



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 11, 2024 – 10:56 PM EST

PDB ID : 2PIF
Title : Crystal structure of UPF0317 protein PSPTO_5379 from *Pseudomonas syringae* pv. tomato. NorthEast Structural Genomics target PsR181
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Deposited on : 2007-04-13
Resolution : 2.30 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	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)

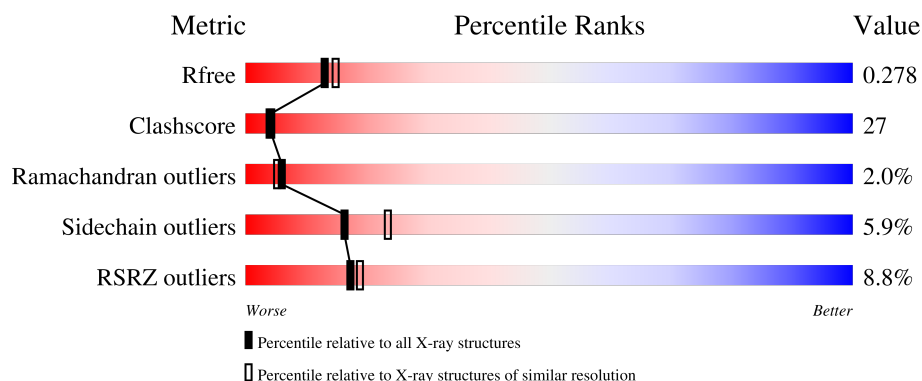
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	164625	5963 (2.30-2.30)
Clashscore	180529	6698 (2.30-2.30)
Ramachandran outliers	177936	6640 (2.30-2.30)
Sidechain outliers	177891	6640 (2.30-2.30)
RSRZ outliers	164620	5963 (2.30-2.30)

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	276	
1	B	276	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3994 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 UPF0317 protein PSPTO_5379.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	252	Total	C	N	O	S	Se	0	0	0
			1898	1191	340	354	5	8			
1	B	253	Total	C	N	O	S	Se	0	0	0
			1906	1195	341	357	5	8			

There are 34 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	modified residue	UNP Q87UC6
A	37	MSE	MET	modified residue	UNP Q87UC6
A	91	MSE	MET	modified residue	UNP Q87UC6
A	117	MSE	MET	modified residue	UNP Q87UC6
A	150	MSE	MET	modified residue	UNP Q87UC6
A	167	MSE	MET	modified residue	UNP Q87UC6
A	171	MSE	MET	modified residue	UNP Q87UC6
A	241	MSE	MET	modified residue	UNP Q87UC6
A	256	MSE	MET	modified residue	UNP Q87UC6
A	269	LEU	-	cloning artifact	UNP Q87UC6
A	270	GLU	-	cloning artifact	UNP Q87UC6
A	271	HIS	-	cloning artifact	UNP Q87UC6
A	272	HIS	-	cloning artifact	UNP Q87UC6
A	273	HIS	-	cloning artifact	UNP Q87UC6
A	274	HIS	-	cloning artifact	UNP Q87UC6
A	275	HIS	-	cloning artifact	UNP Q87UC6
A	276	HIS	-	cloning artifact	UNP Q87UC6
B	1	MSE	MET	modified residue	UNP Q87UC6
B	37	MSE	MET	modified residue	UNP Q87UC6
B	91	MSE	MET	modified residue	UNP Q87UC6
B	117	MSE	MET	modified residue	UNP Q87UC6
B	150	MSE	MET	modified residue	UNP Q87UC6
B	167	MSE	MET	modified residue	UNP Q87UC6
B	171	MSE	MET	modified residue	UNP Q87UC6
B	241	MSE	MET	modified residue	UNP Q87UC6

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Chain	Residue	Modelled	Actual	Comment	Reference
B	256	MSE	MET	modified residue	UNP Q87UC6
B	269	LEU	-	cloning artifact	UNP Q87UC6
B	270	GLU	-	cloning artifact	UNP Q87UC6
B	271	HIS	-	cloning artifact	UNP Q87UC6
B	272	HIS	-	cloning artifact	UNP Q87UC6
B	273	HIS	-	cloning artifact	UNP Q87UC6
B	274	HIS	-	cloning artifact	UNP Q87UC6
B	275	HIS	-	cloning artifact	UNP Q87UC6
B	276	HIS	-	cloning artifact	UNP Q87UC6

