



# Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 08:00 PM GMT

PDB ID : 1H4O  
Title : Monoclinic form of human peroxiredoxin 5  
Authors : Declercq, J.P.; Evrard, C.  
Deposited on : 2001-05-11  
Resolution : 1.95 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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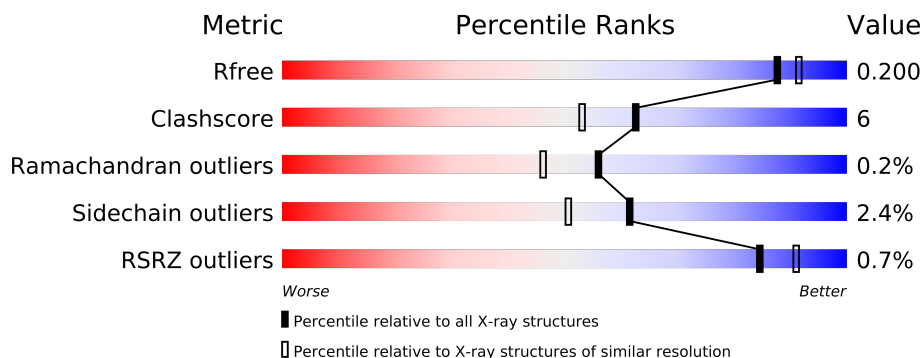
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 1.95 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	66092	1321 (1.96-1.96)
Clashscore	79885	1488 (1.96-1.96)
Ramachandran outliers	78287	1475 (1.96-1.96)
Sidechain outliers	78261	1475 (1.96-1.96)
RSRZ outliers	66119	1321 (1.96-1.96)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	161	
1	B	161	
1	C	161	
1	D	161	
1	E	161	
1	F	161	
1	G	161	
1	H	161	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	BEZ	A	1162	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
2	BEZ	G	1162	-	X

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10805 atoms, of which 0 are hydrogen and 0 are deuterium.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

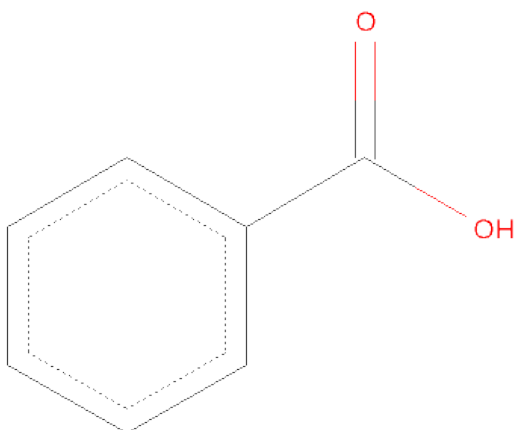
- Molecule 1 is a protein called PEROXIREDOXIN 5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	161	Total	C	N	O	S	0	0	0
			1190	760	204	222	4			
1	B	161	Total	C	N	O	S	0	0	0
			1190	760	204	222	4			
1	C	161	Total	C	N	O	S	0	0	0
			1190	760	204	222	4			
1	D	161	Total	C	N	O	S	0	0	0
			1190	760	204	222	4			
1	E	161	Total	C	N	O	S	0	0	0
			1190	760	204	222	4			
1	F	161	Total	C	N	O	S	0	0	0
			1190	760	204	222	4			
1	G	161	Total	C	N	O	S	0	0	0
			1190	760	204	222	4			
1	H	161	Total	C	N	O	S	0	0	0
			1190	760	204	222	4			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	88	HIS	THR	VARIANT	UNP P30044
B	88	HIS	THR	VARIANT	UNP P30044
C	88	HIS	THR	VARIANT	UNP P30044
D	88	HIS	THR	VARIANT	UNP P30044
E	88	HIS	THR	VARIANT	UNP P30044
F	88	HIS	THR	VARIANT	UNP P30044
G	88	HIS	THR	VARIANT	UNP P30044
H	88	HIS	THR	VARIANT	UNP P30044

