



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:00 AM UTC

PDB ID : 1JPL / pdb_00001jpl
Title : GGA3 VHS domain complexed with C-terminal peptide from cation-independent mannose 6-phosphate receptor
Authors : Misra, S.; Puertollano, R.; Bonifacino, J.S.; Hurley, J.H.
Deposited on : 2001-08-02
Resolution : 2.40 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

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>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (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

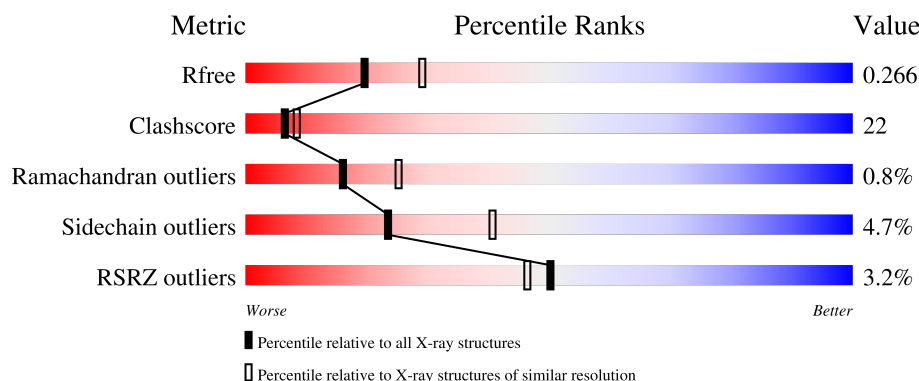
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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.



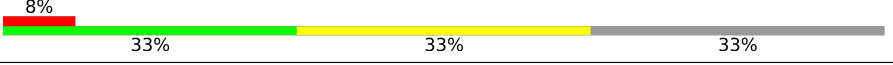
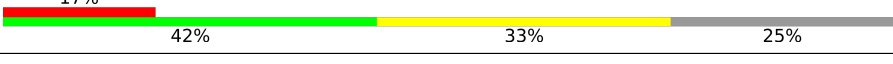
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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 ≥ 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	171	3% 64% 27% 6%
1	B	171	2% 54% 33% 5% 7%
1	C	171	0% 57% 35% 7%
1	D	171	2% 57% 34% 6%
2	E	12	17% 33% 33%

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Mol	Chain	Length	Quality of chain
2	F	12	
2	G	12	
2	H	12	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5655 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 ADP-RIBOSYLATION FACTOR BINDING PROTEIN GGA3.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	161	Total	C	N	O	S	Se	0	0	0
			1293	822	223	240	3	5			
1	B	159	Total	C	N	O	S	Se	0	0	0
			1280	813	221	238	3	5			
1	C	159	Total	C	N	O	S	Se	0	0	0
			1280	813	221	238	3	5			
1	D	161	Total	C	N	O	S	Se	0	0	0
			1293	822	223	240	3	5			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	201	GLY	-	cloning artifact	UNP Q9NZ52
A	202	ALA	-	cloning artifact	UNP Q9NZ52
A	203	MSE	-	cloning artifact	UNP Q9NZ52
A	204	GLY	-	cloning artifact	UNP Q9NZ52
A	205	SER	-	cloning artifact	UNP Q9NZ52
A	1	MSE	MET	modified residue	UNP Q9NZ52
A	73	MSE	MET	modified residue	UNP Q9NZ52
A	123	MSE	MET	modified residue	UNP Q9NZ52
A	137	MSE	MET	modified residue	UNP Q9NZ52
B	201	GLY	-	cloning artifact	UNP Q9NZ52
B	202	ALA	-	cloning artifact	UNP Q9NZ52
B	203	MSE	-	cloning artifact	UNP Q9NZ52
B	204	GLY	-	cloning artifact	UNP Q9NZ52
B	205	SER	-	cloning artifact	UNP Q9NZ52
B	1	MSE	MET	modified residue	UNP Q9NZ52
B	73	MSE	MET	modified residue	UNP Q9NZ52
B	123	MSE	MET	modified residue	UNP Q9NZ52
B	137	MSE	MET	modified residue	UNP Q9NZ52
C	201	GLY	-	cloning artifact	UNP Q9NZ52
C	202	ALA	-	cloning artifact	UNP Q9NZ52
C	203	MSE	-	cloning artifact	UNP Q9NZ52

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Chain	Residue	Modelled	Actual	Comment	Reference
C	204	GLY	-	cloning artifact	UNP Q9NZ52
C	205	SER	-	cloning artifact	UNP Q9NZ52
C	1	MSE	MET	modified residue	UNP Q9NZ52
C	73	MSE	MET	modified residue	UNP Q9NZ52
C	123	MSE	MET	modified residue	UNP Q9NZ52
C	137	MSE	MET	modified residue	UNP Q9NZ52
D	201	GLY	-	cloning artifact	UNP Q9NZ52
D	202	ALA	-	cloning artifact	UNP Q9NZ52
D	203	MSE	-	cloning artifact	UNP Q9NZ52
D	204	GLY	-	cloning artifact	UNP Q9NZ52
D	205	SER	-	cloning artifact	UNP Q9NZ52
D	1	MSE	MET	modified residue	UNP Q9NZ52
D	73	MSE	MET	modified residue	UNP Q9NZ52
D	123	MSE	MET	modified residue	UNP Q9NZ52
D	137	MSE	MET	modified residue	UNP Q9NZ52

