



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:15 AM UTC

PDB ID : 1VI6 / pdb_00001vi6
Title : Crystal structure of ribosomal protein S2P
Authors : Structural GenomiX
Deposited on : 2003-12-01
Resolution : 1.95 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
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

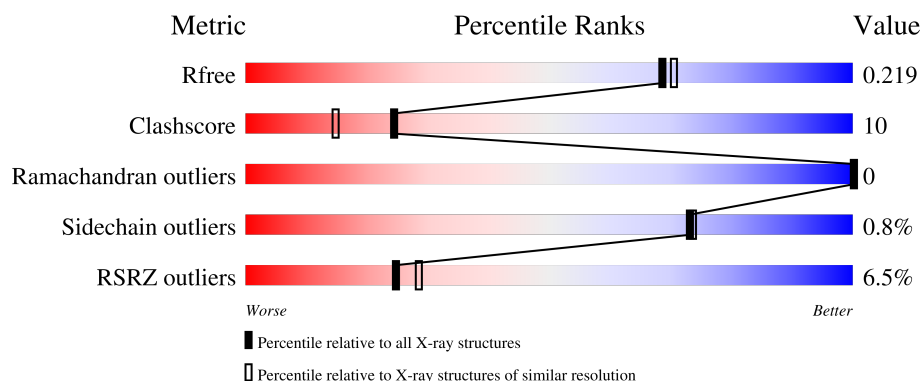
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	180053	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (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 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	208	4% 74% 18% • 7%
1	B	208	7% 76% 16% • 7%
1	C	208	6% 74% 19% 7%
1	D	208	8% 72% 20% 7%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6687 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 30S ribosomal protein S2P.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	193	Total	C	N	O	S	0	3	0
			1534	988	258	284	4			
1	B	193	Total	C	N	O	S	0	3	0
			1524	981	255	284	4			
1	C	194	Total	C	N	O	S	0	3	0
			1537	988	262	283	4			
1	D	193	Total	C	N	O	S	0	2	0
			1524	980	257	283	4			

There are 52 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	cloning artifact	UNP O29132
A	1	SER	-	cloning artifact	UNP O29132
A	2	LEU	-	cloning artifact	UNP O29132
A	198	GLU	-	cloning artifact	UNP O29132
A	199	GLY	-	cloning artifact	UNP O29132
A	200	GLY	-	cloning artifact	UNP O29132
A	201	SER	-	cloning artifact	UNP O29132
A	202	HIS	-	cloning artifact	UNP O29132
A	203	HIS	-	cloning artifact	UNP O29132
A	204	HIS	-	cloning artifact	UNP O29132
A	205	HIS	-	cloning artifact	UNP O29132
A	206	HIS	-	cloning artifact	UNP O29132
A	207	HIS	-	cloning artifact	UNP O29132
B	0	MET	-	cloning artifact	UNP O29132
B	1	SER	-	cloning artifact	UNP O29132
B	2	LEU	-	cloning artifact	UNP O29132
B	198	GLU	-	cloning artifact	UNP O29132
B	199	GLY	-	cloning artifact	UNP O29132
B	200	GLY	-	cloning artifact	UNP O29132
B	201	SER	-	cloning artifact	UNP O29132
B	202	HIS	-	cloning artifact	UNP O29132