- Molecule 2 is BENZOIC ACID (three-letter code: BEZ) (formula: C<sub>7</sub>H<sub>6</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			9	7	2		
2	B	1	Total	C	O	0	0
			9	7	2		
2	C	1	Total	C	O	0	0
			9	7	2		
2	D	1	Total	C	O	0	0
			9	7	2		
2	E	1	Total	C	O	0	0
			9	7	2		
2	F	1	Total	C	O	0	0
			9	7	2		
2	G	1	Total	C	O	0	0
			9	7	2		
2	H	1	Total	C	O	0	0
			9	7	2		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	166	Total	O	0	0
			166	166		
3	B	126	Total	O	0	0
			126	126		
3	C	188	Total	O	0	0
			188	188		
3	D	184	Total	O	0	0
			184	184		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	128	Total 128	O 128	0	0
3	F	160	Total 160	O 160	0	0
3	G	139	Total 139	O 139	0	0
3	H	122	Total 122	O 122	0	0

### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: PEROXIREDOXIN 5

Chain A: 



- Molecule 1: PEROXIREDOXIN 5

Chain B: 



- Molecule 1: PEROXIREDOXIN 5

Chain C: 



- Molecule 1: PEROXIREDOXIN 5

Chain D: 



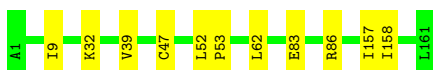
- Molecule 1: PEROXIREDOXIN 5

Chain E: 

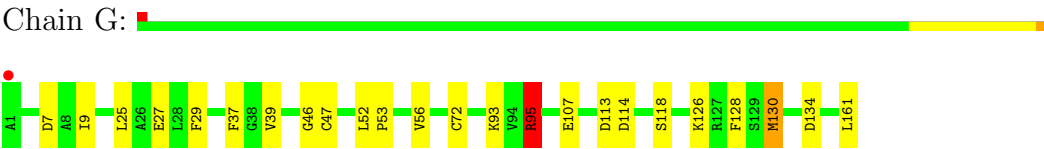


- Molecule 1: PEROXIREDOXIN 5

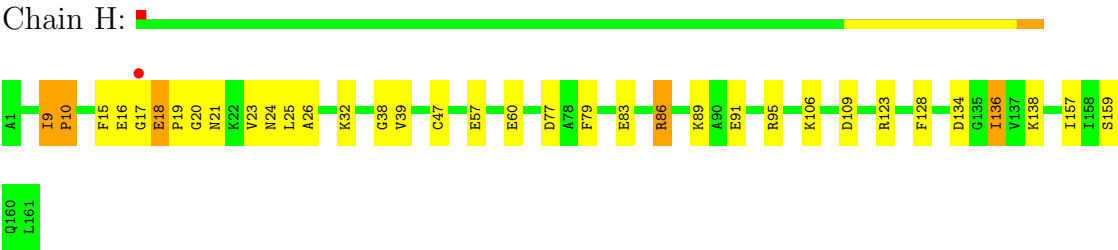
Chain F: 



● Molecule 1: PEROXIREDOXIN 5



● Molecule 1: PEROXIREDOXIN 5





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	130.79Å 66.49Å 141.24Å 90.00° 117.54° 90.00°	Depositor
Resolution (Å)	24.00 – 1.95 24.00 – 1.95	Depositor EDS
% Data completeness (in resolution range)	94.1 (24.00-1.95) 94.1 (24.00-1.95)	Depositor EDS
$R_{merge}$	0.03	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.92 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.0	Depositor
R, $R_{free}$	0.165 , 0.190 0.176 , 0.200	Depositor DCC
$R_{free}$ test set	7402 reflections (5.27%)	DCC
Wilson B-factor (Å <sup>2</sup> )	24.1	Xtriage
Anisotropy	0.041	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.41 , 47.5	EDS
Estimated twinning fraction	0.093 for h,-k,-h-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	0 of 147800 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	10805	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.80% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 5 Model quality ⓘ

