



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 09:26 PM UTC

PDB ID : 4PUF / pdb_00004puf
Title : Complex between the Salmonella T3SS effector SlrP and its human target thioredoxin-1
Authors : Zouhir, S.; Bernal-Bayard, J.; Cordero-Alba, M.; Cardenal-Munoz, E.; Guimaraes, B.; Lazar, N.; Ramos-Morales, F.; Nessler, S.
Deposited on : 2014-03-13
Resolution : 3.30 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	:	4-5-2 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

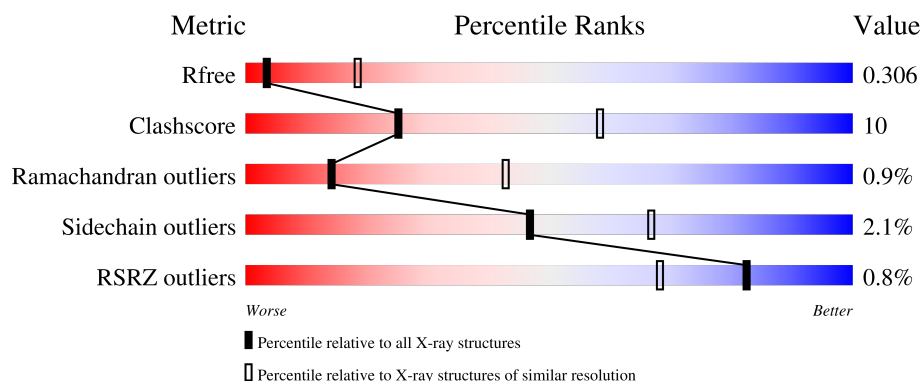
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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}	180053	1169 (3.32-3.28)
Clashscore	190562	1209 (3.32-3.28)
Ramachandran outliers	187476	1188 (3.32-3.28)
Sidechain outliers	187428	1187 (3.32-3.28)
RSRZ outliers	180081	1169 (3.32-3.28)

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	637	
1	B	637	
2	C	117	
2	D	117	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11382 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 E3 ubiquitin-protein ligase SlrP.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	599	Total	C	N	O	S	0	0	0
			4815	3039	826	928	22			
1	B	598	Total	C	N	O	S	0	0	0
			4807	3033	825	927	22			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	129	MET	-	expression tag	UNP D0ZRB2
A	130	ARG	-	expression tag	UNP D0ZRB2
A	131	GLY	-	expression tag	UNP D0ZRB2
A	132	SER	-	expression tag	UNP D0ZRB2
A	133	HIS	-	expression tag	UNP D0ZRB2
A	134	HIS	-	expression tag	UNP D0ZRB2
A	135	HIS	-	expression tag	UNP D0ZRB2
A	136	HIS	-	expression tag	UNP D0ZRB2
A	137	HIS	-	expression tag	UNP D0ZRB2
A	138	HIS	-	expression tag	UNP D0ZRB2
A	139	GLY	-	expression tag	UNP D0ZRB2
A	140	SER	-	expression tag	UNP D0ZRB2
B	129	MET	-	expression tag	UNP D0ZRB2
B	130	ARG	-	expression tag	UNP D0ZRB2
B	131	GLY	-	expression tag	UNP D0ZRB2
B	132	SER	-	expression tag	UNP D0ZRB2
B	133	HIS	-	expression tag	UNP D0ZRB2
B	134	HIS	-	expression tag	UNP D0ZRB2
B	135	HIS	-	expression tag	UNP D0ZRB2
B	136	HIS	-	expression tag	UNP D0ZRB2
B	137	HIS	-	expression tag	UNP D0ZRB2
B	138	HIS	-	expression tag	UNP D0ZRB2
B	139	GLY	-	expression tag	UNP D0ZRB2
B	140	SER	-	expression tag	UNP D0ZRB2

- Molecule 2 is a protein called Thioredoxin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	112	Total	C	N	O	S	0	0	0
			880	560	145	167	8			
2	D	112	Total	C	N	O	S	0	0	0
			880	560	145	167	8			

