is approximately 90 seconds from sampling to separation and report generated. If continuous flow mode is used, the analysis cycle time can be further reduced to 60s.

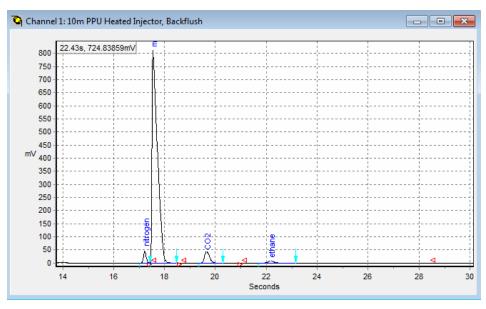


Figure 2a. Chromatogram of  $N_2$ /methane/CO<sub>2</sub>/ethane on 10 m Agilent J&W CP-PoraPLOT U backflush channel.

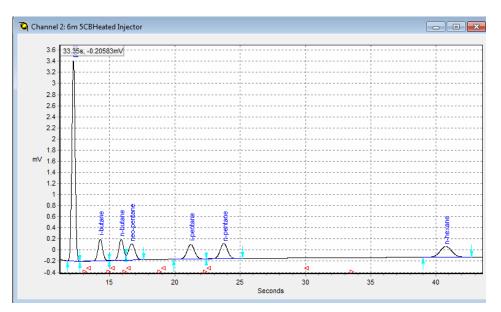


Figure 2b. Chromatogram of C3-C6 hydrocarbons on the 6 m Agilent J&W CP-Sil 5 CB channel.

M N	ormalizat	ion Table									
	Synchron	ize									
#	Active	Peak Name	Channel	Ignore	Bridge Comp #	Estimate	Estim.Conc	Test.Conc	RefConcPeak#	RefPeakConc%	Group#
1	<	nitrogen	1		0. None		0	0	0	0	0
2	$\checkmark$	methane	1		0. None		0	0	0	0	0
3		C02	1		0. None		0	0	0	0	0
4	$\checkmark$	ethane	1		0. None		0	0	0	0	0
5	$\checkmark$	propane	2		0. None		0	0	0	0	0
6	$\checkmark$	i-butane	2		0. None		0	0	0	0	0
7	$\checkmark$	n-butane	2		0. None		0	0	0	0	0
8	$\checkmark$	neo-pentane	2		0. None		0	0	0	0	0
9	$\checkmark$	i-pentane	2		0. None		0	0	0	0	0
10	✓	n-pentane	2		0. None		0	0	0	0	0
11	✓	n-hexane	2		0. None		0	0	0	0	0

Figure 3. Normalization table settings for natural gas analysis in this work.