- Molecule 2 is a protein called Cation-Independent Mannose 6-phosphate receptor.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	8	Total	C	N	O	0	0	0
			66	40	10	16			
2	F	8	Total	C	N	O	0	0	0
			66	40	10	16			
2	G	8	Total	C	N	O	0	0	0
			66	40	10	16			
2	H	9	Total	C	N	O	0	0	0
			74	44	11	19			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	53	Total	O	0	0
			53	53		
3	E	4	Total	O	0	0
			4	4		
3	B	60	Total	O	0	0
			60	60		
3	F	4	Total	O	0	0
			4	4		
3	C	66	Total	O	0	0
			66	66		
3	G	1	Total	O	0	0
			1	1		

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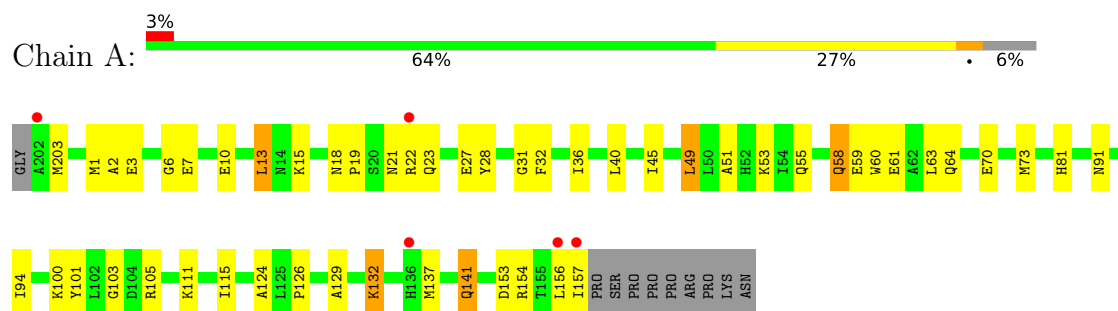
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	48	Total	O	0	0
			48	48		
3	H	1	Total	O	0	0
			1	1		

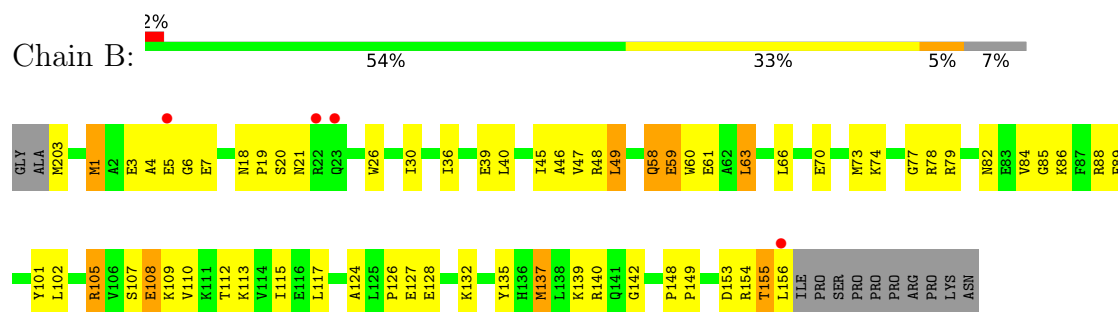
3 Residue-property plots [i](#)

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: ADP-RIBOSYLATION FACTOR BINDING PROTEIN GGA3



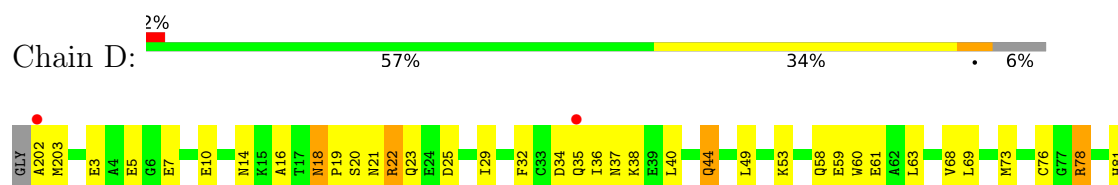
• Molecule 1: ADP-RIBOSYLATION FACTOR BINDING PROTEIN GGA3