### 5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: BEZ

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.81	0/1211	0.92	1/1637 (0.1%)
1	B	0.82	0/1211	0.99	3/1637 (0.2%)
1	C	0.82	0/1211	0.93	2/1637 (0.1%)
1	D	0.84	0/1211	0.90	2/1637 (0.1%)
1	E	0.76	0/1211	0.93	2/1637 (0.1%)
1	F	0.78	0/1211	0.90	1/1637 (0.1%)
1	G	0.82	1/1211 (0.1%)	0.94	6/1637 (0.4%)
1	H	0.79	0/1211	0.93	4/1637 (0.2%)
All	All	0.80	1/9688 (0.0%)	0.93	21/13096 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	130	MET	SD-CE	-7.04	1.38	1.77

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	114	ASP	CB-CG-OD2	7.15	124.73	118.30
1	E	95	ARG	NE-CZ-NH2	-6.31	117.14	120.30
1	A	113	ASP	CB-CG-OD1	6.21	123.89	118.30
1	H	86	ARG	NE-CZ-NH1	6.17	123.39	120.30
1	B	114	ASP	CB-CG-OD2	5.94	123.64	118.30
1	B	134	ASP	CB-CG-OD2	5.81	123.53	118.30
1	F	32	LYS	CB-CA-C	-5.77	98.87	110.40
1	G	95	ARG	NE-CZ-NH1	-5.66	117.47	120.30
1	G	114	ASP	CB-CG-OD2	5.62	123.35	118.30
1	G	113	ASP	CB-CG-OD1	5.43	123.18	118.30
1	H	134	ASP	CB-CG-OD2	5.42	123.18	118.30
1	B	19	PRO	N-CA-C	5.40	126.14	112.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	124	ARG	NE-CZ-NH2	-5.36	117.62	120.30
1	E	95	ARG	NE-CZ-NH1	5.26	122.93	120.30
1	G	7	ASP	CB-CG-OD2	5.22	123.00	118.30
1	H	109	ASP	CB-CG-OD2	5.22	123.00	118.30
1	G	25	LEU	CB-CG-CD1	-5.21	102.14	111.00
1	H	77	ASP	CB-CG-OD2	5.04	122.84	118.30
1	D	130	MET	CB-CA-C	-5.02	100.36	110.40
1	C	86	ARG	NE-CZ-NH2	-5.01	117.79	120.30
1	G	95	ARG	CG-CD-NE	5.01	122.33	111.80

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1190	0	1224	6	0
1	B	1190	0	1224	13	0
1	C	1190	0	1224	5	0
1	D	1190	0	1224	6	0
1	E	1190	0	1224	19	0
1	F	1190	0	1224	6	0
1	G	1190	0	1224	19	0
1	H	1190	0	1224	37	0
2	A	9	0	5	1	0
2	B	9	0	5	0	0
2	C	9	0	5	0	0
2	D	9	0	5	0	0
2	E	9	0	5	0	0
2	F	9	0	5	0	0
2	G	9	0	5	1	0
2	H	9	0	5	0	0
3	A	166	0	0	2	0
3	B	126	0	0	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	188	0	0	2	0
3	D	184	0	0	1	0
3	E	128	0	0	4	0
3	F	160	0	0	0	0
3	G	139	0	0	6	0
3	H	122	0	0	14	0
All	All	10805	0	9832	109	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 6.