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Chain	Residue	Modelled	Actual	Comment	Reference
B	203	HIS	-	cloning artifact	UNP O29132
B	204	HIS	-	cloning artifact	UNP O29132
B	205	HIS	-	cloning artifact	UNP O29132
B	206	HIS	-	cloning artifact	UNP O29132
B	207	HIS	-	cloning artifact	UNP O29132
C	0	MET	-	cloning artifact	UNP O29132
C	1	SER	-	cloning artifact	UNP O29132
C	2	LEU	-	cloning artifact	UNP O29132
C	198	GLU	-	cloning artifact	UNP O29132
C	199	GLY	-	cloning artifact	UNP O29132
C	200	GLY	-	cloning artifact	UNP O29132
C	201	SER	-	cloning artifact	UNP O29132
C	202	HIS	-	cloning artifact	UNP O29132
C	203	HIS	-	cloning artifact	UNP O29132
C	204	HIS	-	cloning artifact	UNP O29132
C	205	HIS	-	cloning artifact	UNP O29132
C	206	HIS	-	cloning artifact	UNP O29132
C	207	HIS	-	cloning artifact	UNP O29132
D	0	MET	-	cloning artifact	UNP O29132
D	1	SER	-	cloning artifact	UNP O29132
D	2	LEU	-	cloning artifact	UNP O29132
D	198	GLU	-	cloning artifact	UNP O29132
D	199	GLY	-	cloning artifact	UNP O29132
D	200	GLY	-	cloning artifact	UNP O29132
D	201	SER	-	cloning artifact	UNP O29132
D	202	HIS	-	cloning artifact	UNP O29132
D	203	HIS	-	cloning artifact	UNP O29132
D	204	HIS	-	cloning artifact	UNP O29132
D	205	HIS	-	cloning artifact	UNP O29132
D	206	HIS	-	cloning artifact	UNP O29132
D	207	HIS	-	cloning artifact	UNP O29132

- Molecule 2 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	3	Total Na 3 3	0	0
2	B	2	Total Na 2 2	0	0
2	C	3	Total Na 3 3	0	0
2	D	2	Total Na 2 2	0	0

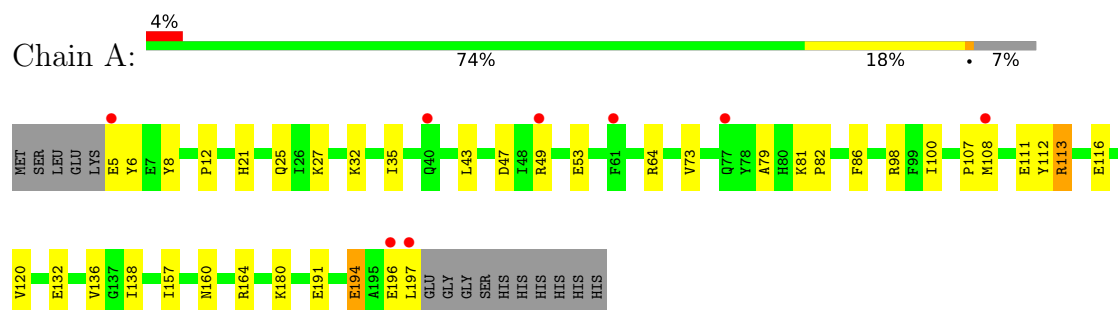
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	156	Total 156	O 156	0	0
3	B	132	Total 132	O 132	0	0
3	C	147	Total 147	O 147	0	0
3	D	123	Total 123	O 123	0	0

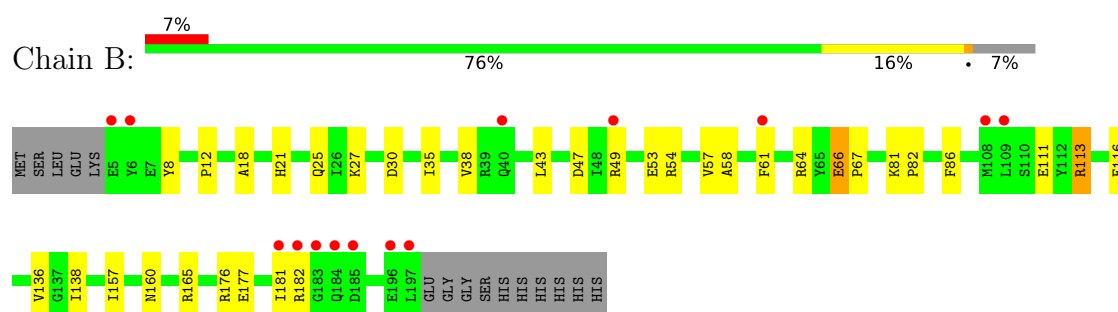
3 Residue-property plots [i](#)

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.