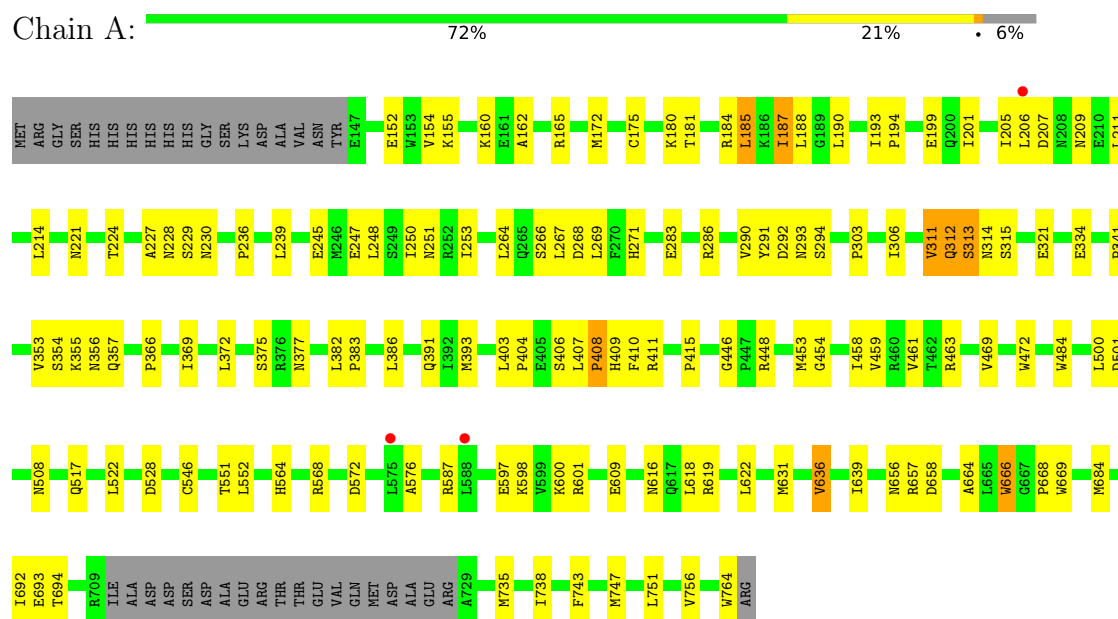
There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-11	MET	-	expression tag	UNP P10599
C	-10	ARG	-	expression tag	UNP P10599
C	-9	GLY	-	expression tag	UNP P10599
C	-8	SER	-	expression tag	UNP P10599
C	-7	HIS	-	expression tag	UNP P10599
C	-6	HIS	-	expression tag	UNP P10599
C	-5	HIS	-	expression tag	UNP P10599
C	-4	HIS	-	expression tag	UNP P10599
C	-3	HIS	-	expression tag	UNP P10599
C	-2	HIS	-	expression tag	UNP P10599
C	-1	GLY	-	expression tag	UNP P10599
C	0	SER	-	expression tag	UNP P10599
D	-11	MET	-	expression tag	UNP P10599
D	-10	ARG	-	expression tag	UNP P10599
D	-9	GLY	-	expression tag	UNP P10599
D	-8	SER	-	expression tag	UNP P10599
D	-7	HIS	-	expression tag	UNP P10599
D	-6	HIS	-	expression tag	UNP P10599
D	-5	HIS	-	expression tag	UNP P10599
D	-4	HIS	-	expression tag	UNP P10599
D	-3	HIS	-	expression tag	UNP P10599
D	-2	HIS	-	expression tag	UNP P10599
D	-1	GLY	-	expression tag	UNP P10599
D	0	SER	-	expression tag	UNP P10599