All (109) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:136:ILE:HD11	1:E:138:LYS:HE3	1.32	1.10
1:H:83:GLU:CD	1:H:86:ARG:HH12	1.73	0.92
1:E:136:ILE:HD11	1:E:138:LYS:CE	2.00	0.90
1:H:83:GLU:CD	1:H:86:ARG:NH1	2.26	0.88
1:E:83:GLU:CD	1:E:86:ARG:HH12	1.80	0.85
1:E:83:GLU:OE2	1:E:86:ARG:NH1	2.16	0.78
1:G:9:ILE:HD11	1:G:130:MET:HE1	1.64	0.77
1:H:16:GLU:O	1:H:16:GLU:CG	2.33	0.77
1:H:16:GLU:HG2	1:H:16:GLU:O	1.86	0.76
1:E:83:GLU:CD	1:E:86:ARG:NH1	2.39	0.75
1:E:31:GLY:N	1:E:134:ASP:OD1	2.17	0.75
1:B:19:PRO:HB3	1:B:78:ALA:HB3	1.72	0.71
1:E:161:LEU:OXT	3:E:2127:HOH:O	2.10	0.69
1:G:29:PHE:HB3	3:G:2032:HOH:O	1.92	0.68
1:D:18:GLU:HG2	1:D:21:ASN:ND2	2.07	0.68
1:B:22:LYS:HD2	3:B:2019:HOH:O	1.94	0.68
1:C:29:PHE:HB3	3:C:2052:HOH:O	1.92	0.68
1:B:57:GLU:HB3	3:B:2046:HOH:O	1.94	0.67
1:H:16:GLU:CD	1:H:95:ARG:HH21	1.98	0.66
1:H:19:PRO:HA	3:H:2012:HOH:O	1.98	0.63
1:D:86:ARG:NH2	3:D:2109:HOH:O	2.31	0.62
1:B:144:PRO:O	1:C:31:GLY:HA2	1.98	0.62
1:B:69:VAL:HG23	1:B:69:VAL:O	1.98	0.62
1:H:18:GLU:HA	1:H:18:GLU:OE1	2.00	0.62
1:H:89:LYS:HE3	3:H:2009:HOH:O	2.00	0.61
1:H:57:GLU:HG3	3:H:2044:HOH:O	1.99	0.61
1:H:17:GLY:O	1:H:19:PRO:N	2.34	0.60
1:H:24:ASN:ND2	3:H:2019:HOH:O	2.34	0.59
1:H:20:GLY:N	3:H:2012:HOH:O	2.35	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:32:LYS:HE3	3:B:2056:HOH:O	2.05	0.57
1:G:161:LEU:O	3:G:2138:HOH:O	2.19	0.53
1:G:126:LYS:NZ	3:G:2116:HOH:O	2.43	0.52
1:C:161:LEU:O	3:C:2187:HOH:O	2.19	0.52
1:H:91:GLU:HB2	3:H:2075:HOH:O	2.09	0.51
1:G:9:ILE:HD11	1:G:130:MET:CE	2.38	0.51
1:G:93:LYS:HE3	3:G:2056:HOH:O	2.10	0.50
1:G:27:GLU:HG2	3:G:2008:HOH:O	2.11	0.50
1:E:136:ILE:CD1	1:E:138:LYS:HE3	2.22	0.50
1:D:54:GLY:HA2	1:D:57:GLU:HG2	1.94	0.50
1:H:123:ARG:NH1	3:H:2106:HOH:O	2.44	0.49
1:G:134:ASP:O	1:G:134:ASP:OD1	2.30	0.49
1:A:29:PHE:HB3	3:A:2053:HOH:O	2.11	0.49
1:E:16:GLU:OE1	1:E:95:ARG:NH1	2.43	0.49
1:F:62:LEU:HD23	1:F:158:ILE:HD13	1.94	0.48
1:H:16:GLU:OE1	1:H:95:ARG:NH2	2.44	0.48
1:A:4:LYS:HE2	3:A:2006:HOH:O	2.13	0.48
1:B:107:GLU:HG2	3:B:2090:HOH:O	2.13	0.48
1:H:89:LYS:HE2	3:H:2027:HOH:O	2.13	0.48
1:B:134:ASP:O	1:B:134:ASP:CG	2.50	0.48
1:B:19:PRO:HG2	1:B:79:PHE:CE1	2.48	0.47
1:H:83:GLU:OE2	1:H:86:ARG:NH1	2.41	0.47
1:H:26:ALA:O	3:H:2021:HOH:O	2.20	0.47
1:H:15:PHE:CE2	1:H:20:GLY:HA2	2.50	0.47
1:E:52:LEU:HB3	1:E:53:PRO:HD3	1.97	0.47
1:D:18:GLU:HG2	1:D:21:ASN:HD21	1.80	0.46
1:H:83:GLU:CG	1:H:86:ARG:HH12	2.27	0.46
1:B:39:VAL:HB	1:B:47:CYS:SG	2.56	0.45
1:H:16:GLU:OE2	1:H:95:ARG:NE	2.42	0.45
1:H:39:VAL:HB	1:H:47:CYS:SG	2.56	0.45
1:D:39:VAL:HB	1:D:47:CYS:SG	2.55	0.45