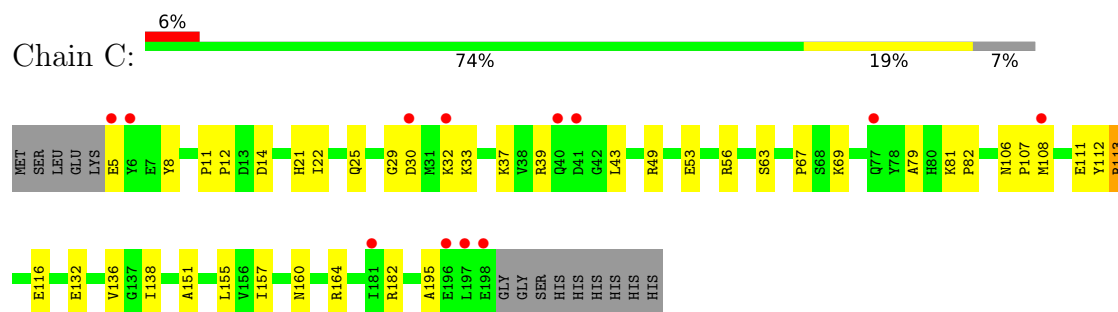
- Molecule 1: 30S ribosomal protein S2P



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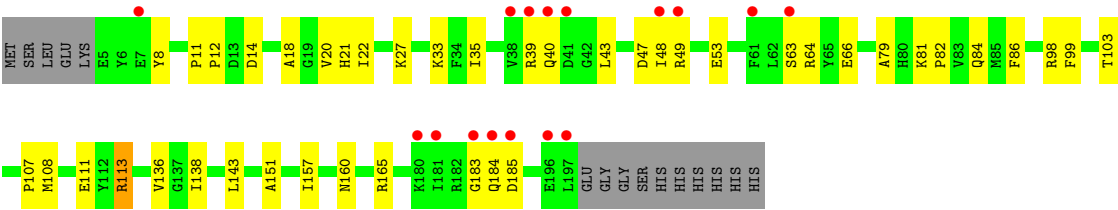


- Molecule 1: 30S ribosomal protein S2P



- Molecule 1: 30S ribosomal protein S2P





4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	123.83Å 123.83Å 131.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.48 – 1.95 36.48 – 1.95	Depositor EDS
% Data completeness (in resolution range)	(Not available) (36.48-1.95) 100.0 (36.48-1.95)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.48 (at 1.95Å)	Xtriage
Refinement program	REFMAC 4.0	Depositor
R, R_{free}	0.208 , 0.243 0.192 , 0.219	Depositor DCC
R_{free} test set	3748 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	21.1	Xtriage
Anisotropy	0.335	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 52.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6687	wwPDB-VP
Average B, all atoms (Å ²)	24.0	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 51.46 % 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 5.5957e-05. 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

5.1 Standard geometry

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

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	0.38	0/1583	0.89	3/2149 (0.1%)
1	B	0.37	0/1571	0.91	3/2134 (0.1%)
1	C	0.35	0/1586	0.88	1/2152 (0.0%)
1	D	0.37	0/1566	0.90	6/2126 (0.3%)
All	All	0.37	0/6306	0.89	13/8561 (0.2%)

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	157	ILE	N-CA-C	-7.65	100.58	108.15
1	D	66	GLU	N-CA-C	-7.24	99.06	109.48
1	A	157	ILE	N-CA-C	-7.23	101.00	108.15
1	B	160	ASN	N-CA-C	-6.60	99.89	109.59
1	D	157	ILE	N-CA-C	-6.17	102.04	108.15
1	A	160	ASN	N-CA-C	-6.11	100.61	109.59
1	D	183	GLY	N-CA-C	-5.66	107.17	115.72
1	D	143	LEU	N-CA-C	-5.19	101.51	109.76
1	C	157	ILE	N-CA-C	-5.16	103.04	108.15
1	D	160	ASN	N-CA-C	-5.16	102.01	109.59
1	D	22	ILE	N-CA-C	5.14	114.67	107.37
1	A	194	GLU	N-CA-C	5.12	117.64	109.96
1	B	66	GLU	N-CA-C	-5.12	100.17	109.44

