



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:15 PM UTC

PDB ID : 7TPB / pdb\_00007tpb  
Title : p120RasGAP SH3 domain in complex with DLC1 RhoGAP domain  
Authors : Stiegler, A.L.; Boggon, T.J.; Chau, J.E.; Vish, K.J.  
Deposited on : 2022-01-25  
Resolution : 3.20 Å(reported)

This is a Full wwPDB X-ray Structure 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0  
Xtrriage (Phenix) : 2.0  
EDS : 3.0  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
CCP4 : 9.0.010 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

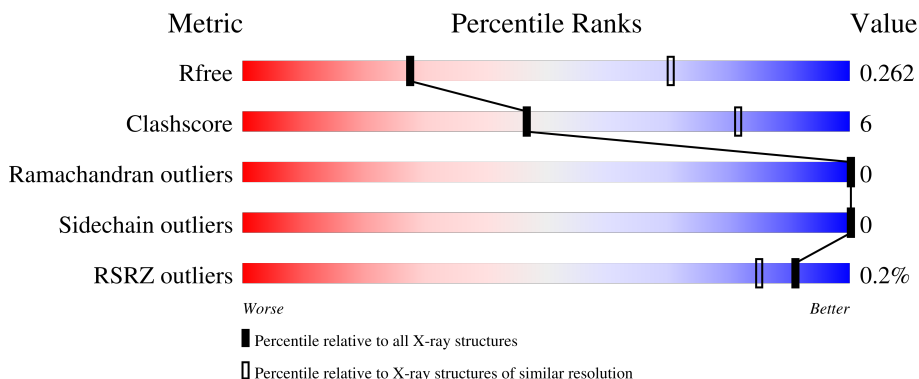
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.20 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))
$R_{free}$	180053	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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 of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	66	 2% 62% 24% 12%
1	C	66	 59% 30% 11%
1	E	66	 65% 24% 11%
1	G	66	 2% 65% 24% 11%
2	B	228	 73% 12% 15%

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Mol	Chain	Length	Quality of chain
2	D	228	 78% 8% 14%
2	F	228	 72% 13% 15%
2	H	228	 76% 9% 15%

## 2 Entry composition i

There are 2 unique types of molecules in this entry. The entry contains 8096 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Ras GTPase-activating protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	58	480	306	79	93	2	0	0	0
1	C	59	491	312	83	94	2	0	0	0
1	E	59	491	312	83	94	2	0	0	0
1	G	59	491	312	83	94	2	0	0	0

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	276	GLY	-	expression tag	UNP B4DTL8
A	277	PRO	-	expression tag	UNP B4DTL8
A	278	LEU	-	expression tag	UNP B4DTL8
A	279	GLY	-	expression tag	UNP B4DTL8
A	280	SER	-	expression tag	UNP B4DTL8
C	276	GLY	-	expression tag	UNP B4DTL8
C	277	PRO	-	expression tag	UNP B4DTL8
C	278	LEU	-	expression tag	UNP B4DTL8
C	279	GLY	-	expression tag	UNP B4DTL8
C	280	SER	-	expression tag	UNP B4DTL8
E	276	GLY	-	expression tag	UNP B4DTL8
E	277	PRO	-	expression tag	UNP B4DTL8
E	278	LEU	-	expression tag	UNP B4DTL8
E	279	GLY	-	expression tag	UNP B4DTL8
E	280	SER	-	expression tag	UNP B4DTL8
G	276	GLY	-	expression tag	UNP B4DTL8
G	277	PRO	-	expression tag	UNP B4DTL8
G	278	LEU	-	expression tag	UNP B4DTL8
G	279	GLY	-	expression tag	UNP B4DTL8
G	280	SER	-	expression tag	UNP B4DTL8

- Molecule 2 is a protein called Rho GTPase-activating protein 7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	194	Total 1538	C 973	N 266	O 288	S 11	0	0	0
2	D	195	Total 1547	C 979	N 268	O 289	S 11	0	0	0
2	F	193	Total 1529	C 967	N 264	O 287	S 11	0	0	0
2	H	193	Total 1529	C 967	N 264	O 287	S 11	0	0	0