1:G:95:ARG:HD3	1:G:95:ARG:HH11	1.49	0.45
1:F:62:LEU:HD23	1:F:158:ILE:CD1	2.47	0.44
1:H:86:ARG:NH2	3:H:2068:HOH:O	2.33	0.44
1:F:52:LEU:HB3	1:F:53:PRO:HD3	2.00	0.44
1:E:69:VAL:HG23	3:E:2059:HOH:O	2.17	0.44
1:H:17:GLY:O	1:H:18:GLU:C	2.56	0.44
1:C:39:VAL:HB	1:C:47:CYS:SG	2.58	0.44
1:E:18:GLU:HG2	3:E:2016:HOH:O	2.17	0.44
1:E:16:GLU:OE2	1:E:95:ARG:HD2	2.17	0.44
1:E:40:PRO:HD2	1:E:47:CYS:SG	2.58	0.44
1:E:13:GLU:OE2	1:E:22:LYS:HD3	2.18	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:H:106:LYS:NZ	3:H:2091:HOH:O	2.51	0.43
1:B:18:GLU:HG2	3:B:2013:HOH:O	2.17	0.43
1:E:136:ILE:HD13	1:E:136:ILE:C	2.39	0.43
1:A:37:PHE:CZ	1:A:72:CYS:HB2	2.53	0.43
1:G:37:PHE:CZ	1:G:72:CYS:HB2	2.54	0.43
1:H:136:ILE:CG1	1:H:138:LYS:HE2	2.48	0.43
1:H:32:LYS:HE3	3:H:2023:HOH:O	2.18	0.43
1:F:83:GLU:OE2	1:F:86:ARG:NH1	2.51	0.43
1:G:56:VAL:HG13	1:G:93:LYS:HD3	2.01	0.43
1:B:69:VAL:CG2	1:B:69:VAL:O	2.66	0.43
1:E:136:ILE:HD13	1:E:136:ILE:O	2.19	0.43
1:E:69:VAL:HA	3:E:2059:HOH:O	2.18	0.42
1:H:9:ILE:HG23	1:H:10:PRO:HD2	1.99	0.42
1:D:40:PRO:HD2	1:D:47:CYS:SG	2.59	0.42
1:H:15:PHE:HE2	1:H:20:GLY:HA2	1.85	0.42
1:A:52:LEU:N	1:A:53:PRO:HD2	2.35	0.42
1:H:16:GLU:OE2	1:H:95:ARG:NH2	2.53	0.41
1:G:39:VAL:HB	1:G:47:CYS:SG	2.60	0.41
1:G:126:LYS:CE	3:G:2116:HOH:O	2.68	0.41
1:G:46:GLY:H	2:G:1162:BEZ:C	2.33	0.41
1:E:38:GLY:HA3	1:E:128:PHE:CE2	2.56	0.41
1:H:19:PRO:HG2	1:H:79:PHE:CD1	2.55	0.41
1:A:33:LYS:NZ	1:A:65:LYS:O	2.46	0.41
1:B:40:PRO:HD2	1:B:47:CYS:SG	2.60	0.41
1:C:101:THR:O	1:C:106:LYS:HE3	2.21	0.41
1:H:38:GLY:HA3	1:H:128:PHE:CE1	2.55	0.41
1:G:52:LEU:HB3	1:G:53:PRO:HD3	2.03	0.41
1:H:157:ILE:HG12	3:H:2120:HOH:O	2.20	0.41
1:G:118:SER:OG	1:H:60:GLU:OE1	2.31	0.41
1:H:19:PRO:CA	3:H:2012:HOH:O	2.64	0.41
1:F:39:VAL:HB	1:F:47:CYS:SG	2.61	0.40
1:G:161:LEU:HA	1:G:161:LEU:HD23	1.61	0.40
1:G:130:MET:HE2	1:G:130:MET:HB2	1.70	0.40
1:A:46:GLY:H	2:A:1162:BEZ:C	2.34	0.40
1:G:128:PHE:C	1:G:128:PHE:CD1	2.95	0.40
1:H:25:LEU:HA	1:H:25:LEU:HD23	1.95	0.40
1:F:62:LEU:HG	1:F:158:ILE:HD11	2.04	0.40
1:H:9:ILE:HA	1:H:10:PRO:HD3	1.94	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	159/161 (99%)	153 (96%)	6 (4%)	0	100	100
1	B	159/161 (99%)	148 (93%)	9 (6%)	2 (1%)	18	5
1	C	159/161 (99%)	154 (97%)	5 (3%)	0	100	100
1	D	159/161 (99%)	153 (96%)	6 (4%)	0	100	100
1	E	159/161 (99%)	153 (96%)	6 (4%)	0	100	100
1	F	159/161 (99%)	154 (97%)	5 (3%)	0	100	100
1	G	159/161 (99%)	155 (98%)	4 (2%)	0	100	100
1	H	159/161 (99%)	153 (96%)	6 (4%)	0	100	100
All	All	1272/1288 (99%)	1223 (96%)	47 (4%)	2 (0%)	56	46