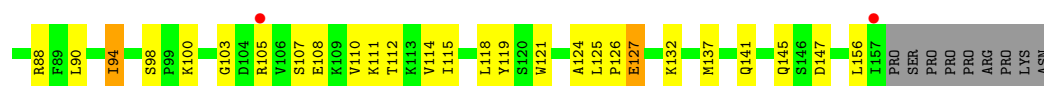


• Molecule 1: ADP-RIBOSYLATION FACTOR BINDING PROTEIN GGA3

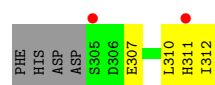


• Molecule 1: ADP-RIBOSYLATION FACTOR BINDING PROTEIN GGA3

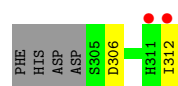




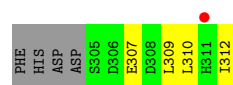
- Molecule 2: Cation-Independent Mannose 6-phosphate receptor



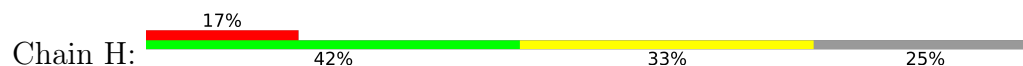
- Molecule 2: Cation-Independent Mannose 6-phosphate receptor



- Molecule 2: Cation-Independent Mannose 6-phosphate receptor



- Molecule 2: Cation-Independent Mannose 6-phosphate receptor



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	108.30Å 125.30Å 64.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.08 – 2.40 35.08 – 2.40	Depositor EDS
% Data completeness (in resolution range)	93.5 (35.08-2.40) 93.8 (35.08-2.40)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.20 (at 2.29Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.213 , 0.257 0.222 , 0.266	Depositor DCC
R_{free} test set	7402 reflections (9.86%)	wwPDB-VP
Wilson B-factor (Å ²)	30.7	Xtriage
Anisotropy	0.394	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 42.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5655	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 22.01 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 6.2297e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical 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	0.44	0/1314	0.90	2/1766 (0.1%)
1	B	0.42	0/1302	0.93	4/1751 (0.2%)
1	C	0.44	0/1302	0.93	3/1751 (0.2%)
1	D	0.41	0/1314	0.89	2/1766 (0.1%)
2	E	0.36	0/66	0.61	0/87
2	F	0.31	0/66	0.59	0/87
2	G	0.33	0/66	0.75	0/87
2	H	0.32	0/74	1.00	1/98 (1.0%)
All	All	0.42	0/5504	0.91	12/7393 (0.2%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	59	GLU	N-CA-C	7.98	120.69	111.11
1	D	105	ARG	N-CA-C	-7.47	104.01	113.72
1	A	59	GLU	N-CA-C	7.11	119.69	111.02
2	H	305	SER	N-CA-C	-6.95	104.54	113.16
1	B	59	GLU	N-CA-C	6.63	119.07	111.11
1	C	71	ALA	N-CA-C	-6.52	104.26	111.36
1	B	149	PRO	N-CA-C	-6.50	101.67	111.41
1	B	77	GLY	N-CA-C	6.49	122.74	111.78
1	C	59	GLU	N-CA-C	6.00	118.31	111.11
1	A	27	GLU	N-CA-C	-5.33	105.64	111.82
1	B	142	GLY	N-CA-C	5.29	122.40	115.47
1	C	122	THR	N-CA-C	-5.07	105.94	111.82

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-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 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	1293	0	1304	60	0
1	B	1280	0	1288	74	0
1	C	1280	0	1288	56	0
1	D	1293	0	1304	56	0
2	E	66	0	58	8	0
2	F	66	0	58	3	0
2	G	66	0	58	8	0
2	H	74	0	62	7	0
3	A	53	0	0	6	0
3	B	60	0	0	6	0
3	C	66	0	0	8	0
3	D	48	0	0	2	0
3	E	4	0	0	1	0
3	F	4	0	0	0	0
3	G	1	0	0	0	0
3	H	1	0	0	0	0
All	All	5655	0	5420	237	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 22.