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1056	MET	-	initiating methionine	UNP Q96QB1-6
B	1057	HIS	-	expression tag	UNP Q96QB1-6
B	1058	HIS	-	expression tag	UNP Q96QB1-6
B	1059	HIS	-	expression tag	UNP Q96QB1-6
B	1060	HIS	-	expression tag	UNP Q96QB1-6
B	1061	HIS	-	expression tag	UNP Q96QB1-6
B	1062	HIS	-	expression tag	UNP Q96QB1-6
B	1063	SER	-	expression tag	UNP Q96QB1-6
B	1064	SER	-	expression tag	UNP Q96QB1-6
B	1065	GLY	-	expression tag	UNP Q96QB1-6
B	1066	ARG	-	expression tag	UNP Q96QB1-6
B	1067	GLU	-	expression tag	UNP Q96QB1-6
B	1068	ASN	-	expression tag	UNP Q96QB1-6
B	1069	LEU	-	expression tag	UNP Q96QB1-6
B	1070	TYR	-	expression tag	UNP Q96QB1-6
B	1071	PHE	-	expression tag	UNP Q96QB1-6
B	1072	GLN	-	expression tag	UNP Q96QB1-6
B	1073	GLY	-	expression tag	UNP Q96QB1-6
D	1056	MET	-	initiating methionine	UNP Q96QB1-6
D	1057	HIS	-	expression tag	UNP Q96QB1-6
D	1058	HIS	-	expression tag	UNP Q96QB1-6
D	1059	HIS	-	expression tag	UNP Q96QB1-6
D	1060	HIS	-	expression tag	UNP Q96QB1-6
D	1061	HIS	-	expression tag	UNP Q96QB1-6
D	1062	HIS	-	expression tag	UNP Q96QB1-6
D	1063	SER	-	expression tag	UNP Q96QB1-6
D	1064	SER	-	expression tag	UNP Q96QB1-6
D	1065	GLY	-	expression tag	UNP Q96QB1-6
D	1066	ARG	-	expression tag	UNP Q96QB1-6
D	1067	GLU	-	expression tag	UNP Q96QB1-6

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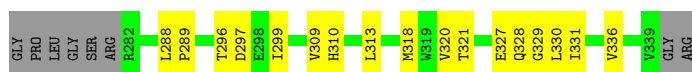
Chain	Residue	Modelled	Actual	Comment	Reference
D	1068	ASN	-	expression tag	UNP Q96QB1-6
D	1069	LEU	-	expression tag	UNP Q96QB1-6
D	1070	TYR	-	expression tag	UNP Q96QB1-6
D	1071	PHE	-	expression tag	UNP Q96QB1-6
D	1072	GLN	-	expression tag	UNP Q96QB1-6
D	1073	GLY	-	expression tag	UNP Q96QB1-6
F	1056	MET	-	initiating methionine	UNP Q96QB1-6
F	1057	HIS	-	expression tag	UNP Q96QB1-6
F	1058	HIS	-	expression tag	UNP Q96QB1-6
F	1059	HIS	-	expression tag	UNP Q96QB1-6
F	1060	HIS	-	expression tag	UNP Q96QB1-6
F	1061	HIS	-	expression tag	UNP Q96QB1-6
F	1062	HIS	-	expression tag	UNP Q96QB1-6
F	1063	SER	-	expression tag	UNP Q96QB1-6
F	1064	SER	-	expression tag	UNP Q96QB1-6
F	1065	GLY	-	expression tag	UNP Q96QB1-6
F	1066	ARG	-	expression tag	UNP Q96QB1-6
F	1067	GLU	-	expression tag	UNP Q96QB1-6
F	1068	ASN	-	expression tag	UNP Q96QB1-6
F	1069	LEU	-	expression tag	UNP Q96QB1-6
F	1070	TYR	-	expression tag	UNP Q96QB1-6
F	1071	PHE	-	expression tag	UNP Q96QB1-6
F	1072	GLN	-	expression tag	UNP Q96QB1-6
F	1073	GLY	-	expression tag	UNP Q96QB1-6
H	1056	MET	-	initiating methionine	UNP Q96QB1-6
H	1057	HIS	-	expression tag	UNP Q96QB1-6
H	1058	HIS	-	expression tag	UNP Q96QB1-6
H	1059	HIS	-	expression tag	UNP Q96QB1-6
H	1060	HIS	-	expression tag	UNP Q96QB1-6
H	1061	HIS	-	expression tag	UNP Q96QB1-6
H	1062	HIS	-	expression tag	UNP Q96QB1-6
H	1063	SER	-	expression tag	UNP Q96QB1-6
H	1064	SER	-	expression tag	UNP Q96QB1-6
H	1065	GLY	-	expression tag	UNP Q96QB1-6
H	1066	ARG	-	expression tag	UNP Q96QB1-6
H	1067	GLU	-	expression tag	UNP Q96QB1-6
H	1068	ASN	-	expression tag	UNP Q96QB1-6
H	1069	LEU	-	expression tag	UNP Q96QB1-6
H	1070	TYR	-	expression tag	UNP Q96QB1-6
H	1071	PHE	-	expression tag	UNP Q96QB1-6
H	1072	GLN	-	expression tag	UNP Q96QB1-6
H	1073	GLY	-	expression tag	UNP Q96QB1-6

### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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: Ras GTPase-activating protein 1

Chain A: 



- Molecule 1: Ras GTPase-activating protein 1

Chain C: 



- Molecule 1: Ras GTPase-activating protein 1

Chain E: 



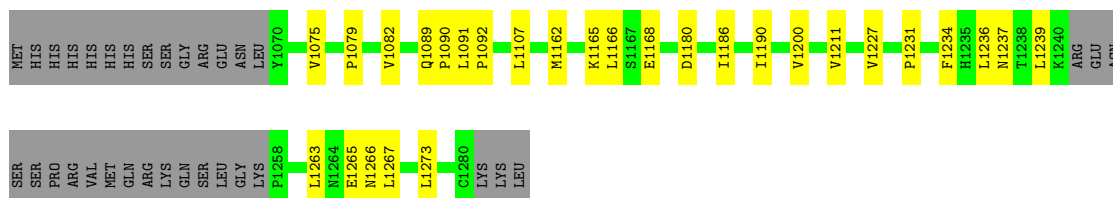
- Molecule 1: Ras GTPase-activating protein 1

Chain G: 




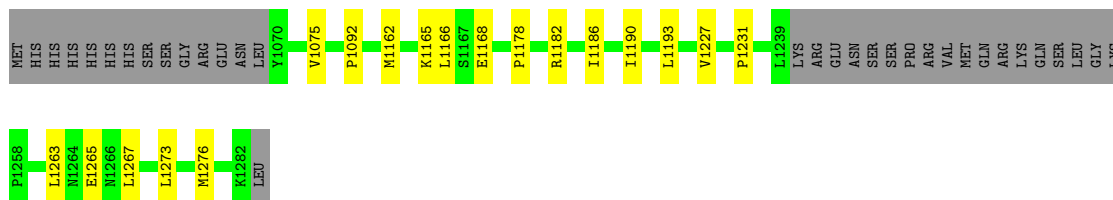
- Molecule 2: Rho GTPase-activating protein 7

Chain B: 



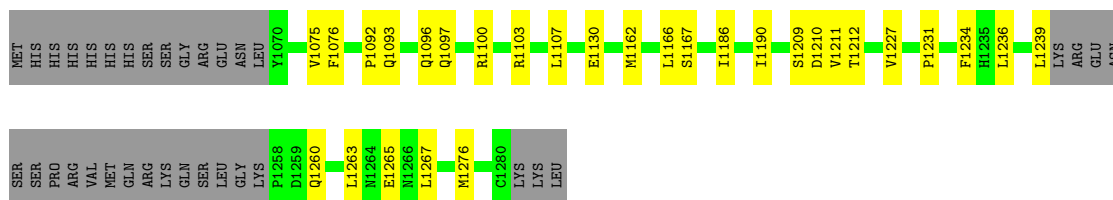
- Molecule 2: Rho GTPase-activating protein 7

Chain D:  78% 8% 14%



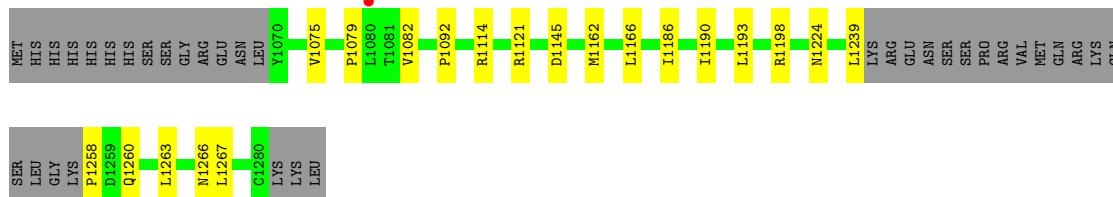
- Molecule 2: Rho GTPase-activating protein 7