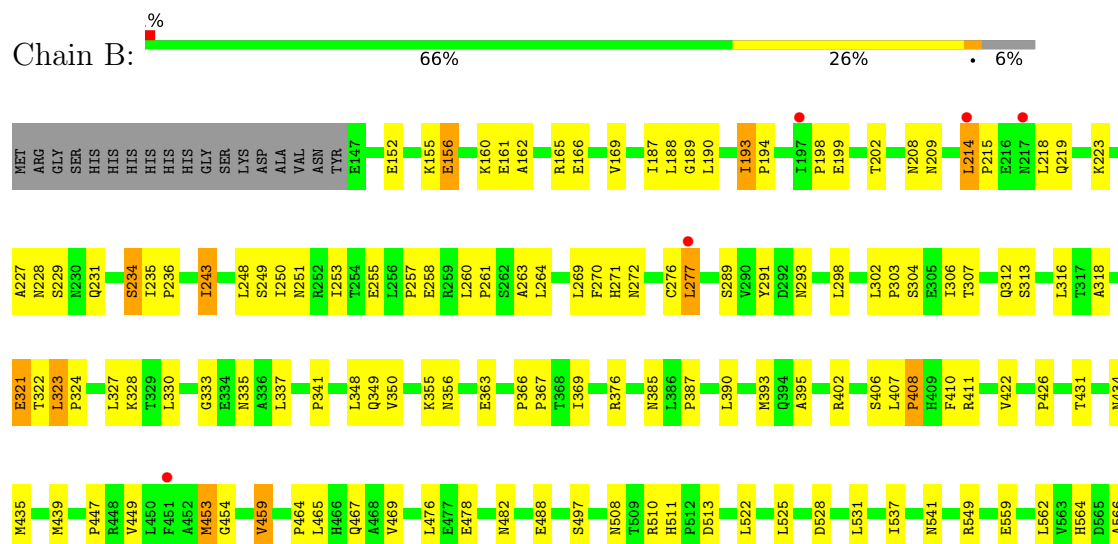
3 Residue-property plots

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: E3 ubiquitin-protein ligase SlrP

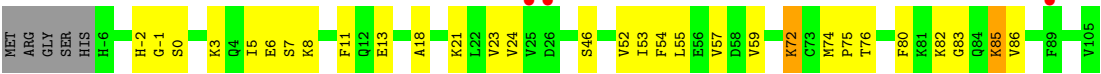


• Molecule 1: E3 ubiquitin-protein ligase SlrP

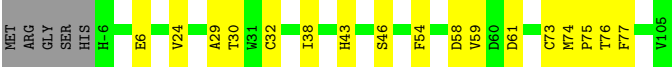
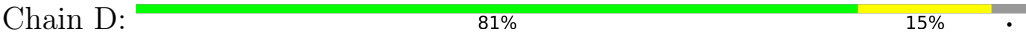




● Molecule 2: Thioredoxin



● Molecule 2: Thioredoxin



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	106.29Å 134.83Å 154.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.85 – 3.30 43.85 – 3.30	Depositor EDS
% Data completeness (in resolution range)	99.3 (43.85-3.30) 99.4 (43.85-3.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	5.60	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.57 (at 3.32Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, R_{free}	0.272 , 0.308 0.279 , 0.306	Depositor DCC
R_{free} test set	1702 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	116.0	Xtriage
Anisotropy	0.552	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 54.4	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	0.93	EDS
Total number of atoms	11382	wwPDB-VP
Average B, all atoms (Å ²)	118.0	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 20.88 % 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 7.9028e-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

5.1 Standard geometry

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	0.29	0/4908	0.73	2/6667 (0.0%)
1	B	0.33	0/4900	0.86	18/6656 (0.3%)
2	C	0.34	0/900	0.79	1/1210 (0.1%)
2	D	0.25	0/900	0.65	0/1210
All	All	0.31	0/11608	0.79	21/15743 (0.1%)