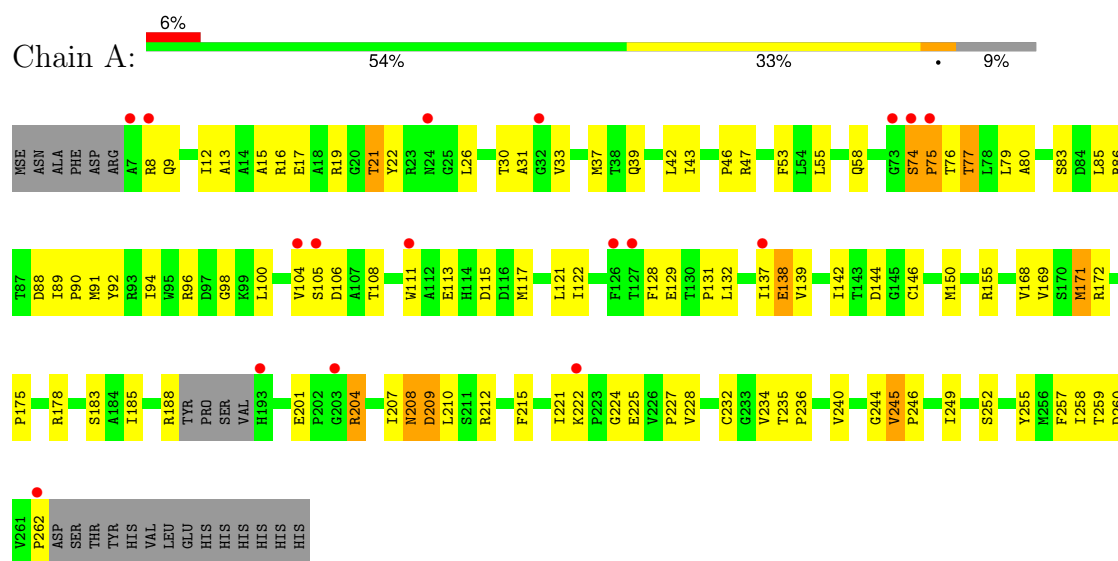
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	99	Total O 99 99	0	0
2	B	91	Total O 91 91	0	0

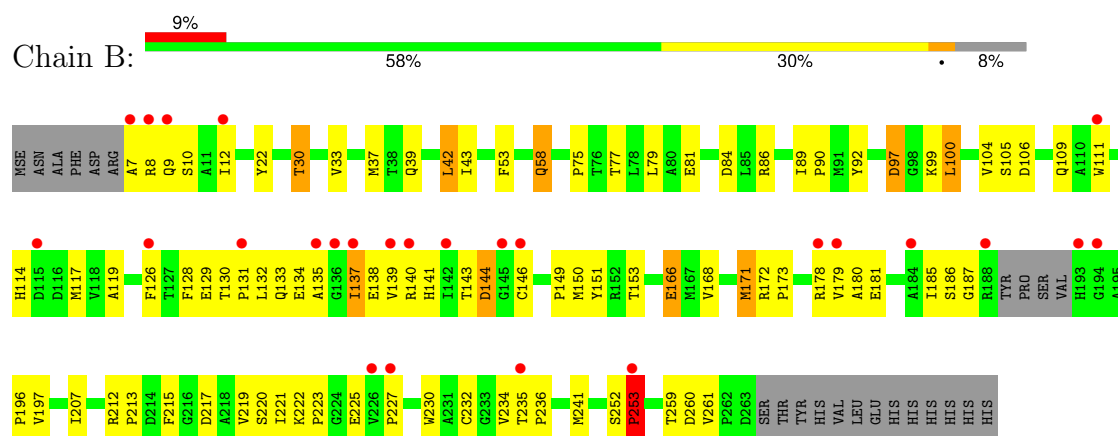
3 Residue-property plots [i](#)

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: UPF0317 protein PSPTO_5379