Chain F:  72% 13% 15%



- Molecule 2: Rho GTPase-activating protein 7

Chain H:  76% 9% 15%



## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	143.75Å 143.75Å 152.78Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.25 – 3.20 48.25 – 3.20	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.25-3.20) 99.9 (48.25-3.20)	Depositor EDS
$R_{merge}$	0.26	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.81 (at 3.19Å)	Xtrriage
Refinement program	PHENIX 1.19.2	Depositor
R, $R_{free}$	0.212 , 0.260 0.213 , 0.262	Depositor DCC
$R_{free}$ test set	920 reflections (4.74%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	80.4	Xtrriage
Anisotropy	0.031	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 49.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	0.064 for h,-h-k,-l	Xtrriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	8096	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	85.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.60% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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	0.07	0/489	0.26	0/664
1	C	0.08	0/500	0.26	0/678
1	E	0.07	0/500	0.25	0/678
1	G	0.09	0/500	0.27	0/678
2	B	0.10	0/1564	0.27	0/2117
2	D	0.10	0/1573	0.26	0/2128
2	F	0.10	0/1555	0.27	0/2106
2	H	0.09	0/1555	0.25	0/2106
All	All	0.09	0/8236	0.26	0/11155

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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 hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	480	0	469	12	0
1	C	491	0	482	14	0
1	E	491	0	482	13	0
1	G	491	0	482	11	0
2	B	1538	0	1548	19	0
2	D	1547	0	1561	13	0
2	F	1529	0	1535	20	0
2	H	1529	0	1535	12	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	8096	0	8094	92	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (92) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:327:GLU:HB2	2:B:1239:LEU:HD11	1.63	0.79
2:B:1162:MET:HB3	2:B:1166:LEU:HD23	1.69	0.73
2:F:1103:ARG:NH1	2:F:1210:ASP:OD2	2.20	0.72
1:G:288:LEU:HD12	1:G:289:PRO:HD2	1.71	0.70
1:A:309:VAL:HA	1:A:320:VAL:HG12	1.73	0.70
2:D:1162:MET:HB3	2:D:1166:LEU:HD23	1.79	0.65
2:F:1162:MET:HB3	2:F:1166:LEU:HD23	1.79	0.65
2:H:1162:MET:HB3	2:H:1166:LEU:HD23	1.81	0.63
1:C:310:HIS:NE2	1:C:321:THR:OG1	2.26	0.62
2:F:1075:VAL:HG13	2:F:1092:PRO:HG2	1.82	0.61
1:C:313:LEU:HD21	2:D:1267:LEU:HD11	1.83	0.61
2:H:1239:LEU:HB3	2:H:1258:PRO:HD2	1.83	0.59
1:G:288:LEU:HD23	1:G:335:LEU:HD22	1.85	0.59
2:B:1180:ASP:HB2	2:F:1093:GLN:HG2	1.85	0.58
1:G:328:GLN:HB2	2:H:1263:LEU:HD13	1.87	0.57
2:F:1167:SER:HB2	2:F:1265:GLU:OE2	2.04	0.56
1:G:313:LEU:HD21	2:H:1267:LEU:HD11	1.87	0.56
2:H:1075:VAL:HG13	2:H:1092:PRO:HG2	1.89	0.55
2:D:1162:MET:HE1	2:D:1276:MET:HE1	1.89	0.54
1:E:309:VAL:HA	1:E:320:VAL:HG12	1.88	0.54
2:F:1162:MET:HE1	2:F:1276:MET:HE1	1.88	0.53
1:E:328:GLN:HB2	2:F:1263:LEU:HD13	1.89	0.53
2:B:1168:GLU:HG2	2:B:1265:GLU:OE2	2.09	0.52
1:C:330:LEU:HD13	2:D:1227:VAL:HG22	1.91	0.52
1:G:309:VAL:HA	1:G:320:VAL:HG12	1.91	0.52
1:E:331:ILE:HD12	1:E:336:VAL:HG11	1.91	0.51
1:G:331:ILE:HD12	1:G:336:VAL:HG11	1.92	0.51
1:E:292:LYS:HG3	1:E:298:GLU:HB2	1.93	0.51
2:F:1097:GLN:HG2	2:F:1100:ARG:HH12	1.77	0.50
1:G:310:HIS:NE2	1:G:321:THR:OG1	2.29	0.50
1:E:328:GLN:NE2	2:F:1260:GLN:OE1	2.45	0.50
2:B:1075:VAL:HG13	2:B:1092:PRO:HG2	1.93	0.49