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	323	LEU	CA-C-N	-13.85	106.11	120.38
1	B	323	LEU	C-N-CA	-13.85	106.11	120.38
1	B	214	LEU	CA-C-N	6.28	126.24	119.78
1	B	214	LEU	C-N-CA	6.28	126.24	119.78
2	C	85	LYS	N-CA-C	6.22	117.73	111.07
1	B	255	GLU	N-CA-C	6.12	117.18	108.74
1	B	764	TRP	CB-CA-C	-5.93	109.75	116.63
1	B	189	GLY	N-CA-C	-5.72	104.98	114.48
1	B	762	ALA	N-CA-C	-5.60	100.79	109.14
1	B	407	LEU	CA-C-N	5.55	126.78	119.84
1	B	407	LEU	C-N-CA	5.55	126.78	119.84
1	B	277	LEU	CA-C-N	5.46	125.47	119.90
1	B	277	LEU	C-N-CA	5.46	125.47	119.90
1	A	315	SER	N-CA-C	5.24	118.94	111.92
1	A	546	CYS	CB-CA-C	-5.17	110.63	116.63
1	B	260	LEU	CA-C-N	-5.17	114.34	119.56
1	B	260	LEU	C-N-CA	-5.17	114.34	119.56
1	B	156	GLU	N-CA-C	-5.16	106.99	113.28
1	B	243	ILE	CB-CA-C	-5.02	106.48	111.80
1	B	193	ILE	CA-C-N	5.01	125.01	119.90
1	B	193	ILE	C-N-CA	5.01	125.01	119.90

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	4815	0	4794	85	1
1	B	4807	0	4783	110	0
2	C	880	0	854	20	0
2	D	880	0	854	13	0
All	All	11382	0	11285	220	1