All (237) 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:B:1:MSE:HE3	1:B:45:ILE:HG23	1.30	1.12
1:B:3:GLU:HG2	1:B:4:ALA:H	1.21	1.02
1:C:94:ILE:HD11	2:G:310:LEU:HD23	1.42	0.99
1:A:21:ASN:HD22	1:A:64:GLN:HE22	0.97	0.92
1:D:137:MSE:HE3	1:D:141:GLN:HG3	1.50	0.91
1:A:13:LEU:HD13	1:A:53:LYS:HG3	1.54	0.90
1:B:1:MSE:HE3	1:B:45:ILE:CG2	2.07	0.84
1:B:5:GLU:CD	1:B:39:GLU:HG2	2.03	0.83
1:A:1:MSE:HE2	1:B:58:GLN:HG2	1.63	0.80
1:A:1:MSE:CE	1:B:58:GLN:HG2	2.10	0.80
1:A:203:MSE:HE3	1:A:2:ALA:HB2	1.64	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:108:GLU:O	1:C:112:THR:HG23	1.81	0.79
1:A:58:GLN:NE2	3:A:206:HOH:O	2.15	0.79
1:A:28:TYR:HA	3:A:235:HOH:O	1.83	0.79
1:A:21:ASN:HD22	1:A:64:GLN:NE2	1.78	0.78
1:D:108:GLU:O	1:D:112:THR:HG23	1.84	0.77
1:B:105:ARG:HD2	1:B:105:ARG:O	1.86	0.76
1:B:154:ARG:C	1:B:156:LEU:H	1.94	0.76
1:A:40:LEU:HD22	1:C:126:PRO:HB3	1.70	0.74
1:C:53:LYS:HE3	3:C:206:HOH:O	1.87	0.73
1:D:78:ARG:HH22	1:D:125:LEU:HD22	1.52	0.73
1:A:58:GLN:CD	3:A:206:HOH:O	2.31	0.73
2:E:307:GLU:HB3	3:E:124:HOH:O	1.89	0.72
1:C:94:ILE:HD11	2:G:310:LEU:CD2	2.19	0.72
1:D:22:ARG:H	1:D:22:ARG:NE	1.87	0.71
1:B:1:MSE:HE2	1:B:6:GLY:HA3	1.70	0.71
1:B:135:TYR:CE2	1:B:139:LYS:HD2	2.25	0.71
1:B:3:GLU:HG2	1:B:4:ALA:N	1.98	0.70
1:D:3:GLU:O	1:D:7:GLU:HG3	1.91	0.70
1:B:155:THR:HG22	1:B:155:THR:O	1.92	0.70
1:A:141:GLN:HG2	2:E:312:ILE:CD1	2.21	0.69
1:D:5:GLU:HG2	3:D:235:HOH:O	1.92	0.69
1:B:124:ALA:O	1:B:126:PRO:HD3	1.93	0.69
1:C:154:ARG:C	1:C:156:LEU:H	2.01	0.69
1:D:44:GLN:H	1:D:44:GLN:NE2	1.92	0.68
1:A:156:LEU:HD12	1:A:157:ILE:H	1.59	0.68
1:A:73:MSE:O	1:A:81:HIS:HE1	1.78	0.67
1:A:1:MSE:HE3	1:A:6:GLY:HA3	1.75	0.67
1:A:141:GLN:HG2	2:E:312:ILE:HD12	1.76	0.66
1:B:74:LYS:HE3	3:B:225:HOH:O	1.95	0.65
1:B:137:MSE:HE1	2:F:312:ILE:HG12	1.78	0.65
1:A:31:GLY:HA3	3:A:235:HOH:O	1.95	0.65
1:A:70:GLU:HG3	3:A:218:HOH:O	1.96	0.65
1:B:36:ILE:CD1	1:B:46:ALA:HB2	2.26	0.65
1:A:101:TYR:HE1	2:E:312:ILE:HG22	1.62	0.65
1:D:29:ILE:HD12	1:D:68:VAL:HG22	1.79	0.65
1:D:126:PRO:HA	1:D:132:LYS:NZ	2.12	0.64
1:D:18:ASN:ND2	1:D:20:SER:H	1.96	0.63
1:D:18:ASN:HD22	1:D:19:PRO:N	1.95	0.63
1:D:18:ASN:HD22	1:D:18:ASN:C	2.05	0.63
1:B:18:ASN:HD22	1:B:19:PRO:CD	2.12	0.63
1:A:21:ASN:ND2	1:A:64:GLN:HE22	1.82	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:MSE:HE3	1:A:1:MSE:O	1.99	0.62
1:B:36:ILE:HD11	1:B:46:ALA:HB2	1.81	0.62
1:B:137:MSE:CG	1:B:140:ARG:HH21	2.13	0.61
1:C:204:GLY:HA3	3:C:265:HOH:O	2.00	0.61
1:D:94:ILE:HD11	2:H:310:LEU:HD23	1.82	0.61
1:B:19:PRO:HG3	1:B:61:GLU:HB2	1.83	0.60
1:C:51:ALA:O	1:C:55:GLN:HG3	2.02	0.60