• Molecule 1: UPF0317 protein PSPTO_5379



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	66.31Å 85.95Å 93.94Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.22 – 2.30 41.22 – 2.30	Depositor EDS
% Data completeness (in resolution range)	91.6 (41.22-2.30) 95.6 (41.22-2.30)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	10.89 (at 2.24Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.241 , 0.272 0.252 , 0.278	Depositor DCC
R_{free} test set	653 reflections (2.76%)	wwPDB-VP
Wilson B-factor (Å ²)	32.5	Xtriage
Anisotropy	0.301	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 37.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	3994	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.55% of the height of the origin peak. No significant pseudotranslation is detected.*

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	0.36	0/1937	0.63	0/2629
1	B	0.35	0/1945	0.62	1/2640 (0.0%)
All	All	0.36	0/3882	0.63	1/5269 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	144	ASP	CB-CG-OD2	5.26	123.04	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	134	GLU	Peptide

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	1898	0	1854	96	0
1	B	1906	0	1858	105	0
2	A	99	0	0	14	0
2	B	91	0	0	16	0
All	All	3994	0	3712	201	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 27.

All (201) 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:168:VAL:H	1:B:235:THR:HG22	1.14	1.10
1:B:104:VAL:HG22	1:B:106:ASP:H	1.22	1.04
1:A:74:SER:HB2	1:A:75:PRO:HA	1.45	0.99
1:B:196:PRO:HG3	1:B:230:TRP:NE1	1.79	0.98
1:B:140:ARG:HG3	1:B:219:VAL:HG21	1.44	0.97
1:A:137:ILE:O	1:A:138:GLU:HB3	1.67	0.94
1:A:208:ASN:O	1:A:209:ASP:HB2	1.68	0.92
1:B:150:MSE:HB2	1:B:215:PHE:HB2	1.53	0.91
1:B:129:GLU:OE1	1:B:139:VAL:HG11	1.72	0.89
1:A:83:SER:HB2	2:A:287:HOH:O	1.72	0.88
1:B:37:MSE:HE1	1:B:260:ASP:C	1.98	0.83
1:A:13:ALA:O	1:A:17:GLU:HG2	1.78	0.83
1:B:137:ILE:HD11	1:B:178:ARG:HD2	1.60	0.83
1:B:140:ARG:HG3	1:B:219:VAL:CG2	2.10	0.82
1:B:171:MSE:HE3	1:B:227:PRO:HA	1.62	0.81
1:A:171:MSE:HE2	1:A:172:ARG:N	1.96	0.81
1:B:130:THR:HB	1:B:131:PRO:HD3	1.64	0.80
1:B:222:LYS:NZ	1:B:222:LYS:HB3	1.97	0.79
1:A:19:ARG:HD2	2:A:287:HOH:O	1.82	0.79
1:B:37:MSE:HE2	1:B:259:THR:O	1.84	0.77
1:B:222:LYS:HB2	1:B:225:GLU:OE2	1.84	0.77
1:A:104:VAL:HG22	1:A:106:ASP:H	1.50	0.76
1:B:140:ARG:CG	1:B:219:VAL:HG21	2.15	0.76
1:B:149:PRO:HG3	1:B:217:ASP:O	1.86	0.76
1:B:171:MSE:CE	1:B:227:PRO:HA	2.16	0.76
1:B:37:MSE:HE1	1:B:261:VAL:N	2.00	0.75
1:B:196:PRO:HG3	1:B:230:TRP:HE1	1.48	0.75