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 (220) 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:293:ASN:H	1:A:313:SER:HB2	1.40	0.85
1:B:235:ILE:HB	1:B:257:PRO:HG2	1.58	0.83
1:A:160:LYS:HZ1	1:A:190:LEU:HA	1.48	0.79
1:B:323:LEU:HG	1:B:324:PRO:HD3	1.64	0.76
1:B:522:LEU:HA	1:B:525:LEU:HD12	1.70	0.74
1:B:152:GLU:HA	1:B:155:LYS:HB2	1.69	0.73
1:B:276:CYS:SG	1:B:277:LEU:N	2.62	0.72
1:A:193:ILE:HD13	1:A:206:LEU:HD21	1.73	0.71
1:A:283:GLU:HA	1:A:303:PRO:HB3	1.73	0.70
1:B:528:ASP:OD1	1:B:531:LEU:N	2.19	0.70
1:A:190:LEU:H	1:A:209:ASN:HD22	1.40	0.70
2:C:82:LYS:HG3	2:C:83:GLY:H	1.57	0.69
1:B:465:LEU:HD13	1:B:488:GLU:HG3	1.74	0.68
1:A:206:LEU:O	1:A:227:ALA:HA	1.93	0.68
1:B:669:TRP:O	1:B:673:LEU:N	2.27	0.66
1:A:313:SER:OG	1:A:314:ASN:N	2.29	0.65
1:B:598:LYS:HE3	1:B:601:ARG:HH21	1.62	0.65
1:B:410:PHE:HB3	1:B:447:PRO:HD3	1.79	0.64
1:A:408:PRO:O	1:A:410:PHE:N	2.27	0.64
1:A:185:LEU:HB2	1:A:188:LEU:HD11	1.77	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:508:ASN:ND2	1:B:637:SER:OG	2.30	0.64
1:A:205:ILE:HD11	2:D:30:THR:HG21	1.79	0.64
1:A:598:LYS:HE3	1:A:601:ARG:HH21	1.62	0.64
1:B:328:LYS:HA	1:B:348:LEU:HA	1.80	0.64
1:B:618:LEU:HB3	1:B:622:LEU:HD13	1.79	0.64
1:B:214:LEU:HD22	1:B:236:PRO:HD2	1.80	0.63
1:B:160:LYS:H	1:B:165:ARG:HH22	1.46	0.63
1:B:188:LEU:HB2	1:B:190:LEU:HB2	1.80	0.63
1:A:372:LEU:HD23	1:A:393:MET:HE2	1.81	0.62
1:A:207:ASP:OD1	1:A:228:ASN:N	2.30	0.62
1:A:459:VAL:HG21	2:C:59:VAL:HG11	1.82	0.62
1:A:286:ARG:HA	1:A:306:ILE:HA	1.82	0.61
1:A:386:LEU:HB2	1:A:410:PHE:HE1	1.65	0.61
1:B:243:ILE:O	1:B:264:LEU:HD12	2.01	0.61
1:B:758:SER:O	1:B:761:SER:HB3	2.01	0.60
1:B:202:THR:HG23	1:B:223:LYS:HB2	1.84	0.59
1:B:537:ILE:O	1:B:541:ASN:ND2	2.32	0.59
1:A:187:ILE:HG12	2:D:6:GLU:HB3	1.83	0.59
2:C:80:PHE:HE1	2:C:85:LYS:HA	1.67	0.58
1:A:656:ASN:OD1	1:A:764:TRP:HB2	2.03	0.58
1:A:375:SER:O	1:A:377:ASN:ND2	2.37	0.58
1:B:478:GLU:O	1:B:482:ASN:ND2	2.36	0.58
1:A:391:GLN:HE22	1:A:415:PRO:HD2	1.67	0.58
1:B:253:ILE:HG22	1:B:272:ASN:HB2	1.84	0.58
1:B:511:HIS:ND1	1:B:513:ASP:OD1	2.36	0.58
1:A:616:ASN:OD1	1:A:619:ARG:NH1	2.36	0.58
1:A:264:LEU:HD21	1:A:267:LEU:HB2	1.86	0.58
1:B:234:SER:HA	1:B:253:ILE:HG13	1.85	0.58
1:B:406:SER:C	1:B:408:PRO:HD3	2.29	0.58
1:B:694:THR:HB	1:B:697:TYR:HB2	1.85	0.58
1:B:228:ASN:OD1	1:B:229:SER:N	2.37	0.57
1:B:333:GLY:O	1:B:335:ASN:ND2	2.37	0.57
2:C:6:GLU:N	2:C:6:GLU:OE1	2.38	0.57
2:D:29:ALA:HB3	2:D:74:MET:HE1	1.85	0.57
1:A:406:SER:C	1:A:408:PRO:HD3	2.30	0.57
1:B:402:ARG:HA	1:B:431:THR:HG21	1.87	0.57
1:A:403:LEU:HD13	1:A:407:LEU:HD12	1.87	0.57
1:B:465:LEU:HD22	1:B:488:GLU:HA	1.86	0.57