1:D:32:PHE:CE2	1:D:36:ILE:HD11	2.36	0.60
1:C:94:ILE:CD1	2:G:312:ILE:HD11	2.32	0.60
1:B:18:ASN:HD22	1:B:19:PRO:HD2	1.65	0.60
1:A:45:ILE:HG22	1:A:49:LEU:HD22	1.85	0.59
1:C:93:LEU:O	1:C:96:VAL:HG22	2.03	0.59
1:C:58:GLN:HE22	1:C:60:TRP:HB3	1.68	0.59
1:C:81:HIS:HD2	3:C:245:HOH:O	1.86	0.59
1:B:153:ASP:OD1	1:B:155:THR:HB	2.03	0.59
1:A:22:ARG:HH21	1:A:23:GLN:CG	2.16	0.59
1:A:101:TYR:CE1	2:E:312:ILE:HG22	2.37	0.58
1:D:19:PRO:HG3	1:D:61:GLU:HB2	1.86	0.58
1:D:202:ALA:O	1:D:203:MSE:HB2	2.02	0.58
1:D:124:ALA:C	1:D:126:PRO:HD3	2.29	0.58
1:D:78:ARG:NH2	1:D:125:LEU:HD22	2.19	0.58
1:A:156:LEU:HD12	1:A:157:ILE:N	2.19	0.57
1:D:25:ASP:O	1:D:29:ILE:HG12	2.04	0.57
1:D:34:ASP:OD1	1:D:38:LYS:HE2	2.04	0.57
1:A:60:TRP:N	3:A:206:HOH:O	2.36	0.57
1:A:105:ARG:NH1	1:B:48:ARG:HD3	2.19	0.57
1:A:154:ARG:HG2	1:A:154:ARG:HH11	1.69	0.57
1:C:94:ILE:HD12	1:C:94:ILE:C	2.29	0.57
1:B:1:MSE:HE2	1:B:6:GLY:CA	2.35	0.57
1:C:117:LEU:HD23	3:C:232:HOH:O	2.05	0.57
1:B:1:MSE:HE2	1:B:1:MSE:HA	1.87	0.56
1:C:79:ARG:HG3	3:C:268:HOH:O	2.05	0.56
1:C:18:ASN:HD22	1:C:18:ASN:C	2.13	0.56
1:B:137:MSE:C	1:B:137:MSE:HE2	2.31	0.56
1:D:88:ARG:HB3	2:H:307:GLU:HG2	1.85	0.56
1:A:124:ALA:HB2	1:A:156:LEU:HD11	1.88	0.56
1:A:58:GLN:HE22	1:B:7:GLU:HG2	1.71	0.56
1:A:18:ASN:HD22	1:A:19:PRO:CD	2.19	0.56
1:A:19:PRO:HG3	1:A:61:GLU:HB2	1.87	0.56
1:B:5:GLU:CD	1:B:39:GLU:CG	2.79	0.55
1:B:58:GLN:HE22	1:B:60:TRP:HB3	1.71	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:MSE:HE2	1:B:58:GLN:CG	2.33	0.55
1:B:155:THR:HG21	3:B:248:HOH:O	2.06	0.55
1:A:111:LYS:O	1:A:115:ILE:HG12	2.06	0.55
1:C:84:VAL:O	1:C:90:LEU:HD13	2.06	0.55
1:C:203:MSE:O	1:C:203:MSE:HG2	2.06	0.55
1:D:137:MSE:HE3	1:D:141:GLN:CG	2.31	0.55
1:A:3:GLU:O	1:A:7:GLU:HG3	2.06	0.55
1:A:100:LYS:O	1:A:100:LYS:HG2	2.07	0.55
1:D:37:ASN:HD21	1:D:76:CYS:HA	1.72	0.54
1:D:58:GLN:HE22	1:D:60:TRP:HB3	1.71	0.54
1:D:73:MSE:O	1:D:81:HIS:HE1	1.89	0.54
1:C:18:ASN:ND2	1:C:20:SER:H	2.06	0.54
1:A:22:ARG:HH21	1:A:23:GLN:HG3	1.71	0.54
1:A:32:PHE:CE2	1:A:36:ILE:HD11	2.43	0.53
1:C:137:MSE:HE3	1:C:141:GLN:HG3	1.91	0.53
1:C:96:VAL:O	1:C:106:VAL:HG21	2.09	0.53
1:A:40:LEU:HD22	1:C:126:PRO:CB	2.37	0.53
1:C:98:SER:OG	2:G:312:ILE:HD12	2.09	0.53
1:C:94:ILE:CD1	2:G:310:LEU:HD23	2.27	0.52
1:B:137:MSE:HE1	2:F:312:ILE:CG1	2.39	0.52
1:D:125:LEU:HA	1:D:127:GLU:OE2	2.09	0.52
1:D:73:MSE:HE2	1:D:121:TRP:CE2	2.45	0.52
1:B:154:ARG:C	1:B:156:LEU:N	2.63	0.52
1:B:108:GLU:O	1:B:112:THR:HG23	2.09	0.52
1:D:126:PRO:HA	1:D:132:LYS:HZ3	1.75	0.51
1:B:109:LYS:HG3	3:B:245:HOH:O	2.10	0.51
1:B:115:ILE:HG22	1:B:148:PRO:HG3	1.93	0.51
1:C:91:ASN:O	1:C:94:ILE:HG13	2.11	0.51