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	1534	0	1551	38	0
1	B	1524	0	1529	28	0
1	C	1537	0	1548	33	0
1	D	1524	0	1532	34	0
2	A	3	0	0	0	0
2	B	2	0	0	0	0
2	C	3	0	0	0	0
2	D	2	0	0	0	0
3	A	156	0	0	6	0
3	B	132	0	0	4	0
3	C	147	0	0	3	0
3	D	123	0	0	5	0
All	All	6687	0	6160	122	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 10.

All (122) 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:D:63:SER:HB2	3:D:303:HOH:O	1.79	0.82
1:A:98:ARG:NH1	1:A:100[A]:ILE:HD13	1.95	0.80
1:C:29:GLY:HA2	1:C:32:LYS:HE3	1.70	0.73
1:C:106:ASN:OD1	1:C:108:MET:HG2	1.87	0.73
1:C:49:ARG:O	1:C:53:GLU:HG3	1.89	0.73
1:C:155:LEU:HD11	3:D:294:HOH:O	1.92	0.69
1:D:21:HIS:HB3	1:D:48:ILE:HD11	1.74	0.69
1:B:54:ARG:HA	1:B:57[A]:VAL:HG22	1.75	0.69
1:D:184:GLN:HG2	1:D:185:ASP:H	1.58	0.69
1:A:111:GLU:O	1:A:113:ARG:HD2	1.94	0.66
1:A:8:TYR:CD2	1:A:12:PRO:HG3	2.31	0.65
1:A:98:ARG:HH12	1:A:100[A]:ILE:HD13	1.59	0.65
1:B:66:GLU:HG2	3:B:252:HOH:O	1.96	0.63
1:B:81:LYS:HB3	1:B:82:PRO:HD3	1.80	0.61
1:A:81:LYS:HD3	1:D:98:ARG:CZ	2.30	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:49:ARG:O	1:A:53:GLU:HG3	2.01	0.61
1:D:81:LYS:HB3	1:D:82:PRO:HD3	1.82	0.60
1:D:184:GLN:HG2	1:D:185:ASP:N	2.16	0.60
1:C:67:PRO:HG3	1:C:182:ARG:NH2	2.17	0.60
1:D:111:GLU:O	1:D:113:ARG:HD2	2.02	0.60
1:C:30:ASP:OD2	1:D:33:LYS:HE2	2.02	0.59
1:C:111:GLU:O	1:C:113:ARG:HD2	2.03	0.58
1:A:136[B]:VAL:HG13	1:A:138:ILE:HG13	1.84	0.58
1:D:136:VAL:HG23	1:D:138:ILE:HG13	1.86	0.58
1:B:111:GLU:O	1:B:113:ARG:HD2	2.02	0.58
1:A:196:GLU:O	1:A:197:LEU:HB2	2.04	0.57
1:C:151:ALA:HA	1:D:33:LYS:HD2	1.87	0.57
1:C:25:GLN:HG2	3:C:354:HOH:O	2.04	0.57
1:A:100[A]:ILE:HD11	1:D:84:GLN:CD	2.30	0.56
1:D:39:ARG:CZ	1:D:43:LEU:HD23	2.35	0.56
1:A:81:LYS:HB3	1:A:82:PRO:HD3	1.88	0.56
1:B:116[B]:GLU:OE1	3:B:218:HOH:O	2.17	0.56
1:C:69:LYS:HG3	1:C:116[A]:GLU:OE2	2.06	0.55
1:D:8:TYR:CD2	1:D:12:PRO:HG3	2.41	0.55
1:C:136[B]:VAL:HG23	1:C:138:ILE:HG13	1.88	0.55
1:B:8:TYR:CD2	1:B:12:PRO:HG3	2.42	0.55
1:A:116[A]:GLU:OE1	3:A:221:HOH:O	2.18	0.54