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	18	GLU
1	B	19	PRO

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	126/126 (100%)	124 (98%)	2 (2%)	75	69
1	B	126/126 (100%)	124 (98%)	2 (2%)	75	69
1	C	126/126 (100%)	123 (98%)	3 (2%)	61	51
1	D	126/126 (100%)	122 (97%)	4 (3%)	51	37
1	E	126/126 (100%)	124 (98%)	2 (2%)	75	69

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	F	126/126 (100%)	124 (98%)	2 (2%)	75	69
1	G	126/126 (100%)	124 (98%)	2 (2%)	75	69
1	H	126/126 (100%)	119 (94%)	7 (6%)	30	13
All	All	1008/1008 (100%)	984 (98%)	24 (2%)	61	51

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	157	ILE
1	A	158	ILE
1	B	73	LEU
1	B	157	ILE
1	C	18	GLU
1	C	144	PRO
1	C	157	ILE
1	D	9	ILE
1	D	91	GLU
1	D	114	ASP
1	D	157	ILE
1	E	13	GLU
1	E	136	ILE
1	F	9	ILE
1	F	157	ILE
1	G	95	ARG
1	G	107	GLU
1	H	9	ILE
1	H	10	PRO
1	H	18	GLU
1	H	21	ASN
1	H	23	VAL
1	H	136	ILE
1	H	159	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	21	ASN



### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

8 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z  > 2$	Counts	RMSZ	# $ Z  > 2$
2	BEZ	A	1162	-	9,9,9	2.29	2 (22%)	11,11,11	2.18	6 (54%)
2	BEZ	B	1162	-	9,9,9	1.88	1 (11%)	11,11,11	1.03	0
2	BEZ	C	1162	-	9,9,9	1.75	1 (11%)	11,11,11	1.22	1 (9%)
2	BEZ	D	1162	-	9,9,9	1.96	2 (22%)	11,11,11	1.20	1 (9%)
2	BEZ	E	1162	-	9,9,9	1.91	1 (11%)	11,11,11	0.89	0
2	BEZ	F	1162	-	9,9,9	1.87	1 (11%)	11,11,11	0.92	0
2	BEZ	G	1162	-	9,9,9	2.17	2 (22%)	11,11,11	1.64	3 (27%)
2	BEZ	H	1162	-	9,9,9	1.72	1 (11%)	11,11,11	1.14	1 (9%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	BEZ	A	1162	-	-	0/4/4/4	0/1/1/1
2	BEZ	B	1162	-	-	0/4/4/4	0/1/1/1
2	BEZ	C	1162	-	-	0/4/4/4	0/1/1/1
2	BEZ	D	1162	-	-	0/4/4/4	0/1/1/1
2	BEZ	E	1162	-	-	0/4/4/4	0/1/1/1
2	BEZ	F	1162	-	-	0/4/4/4	0/1/1/1
2	BEZ	G	1162	-	-	0/4/4/4	0/1/1/1
2	BEZ	H	1162	-	-	0/4/4/4	0/1/1/1