1:B:63:LEU:HD13	1:B:110:VAL:HG22	1.93	0.51
1:D:21:ASN:HB3	1:D:23:GLN:O	2.11	0.51
1:A:129:ALA:O	1:A:132:LYS:HG2	2.11	0.50
1:D:18:ASN:HD22	1:D:19:PRO:CD	2.25	0.50
1:D:40:LEU:HD13	1:D:40:LEU:O	2.12	0.50
1:C:73:MSE:O	1:C:81:HIS:HE1	1.94	0.50
1:C:150:ILE:O	1:C:152:VAL:HG23	2.11	0.50
1:A:10:GLU:CG	1:B:19:PRO:HG2	2.42	0.50
1:A:18:ASN:HD22	1:A:19:PRO:N	2.10	0.50
1:B:203:MSE:HE2	3:B:261:HOH:O	2.11	0.49
1:B:82:ASN:O	1:B:86:LYS:HD3	2.12	0.49
1:C:26:TRP:O	1:C:30:ILE:HG12	2.12	0.49
1:C:84:VAL:HG13	1:C:89:PHE:CE2	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:126:PRO:HA	1:C:132:LYS:NZ	2.27	0.49
1:D:126:PRO:HA	1:D:132:LYS:HZ1	1.77	0.49
1:B:5:GLU:OE2	1:B:39:GLU:HG2	2.13	0.49
1:C:154:ARG:C	1:C:156:LEU:N	2.70	0.49
1:B:153:ASP:OD1	1:B:155:THR:N	2.46	0.49
1:C:81:HIS:CD2	3:C:245:HOH:O	2.63	0.49
1:B:39:GLU:HG3	3:B:254:HOH:O	2.12	0.48
1:A:51:ALA:O	1:A:55:GLN:HG3	2.13	0.48
1:B:155:THR:O	1:B:156:LEU:HG	2.13	0.48
1:C:136:HIS:HB3	3:C:244:HOH:O	2.13	0.48
1:C:95:LYS:HG2	1:C:101:TYR:CD2	2.48	0.48
1:B:36:ILE:HD12	1:B:46:ALA:HB2	1.95	0.48
1:C:18:ASN:HD22	1:C:19:PRO:N	2.11	0.48
1:D:141:GLN:HE21	1:D:141:GLN:HA	1.79	0.48
1:A:19:PRO:HB3	1:A:58:GLN:NE2	2.29	0.48
1:B:5:GLU:CG	1:B:39:GLU:HG2	2.45	0.47
1:D:20:SER:O	1:D:21:ASN:C	2.57	0.47
1:A:203:MSE:CE	1:B:107:SER:HB3	2.45	0.47
1:B:70:GLU:HG2	1:B:117:LEU:HD21	1.96	0.47
1:B:135:TYR:HE2	1:B:139:LYS:HD2	1.78	0.47
1:D:110:VAL:O	1:D:114:VAL:HG23	2.14	0.47
1:C:94:ILE:HD13	2:G:312:ILE:HD11	1.95	0.47
1:A:1:MSE:HE2	1:B:58:GLN:HA	1.97	0.47
1:B:66:LEU:HD13	1:B:113:LYS:CG	2.44	0.47
1:C:124:ALA:C	1:C:126:PRO:HD3	2.39	0.47
1:B:88:ARG:HG3	1:B:88:ARG:HH11	1.80	0.47
1:C:32:PHE:CE2	1:C:36:ILE:HD11	2.49	0.47
1:D:35:GLN:HE21	1:D:35:GLN:HA	1.80	0.47
1:C:90:LEU:O	1:C:94:ILE:HG23	2.15	0.46
1:B:78:ARG:NH2	1:B:127:GLU:HG2	2.30	0.46
1:C:87:PHE:CE1	2:G:309:LEU:HD13	2.50	0.46
1:C:37:ASN:HD21	1:C:76:CYS:HA	1.80	0.46
1:D:18:ASN:ND2	1:D:18:ASN:C	2.69	0.46
1:A:15:LYS:HD2	1:A:28:TYR:CZ	2.50	0.46
1:D:44:GLN:H	1:D:44:GLN:CD	2.22	0.46
1:A:22:ARG:O	1:A:23:GLN:OE1	2.34	0.46
1:B:5:GLU:OE1	1:B:39:GLU:CD	2.58	0.46
1:B:26:TRP:O	1:B:30:ILE:HG12	2.16	0.46
1:B:203:MSE:HE3	1:B:203:MSE:HA	1.97	0.46
1:B:137:MSE:HG2	1:B:140:ARG:HH21	1.81	0.46
1:C:102:LEU:O	1:C:106:VAL:HG13	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:153:ASP:OD1	1:B:153:ASP:C	2.59	0.45
1:D:69:LEU:HD21	1:D:121:TRP:HH2	1.80	0.45
1:A:1:MSE:CE	1:A:1:MSE:O	2.64	0.45
1:B:101:TYR:CE1	2:F:312:ILE:HG22	2.51	0.45
1:A:1:MSE:HE1	1:B:58:GLN:HG2	1.93	0.45
1:A:22:ARG:HH21	1:A:23:GLN:CB	2.30	0.45
1:A:132:LYS:HE3	1:A:132:LYS:HB3	1.80	0.45