1:C:296:THR:HA	2:D:1231:PRO:HG2	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:331:ILE:HD12	1:A:336:VAL:HG11	1.95	0.49
2:B:1089:GLN:HG2	2:B:1090:PRO:HD2	1.95	0.48
1:A:318:MET:HE3	1:A:318:MET:HB2	1.76	0.48
1:A:328:GLN:HB2	2:B:1263:LEU:HD13	1.96	0.48
2:F:1096:GLN:O	2:F:1100:ARG:HG3	2.14	0.47
2:B:1091:LEU:HG	2:B:1200:VAL:HG22	1.97	0.47
2:B:1186:ILE:O	2:B:1190:ILE:HG12	2.14	0.47
2:B:1237:ASN:O	2:B:1266:ASN:ND2	2.48	0.47
1:G:292:LYS:HG3	1:G:298:GLU:HB2	1.97	0.47
1:E:296:THR:HA	2:F:1231:PRO:HG2	1.95	0.47
1:C:331:ILE:HD12	1:C:336:VAL:HG11	1.96	0.47
1:G:297:ASP:OD2	2:H:1266:ASN:HB3	2.15	0.46
1:C:284:VAL:HB	1:C:336:VAL:HB	1.96	0.46
1:A:288:LEU:HD12	1:A:289:PRO:HD2	1.98	0.46
1:C:318:MET:HE1	1:C:336:VAL:HG21	1.98	0.46
2:D:1273:LEU:HA	2:D:1276:MET:HE3	1.98	0.45
1:A:310:HIS:NE2	1:A:321:THR:OG1	2.34	0.45
1:A:299:ILE:HG23	1:A:329:GLY:HA3	1.97	0.45
1:C:286:ALA:HB2	1:C:307:PHE:HE1	1.81	0.45
1:E:288:LEU:HD12	1:E:289:PRO:HD2	1.99	0.45
1:E:330:LEU:HD13	2:F:1227:VAL:HG22	1.99	0.45
1:C:309:VAL:HA	1:C:320:VAL:HG12	1.98	0.45
1:E:318:MET:HE1	1:E:336:VAL:HG21	1.98	0.45
1:E:319:TRP:HZ2	2:F:1263:LEU:HB3	1.81	0.45
2:F:1234:PHE:HB2	2:F:1236:LEU:HG	1.98	0.45
1:A:330:LEU:HD13	2:B:1227:VAL:HG22	1.98	0.44
1:E:313:LEU:HD21	2:F:1267:LEU:HD11	1.99	0.44
2:H:1121:ARG:NE	2:H:1145:ASP:OD2	2.36	0.44
2:D:1168:GLU:HG3	2:D:1265:GLU:OE2	2.17	0.44
1:E:299:ILE:HG23	1:E:329:GLY:HA3	1.98	0.44
1:C:299:ILE:HG23	1:C:329:GLY:HA3	2.00	0.44
2:F:1186:ILE:O	2:F:1190:ILE:HG12	2.18	0.44
1:C:292:LYS:HG3	1:C:298:GLU:HB2	2.00	0.43
1:G:328:GLN:NE2	2:H:1260:GLN:OE1	2.49	0.43
2:F:1107:LEU:HD22	2:F:1211:VAL:HA	1.99	0.43
2:B:1107:LEU:HD22	2:B:1211:VAL:HA	2.00	0.43
2:B:1079:PRO:HG2	2:B:1082:VAL:HG23	2.01	0.43
2:H:1186:ILE:O	2:H:1190:ILE:HG12	2.19	0.43
2:B:1162:MET:HE1	2:B:1273:LEU:HD13	2.01	0.43
2:D:1178:PRO:O	2:D:1182:ARG:HG3	2.20	0.42
1:A:297:ASP:OD2	2:B:1266:ASN:HB3	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:328:GLN:HB2	2:D:1263:LEU:HD13	2.00	0.42
1:C:288:LEU:HD12	1:C:289:PRO:HD2	2.02	0.42
1:G:299:ILE:HG23	1:G:329:GLY:HA3	2.01	0.42
2:B:1234:PHE:HB2	2:B:1236:LEU:HG	2.01	0.42
2:F:1076:PHE:N	2:F:1130:GLU:OE2	2.47	0.42
2:D:1190:ILE:HA	2:D:1193:LEU:HD13	2.01	0.42
1:A:296:THR:HA	2:B:1231:PRO:HG2	2.01	0.42
2:D:1075:VAL:HG13	2:D:1092:PRO:HG2	2.01	0.41
2:H:1079:PRO:HG2	2:H:1082:VAL:HG23	2.02	0.41
2:H:1114:ARG:NH2	2:H:1224:ASN:OD1	2.39	0.41
2:B:1165:LYS:H	2:B:1165:LYS:HD2	1.85	0.41
2:D:1165:LYS:HD2	2:D:1165:LYS:H	1.86	0.41
1:A:313:LEU:HD21	2:B:1267:LEU:HD11	2.02	0.41
2:F:1209:SER:HA	2:F:1212:THR:HG22	2.03	0.41
2:H:1193:LEU:O	2:H:1198:ARG:NH1	2.51	0.41
1:E:299:ILE:HG22	2:F:1239:LEU:HD21	2.03	0.40
2:D:1186:ILE:O	2:D:1190:ILE:HG12	2.21	0.40
1:C:286:ALA:HB2	1:C:307:PHE:CE1	2.56	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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	56/66 (85%)	56 (100%)	0	0	100	100
1	C	57/66 (86%)	57 (100%)	0	0	100	100
1	E	57/66 (86%)	57 (100%)	0	0	100	100
1	G	57/66 (86%)	57 (100%)	0	0	100	100
2	B	190/228 (83%)	188 (99%)	2 (1%)	0	100	100
2	D	191/228 (84%)	189 (99%)	2 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	F	189/228 (83%)	187 (99%)	2 (1%)	0	100	100
2	H	189/228 (83%)	187 (99%)	2 (1%)	0	100	100
All	All	986/1176 (84%)	978 (99%)	8 (1%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