1:B:75:PRO:HB3	1:B:86:ARG:HD2	1.69	0.75
1:B:241:MSE:HG2	2:B:302:HOH:O	1.87	0.74
1:B:30:THR:O	1:B:33:VAL:HG12	1.87	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:252:SER:HB3	1:B:253:PRO:HD2	1.69	0.73
1:A:74:SER:HB2	1:A:75:PRO:CA	2.17	0.73
1:A:232:CYS:SG	1:A:234:VAL:HG22	2.29	0.73
1:A:19:ARG:HB2	1:A:88:ASP:HB3	1.72	0.72
1:A:171:MSE:HE1	1:A:227:PRO:HA	1.73	0.70
1:B:196:PRO:HG3	1:B:230:TRP:CE2	2.26	0.70
1:B:99:LYS:O	1:B:100:LEU:HB3	1.89	0.70
1:A:150:MSE:HB2	1:A:215:PHE:HB2	1.74	0.70
1:A:12:ILE:HD13	1:A:37:MSE:HE1	1.74	0.68
1:B:166:GLU:HG2	2:B:355:HOH:O	1.94	0.68
1:A:262:PRO:HG3	2:A:344:HOH:O	1.94	0.67
1:A:75:PRO:HB2	1:A:85:LEU:H	1.60	0.67
1:B:168:VAL:O	1:B:232:CYS:HB3	1.95	0.66
1:B:75:PRO:CB	1:B:86:ARG:HD2	2.26	0.65
1:B:168:VAL:N	1:B:235:THR:HG22	1.99	0.65
1:A:19:ARG:HA	1:A:22:TYR:CD2	2.33	0.64
1:A:171:MSE:HE2	1:A:172:ARG:H	1.61	0.63
1:B:137:ILE:HD11	1:B:178:ARG:CD	2.28	0.63
1:B:171:MSE:HE3	1:B:227:PRO:CA	2.28	0.62
1:A:79:LEU:HD22	1:A:121:LEU:HD12	1.80	0.62
1:A:74:SER:HA	2:A:282:HOH:O	2.00	0.62
1:B:222:LYS:HB3	1:B:222:LYS:HZ3	1.66	0.61
1:A:89:ILE:HB	1:A:92:TYR:CZ	2.36	0.61
1:A:171:MSE:CE	1:A:228:VAL:H	2.14	0.61
1:B:150:MSE:HE2	1:B:232:CYS:SG	2.41	0.60
1:B:37:MSE:HE3	1:B:37:MSE:HA	1.81	0.60
1:B:241:MSE:HE3	2:B:302:HOH:O	2.02	0.59
1:B:111:TRP:HZ3	2:B:299:HOH:O	1.85	0.59
1:A:37:MSE:CE	1:A:260:ASP:HA	2.31	0.59
1:B:109:GLN:HG2	2:B:352:HOH:O	2.03	0.58
1:A:168:VAL:HG11	1:A:234:VAL:HG23	1.86	0.58
1:B:8:ARG:HG2	1:B:9:GLN:HE21	1.69	0.58
1:B:84:ASP:OD1	1:B:86:ARG:HD3	2.02	0.58
1:B:30:THR:HG23	1:B:39:GLN:OE1	2.03	0.57
1:B:99:LYS:HA	2:B:302:HOH:O	2.03	0.57
1:B:22:TYR:CD1	1:B:90:PRO:HG3	2.40	0.57
1:A:188:ARG:HD3	2:A:361:HOH:O	2.05	0.57
1:B:30:THR:CG2	1:B:39:GLN:OE1	2.53	0.56
1:B:137:ILE:CD1	1:B:178:ARG:HD2	2.34	0.56
1:A:86:ARG:HG2	1:A:111:TRP:CG	2.41	0.56
1:A:235:THR:N	1:A:236:PRO:CD	2.69	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:172:ARG:HG2	1:B:172:ARG:HH11	1.70	0.56
1:B:137:ILE:N	1:B:137:ILE:HD13	2.21	0.56
1:B:151:TYR:HB3	1:B:207:ILE:HD12	1.86	0.56
1:B:128:PHE:C	1:B:131:PRO:HD2	2.26	0.55
1:A:117:MSE:HE2	1:A:249:ILE:HD12	1.88	0.55
1:A:8:ARG:O	1:A:12:ILE:HG12	2.06	0.55
1:A:240:VAL:HG13	1:A:245:VAL:HG22	1.88	0.55
1:B:146:CYS:HB3	1:B:217:ASP:HB2	1.87	0.55
1:A:37:MSE:SE	2:A:344:HOH:O	2.75	0.55
1:A:132:LEU:HD23	1:A:185:ILE:HG21	1.88	0.54
1:A:207:ILE:HG21	1:A:210:LEU:HD23	1.87	0.54
1:A:53:PHE:CZ	1:A:122:ILE:HD11	2.42	0.54
1:B:53:PHE:HE1	1:B:236:PRO:HB3	1.71	0.54
1:B:180:ALA:HB2	2:B:293:HOH:O	2.07	0.54
1:B:53:PHE:CE1	1:B:236:PRO:HB3	2.42	0.54