SAM	PLE				ENERGY							CONDITIONS		
Samp	ling Time		08/10/2019	14:10:50	Calc.Method		ISO 6976-20	16 Dry	Saturat	ed				
Run N	Number		3		Water Mole.		[%]		2.31			ENVIRONMENT		
Run T	Гуре		Analysis		Compressibility		[•]	0.9981	0.9975			Cabinet Temperature	[°C]	34
Calibr	ation Level	1	0		Molar Mass		[kg/kmol]	17.2666	17.283	Э		Ambient Pressure	[kPa]	102.1
Stream	m #		1	(checked)	Relative Density	,Ideal	[•]	0.5961	0.5967					
Sum B	ESTD		1.0238		Relative Density	,Real	[•]	0.5971	0.5980			SITE INFO		
Sum B	Estimates		0.0000		Gas Density,Ide	al	[kg/m3]	0.7178	0.7185			Customer ID		
ium A	Areas		1130262.377	'5	Gas Density,Re	al	[kg/m3]	0.7192	0.7203			Instrument Name		990-PRO Micro GC
otal	Peaks		11		Superior Heatin	g Value (Volume R	eal) [MJ/m3]	35.60	34.79			Serial Number		10001
s Sta	rtup Run		False		Inferior Heating	Value (Volume Re	al) [MJ/m3]	32.01	31.29			Tag Number		
Jnkn	own Peaks	5	6		Superior Heatin	g Value (Volume Id	ea) [MJ/m3]	35.53	34.71			Cylinder 1 Tag		
Curren	nt Stream ‡	ŧ	0		Inferior Heating	Value (Volume Ide	a) [MJ/m3]	31.95	31.21					
					Superior Heatin	g Value(Mass)	[MJ/kg]	49.50	48.30					
					Inferior Heating	Value(Mass)	[MJ/kg]	44.51	43.44					
					Superior Heatin	g Value(Molar)	[kJ/mol]	854.62	834.88					
			Inferior Heating Value(Molar)		[kJ/mol]	768.57	750.82	750.82						
	Hide non A				Wobbe Index (Real )		[MJ/m3]	46.07	44.99					
Hide Ignored Appl.pks			Wobbe Index inferior		[MJ/m3]	41.43	40.46							
#	Channel	Peakna	ame		ESTD Conc.	Norm. Conc.	Retention [s]	Area	Height	Meth-Index	Group#	R.F.		Weight%
1	1	nitroger	n		0.019951	1.948797	17.38	24782.7169	13635198.2879	1	0	8.0504E-07		3.1617
2	1	methan	ie		0.965245	94.283973	17.60	746196.9431	132556588.1529	2	0	1.293553E-06		87.5998
3	1	CO2			0.031328	3.060127	19.75	38813.4314	6589405.0554	3	0	8.071555E-07		7.7997
4	1	ethane			0.005773	0.563867	22.33	7757.2239	1375294.4939	4	0	7.441665E-07		0.9820
5	2	propan	e		0.000837	0.081791	12.24	1714.7951	352927.5260	5	0	4.883076E-07		0.2089
6	2	i-butane	e		0.000107	0.010478	14.29	290.7882	39425.5766	6	0	3.688964E-07		0.0353
7	2	n-butar	ne		0.000106	0.010367	15.89	283.3521	38176.9311	7	0	3.745673E-07		0.0349
8	2	neo-pei	ntane		0.000106	0.010374	16.69	282.8061	29531.1083	8	0	3.755436E-07		0.0433
9	2	i-pentar	ne		0.000097	0.009514	21.22	290.1013	26104.6298	9	0	3.357591E-07		0.0398
10	2	n-penta	ane		0.000110	0.010771	23.74	300.9403	27215.5715	10	0	3.664084E-07		0.0450
11	2	n-hexar	ne		0.000102	0.009941	40.74	332.8352	19101.1810	11	0	3.057677E-07		0.0496

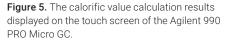
Figure 4. Energy content calculation report generated by the Agilent 990 PRO Micro GC.

The 990 PRO Micro GC works as a "detector" or "sensor" for natural gas stream analysis. PROstation software is used to

- Develop methods, including analytical, qualitative and quantitative method, and energy content calculation method
- Set the automation mode
- Define how to output the results

All these "commands" are written to the mainboard by PROstation. In the real analysis, the Pro Micro GC can run by itself without connection to the PROstation software. The quantitation result and energy content will not be displayed in the format shown here if PROstation software is not connected. Instead, the results can be displayed in a scrollable way on the touch screen of the 990 PRO Micro GC, as shown in Figure 5. Additionally, the result can be output through FTP in .txt file or using Modbus protocol to other terminals for monitoring and recording. Analog output is another approach to produce analysis results in voltage or current signal through a connection to the extension analog board (Figure 6). The conversion of analog signal and the quantitation result or energy contents can be predefined and loaded on the PRO GC mainboard.





#### Conclusion

This Application Note demonstrates natural gas composition analysis and energy content calculation by the Agilent 990 PRO Micro GC. The PRO license and energy content license are activated on the 990 PRO Micro GC to enable automated fuel gas composition analysis and energy content calculation. The analysis process from sampling, separation, quantitation to calorific value calculation and result output, are autonomously executed according to the prewritten method and automation mode on the PRO Micro GC mainboard. The energy calculation methods are developed with compliance to different international standards, including ASTM, ISO, GPA, and GOST standards. All methods are developed in PROstation and then downloaded to the 990 PRO Micro GC for its independent and automated operation. The energy content calculation results can be shown on a local touch screen or outputted through FTP, Modbus, and analog signal for monitoring and recording.

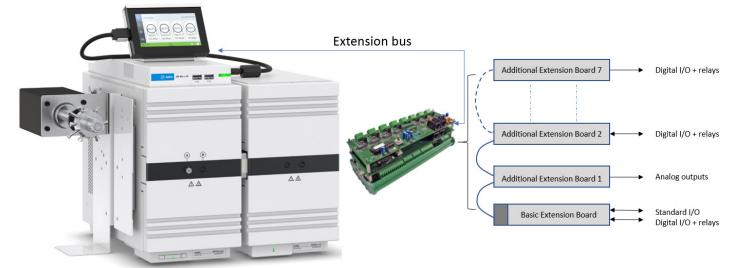


Figure 6. Agilent 990 PRO Micro GC connection with extension board for analog output of analysis results.

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