1:C:11:PRO:HG2	1:C:14:ASP:OD2	2.06	0.54
1:D:49:ARG:O	1:D:53:GLU:HG3	2.08	0.54
1:B:176:ARG:NE	3:B:278:HOH:O	2.40	0.53
1:B:47:ASP:OD2	1:B:49:ARG:HB3	2.09	0.52
1:C:8:TYR:CD2	1:C:12:PRO:HG3	2.44	0.52
1:A:21:HIS:CD2	1:A:21:HIS:H	2.26	0.52
1:C:33:LYS:HD2	1:D:151:ALA:HA	1.91	0.51
1:B:49:ARG:O	1:B:53:GLU:HG3	2.11	0.51
1:A:98:ARG:CZ	1:D:81:LYS:HD3	2.40	0.51
1:C:132:GLU:O	1:C:136[B]:VAL:HG22	2.10	0.51
1:A:79:ALA:C	1:A:82:PRO:HD2	2.36	0.50
1:C:21:HIS:H	1:C:21:HIS:CD2	2.28	0.50
1:D:21:HIS:H	1:D:21:HIS:CD2	2.28	0.50
1:A:107:PRO:HA	1:A:112:TYR:CD2	2.46	0.50
1:B:49:ARG:NH1	1:B:53:GLU:OE2	2.45	0.50
1:A:164:ARG:NH1	1:A:194:GLU:HG2	2.26	0.50
1:B:21:HIS:H	1:B:21:HIS:CD2	2.30	0.49
1:C:63:SER:HB2	3:C:286:HOH:O	2.12	0.49
1:A:6:TYR:CG	1:A:49:ARG:HD2	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:107:PRO:HA	1:C:112:TYR:CG	2.48	0.49
1:A:180:LYS:NZ	3:A:279:HOH:O	2.46	0.49
1:B:61:PHE:O	1:B:64:ARG:HB2	2.13	0.49
1:C:155:LEU:CD1	3:D:294:HOH:O	2.56	0.49
1:A:107:PRO:C	1:A:108:MET:HE2	2.38	0.48
1:B:25:GLN:HE21	1:B:43:LEU:CD2	2.26	0.48
1:D:18:ALA:HA	1:D:165:ARG:HG2	1.95	0.48
1:D:79:ALA:C	1:D:82:PRO:HD2	2.38	0.48
1:C:116[A]:GLU:OE1	1:D:64:ARG:NH2	2.40	0.48
1:A:27:LYS:HD2	1:A:32:LYS:HG2	1.94	0.47
1:A:132:GLU:O	1:A:136[B]:VAL:HG12	2.14	0.47
1:A:100[A]:ILE:HD11	1:D:84:GLN:OE1	2.15	0.46
1:B:54:ARG:HA	1:B:57[A]:VAL:CG2	2.42	0.46
1:C:30:ASP:OD2	1:D:33:LYS:CE	2.64	0.46
1:C:79:ALA:C	1:C:82:PRO:HD2	2.41	0.46
1:D:11:PRO:HG2	1:D:14:ASP:CG	2.40	0.46
1:D:47:ASP:OD2	1:D:49:ARG:HB3	2.16	0.46
1:D:49:ARG:NH1	1:D:53:GLU:OE2	2.49	0.46
1:D:99:PHE:CZ	1:D:103:THR:HB	2.51	0.46
1:A:21:HIS:HE1	3:A:325:HOH:O	1.98	0.46
1:A:47:ASP:OD2	1:A:49:ARG:HB3	2.16	0.46
1:C:81:LYS:HB3	1:C:82:PRO:HD3	1.97	0.46
1:B:136:VAL:HG23	1:B:138:ILE:HG13	1.99	0.45
1:A:27:LYS:HG3	1:A:35:ILE:HD12	1.97	0.45
1:A:107:PRO:HA	1:A:112:TYR:CG	2.52	0.45
1:B:30:ASP:OD2	3:B:238:HOH:O	2.21	0.45
1:C:5:GLU:HG2	1:C:5:GLU:O	2.16	0.45
1:D:20:VAL:HG23	1:D:48:ILE:CD1	2.47	0.45
1:C:106:ASN:HD21	1:C:108:MET:HE2	1.81	0.45