2:C:5:ILE:HG12	2:C:11:PHE:HD1	1.70	0.57
1:B:187:ILE:HA	1:B:208:ASN:HD22	1.70	0.56
1:A:230:ASN:OD1	1:A:251:ASN:ND2	2.38	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:231:GLN:OE1	1:B:231:GLN:N	2.39	0.56
1:A:597:GLU:HA	1:A:600:LYS:HE3	1.86	0.56
1:A:564:HIS:HB3	1:A:568:ARG:HH12	1.69	0.56
1:B:363:GLU:OE2	1:B:385:ASN:ND2	2.37	0.56
1:A:587:ARG:NH1	1:A:622:LEU:O	2.38	0.56
1:A:386:LEU:HD22	1:A:393:MET:HE1	1.88	0.55
1:B:248:LEU:HB3	1:B:251:ASN:HD21	1.71	0.55
1:B:459:VAL:HG21	2:D:59:VAL:HG11	1.88	0.55
1:B:304:SER:HA	1:B:324:PRO:HB3	1.89	0.55
1:A:664:ALA:HA	1:A:669:TRP:HZ3	1.72	0.55
1:B:566:ALA:HB1	1:B:672:VAL:HG23	1.89	0.55
2:D:43:HIS:O	2:D:46:SER:OG	2.21	0.54
1:B:270:PHE:HD1	1:B:291:TYR:HB3	1.73	0.54
1:B:387:PRO:HD2	1:B:390:LEU:HD12	1.89	0.54
1:B:439:MET:HE3	1:B:449:VAL:H	1.73	0.54
1:A:517:GLN:NE2	1:A:631:MET:SD	2.81	0.54
1:B:691:CYS:O	1:B:694:THR:OG1	2.22	0.54
1:A:386:LEU:HB2	1:A:410:PHE:CE1	2.42	0.53
1:B:323:LEU:HG	1:B:324:PRO:CD	2.34	0.53
1:B:214:LEU:HB2	1:B:236:PRO:HG2	1.91	0.53
1:B:751:LEU:HD21	1:B:760:MET:HG2	1.90	0.53
1:A:224:THR:HG22	1:A:245:GLU:HB2	1.91	0.53
1:B:594:LEU:HD21	1:B:650:ARG:HD2	1.91	0.53
1:B:349:GLN:HA	1:B:369:ILE:HA	1.91	0.53
1:A:472:TRP:HE1	1:A:500:LEU:HB3	1.74	0.53
2:D:73:CYS:O	2:D:76:THR:OG1	2.24	0.53
1:A:528:ASP:OD1	1:A:568:ARG:NH2	2.43	0.52
1:A:618:LEU:HB3	1:A:622:LEU:HD13	1.91	0.52
1:B:390:LEU:HD21	1:B:393:MET:HB2	1.91	0.52
1:B:431:THR:HG22	1:B:604:PHE:HB3	1.91	0.52
1:A:657:ARG:NH1	1:A:658:ASP:OD1	2.43	0.51
1:B:754:LYS:HB3	1:B:756:VAL:HG22	1.91	0.51
1:A:175:CYS:HA	1:A:180:LYS:HG2	1.92	0.51
1:B:349:GLN:HG3	1:B:350:VAL:HG23	1.92	0.51
1:B:673:LEU:HD11	1:B:750:ILE:HD12	1.92	0.51
1:A:410:PHE:O	1:A:411:ARG:HG3	2.11	0.51
1:B:559:GLU:HG3	1:B:585:ILE:HD13	1.93	0.51
2:C:46:SER:HA	2:C:54:PHE:CE2	2.45	0.50
1:B:190:LEU:HD23	1:B:209:ASN:HB3	1.94	0.50
1:B:583:ARG:HB3	1:B:655:GLU:OE2	2.11	0.50
1:A:664:ALA:HB2	1:A:747:MET:CE	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:636:VAL:HG12	1:A:639:ILE:HB	1.95	0.49
2:C:5:ILE:HG21	2:C:57:VAL:HG22	1.93	0.49
1:A:155:LYS:HE2	1:A:165:ARG:HH21	1.77	0.49
2:C:23:VAL:HG12	2:C:53:ILE:HB	1.94	0.49
1:A:229:SER:HA	1:A:250:ILE:HB	1.94	0.49
2:C:6:GLU:HG2	2:C:7:SER:H	1.77	0.49
1:A:165:ARG:NH2	1:A:194:PRO:HD3	2.28	0.49
1:A:184:ARG:HH21	2:D:61:ASP:CG	2.20	0.48
1:A:508:ASN:HB2	1:A:609:GLU:HG2	1.94	0.48
1:B:656:ASN:ND2	1:B:763:TYR:O	2.46	0.48
1:B:575:LEU:HD12	1:B:756:VAL:HG21	1.96	0.48
1:A:236:PRO:HG2	1:A:239:LEU:HG	1.95	0.48
1:A:407:LEU:N	1:A:408:PRO:HD3	2.28	0.48
1:B:653:MET:O	1:B:657:ARG:HG2	2.14	0.48
1:A:291:TYR:HA	1:A:312:GLN:O	2.13	0.48