1:B:8:ARG:HD2	2:B:303:HOH:O	2.08	0.53
1:A:12:ILE:HD13	1:A:37:MSE:CE	2.37	0.53
1:A:39:GLN:NE2	1:A:258:ILE:HG12	2.24	0.53
1:B:8:ARG:HB2	2:B:282:HOH:O	2.08	0.53
1:B:222:LYS:HZ3	1:B:223:PRO:HD2	1.73	0.53
1:A:42:LEU:HG	1:A:43:ILE:N	2.24	0.52
1:B:171:MSE:HE3	1:B:227:PRO:CB	2.39	0.52
1:B:221:ILE:HD12	1:B:221:ILE:N	2.24	0.52
1:B:235:THR:N	1:B:236:PRO:HD2	2.24	0.52
1:A:129:GLU:OE2	1:A:139:VAL:HG11	2.10	0.52
1:A:90:PRO:HB2	1:A:255:TYR:HD1	1.75	0.52
1:B:37:MSE:HE1	1:B:261:VAL:C	2.31	0.51
1:B:42:LEU:HG	1:B:43:ILE:N	2.24	0.51
1:B:37:MSE:HE3	1:B:37:MSE:CA	2.40	0.51
1:B:178:ARG:NH1	1:B:181:GLU:OE2	2.41	0.51
1:A:58:GLN:NE2	2:A:277:HOH:O	2.44	0.51
1:B:97:ASP:O	1:B:99:LYS:HG3	2.10	0.51
1:A:175:PRO:HG2	1:A:178:ARG:CG	2.41	0.51
1:A:150:MSE:HE1	2:A:352:HOH:O	2.11	0.50
1:A:55:LEU:HD12	1:A:58:GLN:HG3	1.94	0.50
1:B:81:GLU:HG3	2:B:349:HOH:O	2.11	0.50
1:A:16:ARG:HD3	2:A:348:HOH:O	2.12	0.49
1:A:212:ARG:HG2	2:A:358:HOH:O	2.12	0.49
1:A:79:LEU:HB3	1:A:259:THR:HB	1.93	0.49
1:B:219:VAL:HG22	1:B:220:SER:N	2.26	0.49
1:A:31:ALA:HB2	2:A:365:HOH:O	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:30:THR:HA	1:A:33:VAL:HG21	1.93	0.49
1:A:208:ASN:O	1:A:209:ASP:CB	2.49	0.49
1:B:140:ARG:CG	1:B:219:VAL:CG2	2.83	0.49
1:B:128:PHE:O	1:B:131:PRO:HD2	2.13	0.48
1:B:171:MSE:HE1	1:B:173:PRO:HA	1.94	0.48
1:A:74:SER:OG	1:A:111:TRP:HZ3	1.95	0.48
1:B:104:VAL:HG22	1:B:105:SER:N	2.28	0.48
1:B:133:GLN:HG2	1:B:139:VAL:HG23	1.94	0.48
1:B:141:HIS:O	1:B:144:ASP:O	2.31	0.48
1:A:168:VAL:HG11	1:A:234:VAL:CG2	2.43	0.48
1:A:175:PRO:HG2	1:A:178:ARG:HG2	1.96	0.48
1:A:19:ARG:HA	1:A:22:TYR:HD2	1.79	0.48
1:B:8:ARG:HG3	2:B:279:HOH:O	2.13	0.47
1:A:30:THR:HA	1:A:33:VAL:CG2	2.44	0.47
1:B:37:MSE:CE	1:B:260:ASP:C	2.76	0.47
1:A:128:PHE:O	1:A:131:PRO:HG2	2.14	0.47
1:A:55:LEU:HD12	1:A:58:GLN:CG	2.45	0.47
1:A:257:PHE:HE1	2:A:287:HOH:O	1.96	0.47
1:A:150:MSE:HE2	1:A:232:CYS:SG	2.55	0.46
1:B:111:TRP:HH2	1:B:119:ALA:HB2	1.79	0.46
1:B:171:MSE:HE1	1:B:173:PRO:CA	2.46	0.46
1:A:91:MSE:HG2	1:A:252:SER:OG	2.14	0.46
1:B:171:MSE:HE3	1:B:227:PRO:HB3	1.96	0.46
1:B:79:LEU:CD2	1:B:261:VAL:HG21	2.46	0.46
1:B:185:ILE:C	1:B:187:GLY:H	2.19	0.46
1:B:9:GLN:N	2:B:306:HOH:O	2.49	0.46
1:A:80:ALA:HA	1:A:260:ASP:OD2	2.16	0.46
1:B:7:ALA:HB3	1:B:10:SER:HB2	1.98	0.46
1:A:155:ARG:NH2	2:A:346:HOH:O	2.38	0.46
1:A:240:VAL:HA	1:A:245:VAL:HG13	1.98	0.46
1:B:58:GLN:NE2	2:B:278:HOH:O	2.48	0.46
1:A:37:MSE:HE1	1:A:260:ASP:HA	1.98	0.46
1:B:129:GLU:O	1:B:133:GLN:HG3	2.15	0.45
1:B:234:VAL:C	1:B:236:PRO:HD2	2.36	0.45