1:A:154:ARG:HH11	1:A:154:ARG:CG	2.29	0.44
1:C:203:MSE:HE1	3:D:249:HOH:O	2.16	0.44
1:B:132:LYS:HB3	3:B:214:HOH:O	2.16	0.44
1:C:106:VAL:HG23	1:C:111:LYS:HG3	1.99	0.44
1:D:16:ALA:HB1	1:D:29:ILE:HD13	1.99	0.44
1:D:107:SER:O	1:D:111:LYS:HE2	2.17	0.44
2:E:312:ILE:HD12	2:E:312:ILE:C	2.43	0.44
1:D:111:LYS:O	1:D:115:ILE:HG12	2.18	0.43
1:C:53:LYS:HB3	1:C:65:ALA:HB2	1.99	0.43
1:A:22:ARG:HH21	1:A:23:GLN:HB2	1.83	0.43
1:B:1:MSE:CE	1:B:6:GLY:CA	2.96	0.43
1:B:102:LEU:HD22	1:B:105:ARG:HH21	1.83	0.43
1:C:12:TRP:HE1	1:C:35:GLN:NE2	2.17	0.43
1:A:18:ASN:HD22	1:A:19:PRO:HD2	1.82	0.43
1:B:49:LEU:HD12	1:B:49:LEU:HA	1.84	0.43
1:A:153:ASP:OD1	1:A:153:ASP:C	2.61	0.42
1:B:59:GLU:OE2	1:B:109:LYS:HD2	2.19	0.42
1:C:78:ARG:HG3	1:C:82:ASN:HD21	1.83	0.42
1:C:125:LEU:N	1:C:126:PRO:HD3	2.33	0.42
1:A:137:MSE:HE1	2:E:310:LEU:O	2.19	0.42
1:B:105:ARG:HD2	1:B:105:ARG:C	2.43	0.42
1:B:20:SER:O	1:B:21:ASN:C	2.61	0.42
1:C:137:MSE:HA	1:C:140:ARG:HE	1.85	0.42
1:D:124:ALA:HB2	1:D:156:LEU:HD22	2.02	0.42
1:B:58:GLN:NE2	1:B:61:GLU:H	2.18	0.42
1:B:137:MSE:HE2	1:B:137:MSE:O	2.20	0.42
1:D:94:ILE:CD1	2:H:310:LEU:HD23	2.49	0.42
1:D:90:LEU:O	1:D:94:ILE:HG23	2.20	0.42
1:C:94:ILE:HD13	2:G:312:ILE:CG1	2.50	0.41
1:D:98:SER:OG	2:H:312:ILE:HD12	2.19	0.41
1:D:58:GLN:NE2	1:D:61:GLU:H	2.19	0.41
1:A:137:MSE:HE2	2:E:311:HIS:HA	2.02	0.41
1:C:10:GLU:CG	1:D:19:PRO:HG2	2.50	0.41
1:D:94:ILE:HD11	2:H:310:LEU:HB3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:94:ILE:CD1	2:H:310:LEU:HB3	2.50	0.41
1:A:126:PRO:HA	1:A:132:LYS:NZ	2.36	0.41
1:B:73:MSE:SE	1:B:84:VAL:HG21	2.71	0.41
1:A:18:ASN:HD22	1:A:18:ASN:C	2.29	0.41
1:B:155:THR:O	1:B:155:THR:CG2	2.61	0.41
1:D:202:ALA:O	1:D:203:MSE:CB	2.68	0.41
1:D:14:ASN:ND2	1:D:53:LYS:NZ	2.69	0.41
1:D:35:GLN:HA	1:D:35:GLN:NE2	2.36	0.41
1:D:100:LYS:HA	1:D:100:LYS:HD3	1.94	0.41
1:C:13:LEU:HA	1:C:13:LEU:HD12	1.83	0.40
1:C:69:LEU:HD21	1:C:121:TRP:HH2	1.87	0.40
1:D:10:GLU:OE1	1:D:49:LEU:HD22	2.21	0.40
1:D:119:TYR:CZ	1:D:147:ASP:HB3	2.56	0.40
2:H:312:ILE:OXT	2:H:312:ILE:HG22	2.21	0.40
1:A:91:ASN:HA	1:A:94:ILE:HD12	2.03	0.40
1:B:47:VAL:HB	1:B:89:PHE:CD1	2.55	0.40
1:B:85:GLY:HA3	1:B:128:GLU:OE1	2.21	0.40
1:C:126:PRO:HD2	3:C:227:HOH:O	2.21	0.40
1:C:137:MSE:HE3	1:C:141:GLN:CG	2.51	0.40
1:C:152:VAL:HG12	1:C:153:ASP:N	2.35	0.40
1:B:18:ASN:HD22	1:B:19:PRO:N	2.19	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	159/171 (93%)	154 (97%)	4 (2%)	1 (1%)	21	32
1	B	157/171 (92%)	148 (94%)	7 (4%)	2 (1%)	9	14
1	C	157/171 (92%)	153 (98%)	4 (2%)	0	100	100
1	D	159/171 (93%)	147 (92%)	11 (7%)	1 (1%)	21	32