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	54/59 (92%)	54 (100%)	0	100	100
1	C	55/59 (93%)	55 (100%)	0	100	100
1	E	55/59 (93%)	55 (100%)	0	100	100
1	G	55/59 (93%)	55 (100%)	0	100	100
2	B	170/202 (84%)	170 (100%)	0	100	100
2	D	171/202 (85%)	171 (100%)	0	100	100
2	F	169/202 (84%)	169 (100%)	0	100	100
2	H	169/202 (84%)	169 (100%)	0	100	100
All	All	898/1044 (86%)	898 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	311	ASN
2	B	1109	GLN
2	B	1175	GLN
1	C	311	ASN
2	D	1129	ASN
1	E	311	ASN
2	F	1175	GLN

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Mol	Chain	Res	Type
1	G	311	ASN
2	H	1129	ASN
2	H	1175	GLN
2	H	1181	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

### 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

### 5.7 Other polymers [i](#)

There are no such residues in this entry.

### 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	58/66 (87%)	-0.29	0 <a href="#">100</a> <a href="#">100</a>	65, 82, 112, 117	0
1	C	59/66 (89%)	-0.13	0 <a href="#">100</a> <a href="#">100</a>	67, 87, 132, 146	0
1	E	59/66 (89%)	-0.20	0 <a href="#">100</a> <a href="#">100</a>	81, 98, 125, 150	0
1	G	59/66 (89%)	-0.11	1 (1%) <a href="#">69</a> <a href="#">49</a>	78, 100, 142, 198	0
2	B	194/228 (85%)	-0.35	0 <a href="#">100</a> <a href="#">100</a>	52, 68, 106, 140	0
2	D	195/228 (85%)	-0.30	0 <a href="#">100</a> <a href="#">100</a>	58, 77, 109, 154	0
2	F	193/228 (84%)	-0.29	0 <a href="#">100</a> <a href="#">100</a>	52, 79, 123, 168	0
2	H	193/228 (84%)	-0.24	1 (0%) <a href="#">87</a> <a href="#">76</a>	62, 88, 121, 160	0
All	All	1010/1176 (85%)	-0.27	2 (0%) <a href="#">91</a> <a href="#">85</a>	52, 82, 121, 198	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	288	LEU	4.1
2	H	1080	LEU	2.1

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

### 6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

### 6.4 Ligands [i](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.