1:B:253:PRO:HB3	2:B:310:HOH:O	2.15	0.45
1:A:21:THR:HG21	1:A:26:LEU:HD22	1.98	0.45
1:A:171:MSE:CE	1:A:227:PRO:HA	2.43	0.45
1:B:132:LEU:O	1:B:135:ALA:N	2.45	0.45
1:A:37:MSE:HE3	1:A:260:ASP:HA	1.97	0.45
1:A:96:ARG:HA	1:A:246:PRO:O	2.17	0.45
1:A:77:THR:HG21	1:A:121:LEU:HD11	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:171:MSE:HE2	1:B:172:ARG:O	2.17	0.44
1:A:43:ILE:O	1:A:43:ILE:HG23	2.17	0.44
1:A:94:ILE:HD11	1:A:104:VAL:HG11	2.00	0.44
1:A:171:MSE:CE	1:A:228:VAL:N	2.80	0.44
1:A:204:ARG:HD2	2:A:337:HOH:O	2.17	0.44
1:B:139:VAL:O	1:B:140:ARG:C	2.55	0.44
1:B:234:VAL:HG12	1:B:234:VAL:O	2.18	0.44
1:A:46:PRO:O	1:A:47:ARG:C	2.56	0.43
1:B:8:ARG:CG	1:B:9:GLN:HE21	2.30	0.43
1:B:43:ILE:HG23	1:B:43:ILE:O	2.18	0.43
1:A:53:PHE:HZ	1:A:122:ILE:HD11	1.83	0.43
1:A:142:ILE:C	1:A:144:ASP:N	2.72	0.43
1:A:12:ILE:HA	1:A:37:MSE:HE2	2.00	0.43
1:B:222:LYS:HB3	1:B:222:LYS:HZ2	1.76	0.43
1:A:222:LYS:CG	1:A:225:GLU:HG3	2.48	0.43
1:B:79:LEU:HD23	1:B:261:VAL:HG21	2.00	0.43
1:B:151:TYR:CE1	1:B:213:PRO:HA	2.53	0.43
1:A:96:ARG:NH2	1:A:113:GLU:OE1	2.52	0.43
1:B:89:ILE:HB	1:B:92:TYR:CZ	2.53	0.43
1:B:126:PHE:N	1:B:126:PHE:CD2	2.86	0.43
1:B:153:THR:HA	2:B:326:HOH:O	2.18	0.43
1:A:98:GLY:HA3	1:A:244:GLY:HA2	2.01	0.43
1:A:207:ILE:HG21	1:A:210:LEU:CD2	2.49	0.43
1:A:171:MSE:HE1	1:A:228:VAL:H	1.84	0.42
1:A:74:SER:OG	1:A:111:TRP:CZ3	2.70	0.42
1:B:232:CYS:O	1:B:235:THR:HG23	2.19	0.42
1:A:208:ASN:HD22	1:A:208:ASN:HA	1.68	0.42
1:A:86:ARG:NH1	1:A:108:THR:HG23	2.35	0.42
1:A:37:MSE:HE3	1:A:259:THR:O	2.20	0.42
1:A:86:ARG:HG2	1:A:111:TRP:CD1	2.55	0.42
1:B:172:ARG:HB2	1:B:230:TRP:HZ3	1.85	0.42
1:A:222:LYS:HG2	1:A:224:GLY:H	1.85	0.42
1:B:137:ILE:CD1	1:B:178:ARG:CD	2.95	0.42
1:B:8:ARG:C	1:B:10:SER:H	2.23	0.41
1:B:114:HIS:O	1:B:117:MSE:HG2	2.20	0.41
1:B:137:ILE:HD11	1:B:178:ARG:NE	2.34	0.41
1:A:201:GLU:HG3	1:A:204:ARG:NH1	2.36	0.41
1:A:221:ILE:HG13	1:A:225:GLU:HB2	2.02	0.41
1:A:235:THR:HB	1:A:236:PRO:HD3	2.03	0.41
1:B:212:ARG:HD2	2:B:322:HOH:O	2.20	0.41
1:A:86:ARG:HD2	1:A:108:THR:HA	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:128:PHE:O	1:A:132:LEU:HG	2.21	0.40
1:A:222:LYS:HG2	1:A:225:GLU:HG3	2.02	0.40
1:A:15:ALA:O	1:A:19:ARG:HG2	2.21	0.40
1:B:172:ARG:HH11	1:B:172:ARG:CG	2.35	0.40
1:A:104:VAL:HG22	1:A:105:SER:N	2.36	0.40
1:A:150:MSE:HA	1:A:169:VAL:O	2.22	0.40
1:B:30:THR:O	1:B:33:VAL:CG1	2.64	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	248/276 (90%)	230 (93%)	13 (5%)	5 (2%)	6	5
1	B	249/276 (90%)	230 (92%)	14 (6%)	5 (2%)	6	5
All	All	497/552 (90%)	460 (93%)	27 (5%)	10 (2%)	6	5