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	E	6/12 (50%)	5 (83%)	1 (17%)	0	100	100
2	F	6/12 (50%)	5 (83%)	0	1 (17%)	0	0
2	G	6/12 (50%)	6 (100%)	0	0	100	100
2	H	7/12 (58%)	7 (100%)	0	0	100	100
All	All	657/732 (90%)	625 (95%)	27 (4%)	5 (1%)	16	25

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	103	GLY
1	B	105	ARG
2	F	306	ASP
1	A	103	GLY
1	B	155	THR

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	141/145 (97%)	135 (96%)	6 (4%)	26	44
1	B	140/145 (97%)	132 (94%)	8 (6%)	18	33
1	C	140/145 (97%)	136 (97%)	4 (3%)	37	60
1	D	141/145 (97%)	132 (94%)	9 (6%)	16	28
2	E	8/12 (67%)	8 (100%)	0	100	100
2	F	8/12 (67%)	8 (100%)	0	100	100
2	G	8/12 (67%)	7 (88%)	1 (12%)	4	6
2	H	9/12 (75%)	9 (100%)	0	100	100
All	All	595/628 (95%)	567 (95%)	28 (5%)	23	41

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	LEU
1	A	49	LEU
1	A	58	GLN
1	A	63	LEU
1	A	132	LYS
1	A	141	GLN
1	B	1	MSE
1	B	40	LEU
1	B	49	LEU
1	B	58	GLN
1	B	63	LEU
1	B	79	ARG
1	B	108	GLU
1	B	137	MSE
1	C	58	GLN
1	C	90	LEU
1	C	94	ILE
1	C	105	ARG
2	G	307	GLU
1	D	18	ASN
1	D	22	ARG
1	D	44	GLN
1	D	63	LEU
1	D	78	ARG
1	D	94	ILE
1	D	118	LEU
1	D	127	GLU
1	D	145	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (35) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	ASN
1	A	23	GLN
1	A	58	GLN
1	A	64	GLN
1	A	81	HIS
1	A	82	ASN
1	B	18	ASN
1	B	21	ASN
1	B	58	GLN
1	B	75	ASN
1	B	82	ASN

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Mol	Chain	Res	Type
1	B	141	GLN
2	F	311	HIS
1	C	14	ASN
1	C	18	ASN
1	C	21	ASN
1	C	35	GLN
1	C	37	ASN
1	C	58	GLN
1	C	64	GLN
1	C	75	ASN
1	C	81	HIS
1	C	82	ASN
1	C	141	GLN
1	D	14	ASN
1	D	18	ASN
1	D	21	ASN
1	D	35	GLN
1	D	37	ASN
1	D	44	GLN
1	D	55	GLN
1	D	58	GLN
1	D	64	GLN
1	D	81	HIS
1	D	141	GLN

5.3.3 RNA ⓘ

There are no RNA molecules 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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

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, 95th 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 ²)	Q<0.9
1	A	156/171 (91%)	0.10	5 (3%) 50 46	21, 35, 57, 77	0
1	B	154/171 (90%)	0.03	4 (2%) 57 53	24, 33, 53, 61	0
1	C	154/171 (90%)	-0.01	1 (0%) 85 83	22, 35, 50, 61	0
1	D	156/171 (91%)	0.00	4 (2%) 57 53	21, 36, 53, 74	0
2	E	8/12 (66%)	1.61	2 (25%) 2 1	51, 58, 63, 69	0
2	F	8/12 (66%)	1.29	2 (25%) 2 1	51, 57, 68, 69	0
2	G	8/12 (66%)	0.86	1 (12%) 8 6	45, 52, 66, 66	0
2	H	9/12 (75%)	1.24	2 (22%) 2 2	47, 53, 68, 69	0
All	All	653/732 (89%)	0.09	21 (3%) 50 46	21, 35, 57, 77	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	202	ALA	5.6
1	A	157	ILE	4.8
1	D	157	ILE	4.7
2	E	311	HIS	3.8
1	B	5	GLU	3.5
1	A	202	ALA	3.5
1	A	156	LEU	2.9
1	B	156	LEU	2.8
1	B	22	ARG	2.8
1	B	23	GLN	2.7
2	H	312	ILE	2.5
1	D	35	GLN	2.5
1	D	105	ARG	2.4
2	G	311	HIS	2.4
2	H	311	HIS	2.3
2	F	311	HIS	2.2

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Mol	Chain	Res	Type	RSRZ
2	E	305	SER	2.2
1	A	136	HIS	2.2
1	C	204	GLY	2.2
1	A	22	ARG	2.1
2	F	312	ILE	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.