1:B:177:GLU:O	1:B:181:ILE:HG13	2.17	0.44
1:C:39[A]:ARG:CZ	1:C:43:LEU:HD23	2.46	0.44
1:B:67:PRO:HG3	1:B:182:ARG:NH2	2.32	0.44
1:B:49:ARG:HH11	1:B:49:ARG:HG3	1.82	0.44
1:B:54:ARG:CA	1:B:57[A]:VAL:HG22	2.46	0.44
1:D:81:LYS:NZ	3:D:305:HOH:O	2.48	0.44
1:A:191:GLU:HG2	3:A:353:HOH:O	2.18	0.43
1:C:164:ARG:HD2	1:C:195:ALA:O	2.18	0.43
1:A:180:LYS:HG3	3:A:267:HOH:O	2.17	0.43
1:C:37:LYS:HB2	3:C:328:HOH:O	2.18	0.43
1:A:86:PHE:CD1	1:A:86:PHE:C	2.97	0.43
1:C:151:ALA:HB2	1:D:33:LYS:HB2	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:18:ALA:HA	1:D:165:ARG:CG	2.49	0.42
1:B:49:ARG:HH11	1:B:49:ARG:CG	2.32	0.42
1:B:86:PHE:CD1	1:B:86:PHE:C	2.96	0.42
1:D:86:PHE:C	1:D:86:PHE:CD1	2.97	0.42
1:D:107:PRO:C	1:D:108:MET:HE2	2.43	0.42
1:A:27:LYS:HG3	1:A:35:ILE:CD1	2.49	0.42
1:A:25:GLN:HE21	1:A:43:LEU:CD2	2.33	0.42
1:B:38:VAL:O	1:B:38:VAL:HG13	2.20	0.41
1:A:64:ARG:HH22	1:B:116[B]:GLU:CD	2.28	0.41
1:B:18:ALA:HA	1:B:165:ARG:HD3	2.02	0.41
1:D:27:LYS:HG3	1:D:35:ILE:CD1	2.49	0.41
1:A:73:VAL:HB	1:A:120:VAL:HG12	2.02	0.41
1:C:56:ARG:HB3	3:D:280:HOH:O	2.20	0.41
1:C:107:PRO:HA	1:C:112:TYR:CD2	2.56	0.41
1:B:54:ARG:O	1:B:57[A]:VAL:HG22	2.21	0.41
1:D:11:PRO:HA	1:D:12:PRO:HD3	1.96	0.41
1:B:27:LYS:HG3	1:B:35:ILE:HD12	2.01	0.41
1:A:5:GLU:HG3	1:A:5:GLU:O	2.21	0.41
1:A:81:LYS:NZ	3:A:242:HOH:O	2.46	0.40
1:B:57[A]:VAL:HG23	1:B:58:ALA:N	2.37	0.40
1:A:196:GLU:O	1:A:197:LEU:CB	2.69	0.40
1:C:22:ILE:O	1:C:160:ASN:HB2	2.21	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	194/208 (93%)	188 (97%)	6 (3%)	0	100	100
1	B	194/208 (93%)	189 (97%)	5 (3%)	0	100	100
1	C	195/208 (94%)	188 (96%)	7 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	193/208 (93%)	188 (97%)	5 (3%)	0	100	100
All	All	776/832 (93%)	753 (97%)	23 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	167/178 (94%)	166 (99%)	1 (1%)	78	79
1	B	164/178 (92%)	163 (99%)	1 (1%)	78	79
1	C	166/178 (93%)	165 (99%)	1 (1%)	78	79
1	D	164/178 (92%)	162 (99%)	2 (1%)	63	61
All	All	661/712 (93%)	656 (99%)	5 (1%)	73	74