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1162	BEZ	O1-C	6.10	1.41	1.23
2	G	1162	BEZ	O1-C	5.99	1.41	1.23
2	E	1162	BEZ	O1-C	5.62	1.40	1.23
2	B	1162	BEZ	O1-C	5.41	1.39	1.23
2	D	1162	BEZ	O1-C	5.41	1.39	1.23
2	F	1162	BEZ	O1-C	5.38	1.39	1.23
2	H	1162	BEZ	O1-C	5.04	1.38	1.23
2	C	1162	BEZ	O1-C	5.00	1.38	1.23
2	A	1162	BEZ	C1-C	-2.52	1.44	1.49
2	G	1162	BEZ	C1-C	-2.02	1.45	1.49
2	D	1162	BEZ	C1-C	-2.01	1.45	1.49

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1162	BEZ	C6-C1-C2	3.96	124.30	118.63
2	G	1162	BEZ	C6-C1-C2	3.03	122.96	118.63
2	A	1162	BEZ	C5-C6-C1	-2.88	116.57	120.32
2	A	1162	BEZ	C3-C2-C1	-2.59	116.95	120.32
2	G	1162	BEZ	C5-C6-C1	-2.48	117.11	120.32
2	A	1162	BEZ	O2-C-O1	2.37	128.72	123.35
2	D	1162	BEZ	C6-C1-C2	2.32	121.94	118.63
2	C	1162	BEZ	C6-C1-C2	2.26	121.86	118.63
2	A	1162	BEZ	O1-C-C1	-2.24	115.25	121.25
2	H	1162	BEZ	C6-C1-C2	2.21	121.79	118.63
2	G	1162	BEZ	C3-C2-C1	-2.09	117.60	120.32
2	A	1162	BEZ	C2-C1-C	-2.08	116.62	120.44

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	161/161 (100%)	-0.28	1 (0%) 86 92	10, 15, 22, 27	0
1	B	161/161 (100%)	-0.11	3 (1%) 64 72	10, 15, 24, 32	0
1	C	161/161 (100%)	-0.14	0 100 100	10, 15, 23, 27	0
1	D	161/161 (100%)	-0.21	0 100 100	10, 15, 22, 27	0
1	E	161/161 (100%)	-0.08	3 (1%) 64 72	10, 15, 22, 27	0
1	F	161/161 (100%)	-0.30	0 100 100	10, 15, 22, 27	0
1	G	161/161 (100%)	-0.36	1 (0%) 86 92	10, 15, 22, 27	0
1	H	161/161 (100%)	-0.16	1 (0%) 86 92	10, 15, 23, 30	0
All	All	1288/1288 (100%)	-0.21	9 (0%) 84 91	10, 15, 23, 32	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	17	GLY	3.9
1	A	161	LEU	3.2
1	E	31	GLY	3.2
1	E	1	ALA	2.9
1	B	1	ALA	2.6
1	B	18	GLU	2.4
1	G	1	ALA	2.2
1	E	30	LYS	2.2
1	H	17	GLY	2.1

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

### 6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	BEZ	G	1162	9/9	0.15	4.34	18,23,25,25	0
2	BEZ	A	1162	9/9	0.13	3.33	18,22,25,26	0
2	BEZ	F	1162	9/9	0.11	1.80	18,23,25,25	0
2	BEZ	B	1162	9/9	0.10	1.45	18,23,24,25	0
2	BEZ	E	1162	9/9	0.10	1.28	18,23,25,25	0
2	BEZ	D	1162	9/9	0.10	0.85	18,23,25,25	0
2	BEZ	C	1162	9/9	0.11	0.06	18,23,24,25	0
2	BEZ	H	1162	9/9	0.09	-0.13	18,23,24,25	0

### 6.5 Other polymers ⓘ

There are no such residues in this entry.