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	74	SER
1	A	75	PRO
1	A	209	ASP
1	B	253	PRO
1	A	76	THR
1	A	138	GLU
1	B	138	GLU
1	B	100	LEU
1	B	186	SER
1	B	179	VAL

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	196/210 (93%)	185 (94%)	11 (6%)	17	26
1	B	197/210 (94%)	185 (94%)	12 (6%)	15	22
All	All	393/420 (94%)	370 (94%)	23 (6%)	16	23

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	GLN
1	A	21	THR
1	A	77	THR
1	A	100	LEU
1	A	115	ASP
1	A	146	CYS
1	A	171	MSE
1	A	183	SER
1	A	204	ARG
1	A	208	ASN
1	A	245	VAL
1	B	12	ILE
1	B	30	THR
1	B	42	LEU
1	B	58	GLN
1	B	77	THR
1	B	97	ASP
1	B	137	ILE
1	B	143	THR
1	B	166	GLU
1	B	171	MSE
1	B	197	VAL
1	B	253	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	ASN
1	A	39	GLN
1	A	58	GLN
1	A	109	GLN
1	A	114	HIS
1	A	208	ASN
1	B	9	GLN
1	B	58	GLN
1	B	164	HIS

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

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	244/276 (88%)	0.51	17 (6%) 24 25	17, 36, 52, 65	0
1	B	245/276 (88%)	0.74	26 (10%) 13 14	20, 38, 64, 74	0
All	All	489/552 (88%)	0.62	43 (8%) 17 19	17, 37, 59, 74	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	7	ALA	5.8
1	A	7	ALA	4.5
1	A	74	SER	4.4
1	B	139	VAL	4.1
1	B	146	CYS	3.9
1	A	111	TRP	3.8
1	A	73	GLY	3.7
1	B	137	ILE	3.5
1	B	193	HIS	3.3
1	B	145	GLY	3.3
1	A	126	PHE	3.2
1	B	136	GLY	3.2
1	B	9	GLN	3.1
1	B	126	PHE	3.1
1	A	222	LYS	3.1
1	B	8	ARG	3.0
1	B	140	ARG	2.9
1	B	188	ARG	2.8
1	A	105	SER	2.7
1	B	111	TRP	2.6
1	B	253	PRO	2.6
1	A	127	THR	2.6
1	B	178	ARG	2.5
1	A	137	ILE	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	115	ASP	2.4
1	A	24	ASN	2.4
1	A	75	PRO	2.3
1	A	32	GLY	2.3
1	A	8	ARG	2.3
1	B	142	ILE	2.2
1	B	184	ALA	2.2
1	B	235	THR	2.2
1	B	194	GLY	2.2
1	B	179	VAL	2.2
1	A	262	PRO	2.2
1	B	135	ALA	2.2
1	B	12	ILE	2.1
1	A	104	VAL	2.1
1	A	193	HIS	2.1
1	B	227	PRO	2.1
1	B	226	VAL	2.0
1	A	203	GLY	2.0
1	B	131	PRO	2.0

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 monosaccharides 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.