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	113	ARG
1	B	113	ARG
1	C	113	ARG
1	D	40	GLN
1	D	113	ARG

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	21	HIS
1	A	25	GLN
1	A	77	GLN
1	B	21	HIS
1	B	25	GLN
1	B	77	GLN

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Mol	Chain	Res	Type
1	C	21	HIS
1	C	84	GLN
1	D	21	HIS
1	D	40	GLN
1	D	77	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](#)

Of 10 ligands modelled in this entry, 10 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	193/208 (92%)	0.10	8 (4%)	41 48	11, 19, 39, 59	3 (1%)
1	B	193/208 (92%)	0.18	14 (7%)	21 24	12, 20, 38, 61	3 (1%)
1	C	194/208 (93%)	0.33	12 (6%)	26 31	13, 22, 38, 69	3 (1%)
1	D	193/208 (92%)	0.31	16 (8%)	17 20	13, 22, 42, 59	2 (1%)
All	All	773/832 (92%)	0.23	50 (6%)	25 28	11, 21, 40, 69	11 (1%)

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	197	LEU	6.1
1	A	197	LEU	5.7
1	B	185	ASP	5.7
1	D	197	LEU	5.4
1	B	197	LEU	5.4
1	D	185	ASP	4.8
1	C	198	GLU	4.0
1	B	183	GLY	3.8
1	D	183	GLY	3.8
1	B	5	GLU	3.6
1	D	38	VAL	3.6
1	A	5	GLU	3.3
1	C	6	TYR	3.1
1	D	39	ARG	3.1
1	C	32	LYS	3.1
1	B	40	GLN	3.0
1	C	5	GLU	3.0
1	C	181	ILE	2.8
1	C	196	GLU	2.8
1	C	40	GLN	2.8
1	C	77	GLN	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	49	ARG	2.6
1	B	61	PHE	2.6
1	D	61	PHE	2.6
1	D	196	GLU	2.6
1	D	184	GLN	2.5
1	D	41	ASP	2.5
1	C	108	MET	2.5
1	C	30	ASP	2.5
1	C	41	ASP	2.5
1	B	109	LEU	2.4
1	D	40	GLN	2.4
1	A	196	GLU	2.4
1	A	40	GLN	2.4
1	B	196	GLU	2.3
1	A	77	GLN	2.3
1	D	181	ILE	2.3
1	D	180	LYS	2.2
1	D	7	GLU	2.2
1	D	63	SER	2.2
1	B	6	TYR	2.2
1	A	108	MET	2.2
1	B	108	MET	2.2
1	B	184	GLN	2.2
1	D	49	ARG	2.2
1	B	49	ARG	2.2
1	B	181	ILE	2.1
1	D	48	ILE	2.1
1	A	61	PHE	2.0
1	B	182	ARG	2.0

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 oligosaccharides in this entry.

6.4 Ligands [i](#)

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. The B-factors column lists the minimum, median, 95th 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	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NA	A	210	1/1	0.93	0.12	42,42,42,42	0
2	NA	B	209	1/1	0.93	0.10	29,29,29,29	0
2	NA	D	208	1/1	0.94	0.07	21,21,21,21	0
2	NA	C	210	1/1	0.96	0.07	42,42,42,42	0
2	NA	C	209	1/1	0.96	0.05	24,24,24,24	0
2	NA	A	209	1/1	0.98	0.05	23,23,23,23	0
2	NA	B	208	1/1	0.98	0.06	19,19,19,19	0
2	NA	C	208	1/1	0.99	0.05	22,22,22,22	0
2	NA	A	208	1/1	0.99	0.03	19,19,19,19	0
2	NA	D	209	1/1	0.99	0.10	24,24,24,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.