1:B:223:LYS:HA	1:B:243:ILE:HA	1.96	0.47
1:B:218:LEU:HD12	1:B:219:GLN:N	2.30	0.47
1:B:193:ILE:HG23	1:B:215:PRO:HG3	1.96	0.47
1:A:459:VAL:HG12	2:C:72:LYS:O	2.15	0.47
1:B:248:LEU:HB3	1:B:251:ASN:ND2	2.29	0.47
1:B:227:ALA:HB3	1:B:248:LEU:HD23	1.97	0.47
1:B:229:SER:H	1:B:250:ILE:HB	1.79	0.47
1:B:743:PHE:HA	1:B:746:ILE:HG22	1.97	0.46
1:A:199:GLU:HG3	1:A:221:ASN:HB2	1.98	0.46
1:B:277:LEU:HD12	1:B:298:LEU:HD22	1.98	0.46
1:A:576:ALA:HB2	1:A:756:VAL:HG12	1.98	0.46
1:B:465:LEU:HG	1:B:497:SER:HB2	1.97	0.46
1:A:453:MET:N	1:A:454:GLY:HA3	2.30	0.46
1:B:564:HIS:O	1:B:568:ARG:N	2.47	0.46
1:B:583:ARG:O	1:B:587:ARG:HG3	2.16	0.46
1:A:383:PRO:HG2	1:A:386:LEU:HD23	1.98	0.46
1:A:463:ARG:NE	1:A:501:ASP:OD2	2.48	0.46
1:A:160:LYS:HE3	1:A:162:ALA:HB3	1.98	0.45
1:A:334:GLU:HG2	1:A:355:LYS:HD2	1.98	0.45
1:B:562:LEU:HD12	1:B:585:ILE:HD11	1.98	0.45
1:A:227:ALA:O	1:A:248:LEU:HA	2.15	0.45
1:B:291:TYR:CE2	1:B:312:GLN:HG3	2.51	0.45
1:A:184:ARG:HH22	2:D:58:ASP:CG	2.25	0.45
1:A:735:MET:O	1:A:738:ILE:HG13	2.17	0.45
1:B:307:THR:HA	1:B:327:LEU:HA	1.98	0.45
1:A:564:HIS:HB3	1:A:568:ARG:NH1	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:52:VAL:HG23	2:C:54:PHE:CE2	2.52	0.44
1:A:248:LEU:HB2	1:A:269:LEU:HD23	1.99	0.44
1:B:327:LEU:HD21	1:B:330:LEU:HB2	2.00	0.44
1:B:694:THR:HG22	1:B:696:GLU:H	1.83	0.44
1:B:161:GLU:OE1	1:B:162:ALA:N	2.50	0.44
1:B:549:ARG:NH2	1:B:592:GLU:OE2	2.51	0.44
2:C:3:LYS:HB3	2:C:55:LEU:HD23	2.00	0.44
1:A:354:SER:O	1:A:356:ASN:ND2	2.51	0.44
1:B:355:LYS:HG2	1:B:376:ARG:HD2	2.00	0.44
2:D:32:CYS:HB2	2:D:74:MET:HE3	2.00	0.44
1:A:248:LEU:HD22	1:A:253:ILE:HD11	1.99	0.43
1:B:261:PRO:C	1:B:263:ALA:H	2.26	0.43
1:B:585:ILE:HD12	1:B:666:TRP:CZ2	2.53	0.43
2:D:46:SER:HA	2:D:54:PHE:CE2	2.53	0.43
1:B:152:GLU:O	1:B:156:GLU:N	2.51	0.43
1:B:258:GLU:O	1:B:258:GLU:HG3	2.18	0.43
1:B:469:VAL:HG13	1:B:522:LEU:HD23	1.99	0.43
1:B:249:SER:HB3	1:B:270:PHE:HD2	1.84	0.43
2:C:80:PHE:CE1	2:C:85:LYS:HA	2.52	0.43
1:A:366:PRO:HB2	1:A:369:ILE:HG13	2.00	0.43
1:A:693:GLU:HG3	1:A:694:THR:HG22	2.00	0.43
1:B:303:PRO:HB2	1:B:306:ILE:HG13	2.00	0.43
1:B:582:GLY:HA2	1:B:666:TRP:CH2	2.54	0.43
1:A:684:MET:HE1	1:A:743:PHE:CD1	2.54	0.43
1:B:508:ASN:HB3	1:B:633:PHE:CD1	2.53	0.43
1:B:763:TYR:CG	1:B:764:TRP:N	2.87	0.43
2:C:74:MET:HG3	2:C:75:PRO:HA	2.00	0.43
1:B:579:ILE:O	1:B:583:ARG:HG3	2.18	0.43
2:C:21:LYS:NZ	2:C:53:ILE:HD11	2.34	0.43
1:B:152:GLU:C	1:B:155:LYS:H	2.26	0.42
1:B:271:HIS:HE1	2:C:-2:HIS:HB2	1.84	0.42
1:B:395:ALA:HB3	1:B:422:VAL:HG12	2.01	0.42
2:D:38:ILE:HD13	2:D:77:PHE:CZ	2.54	0.42
1:A:271:HIS:HA	1:A:292:ASP:HB3	2.00	0.42
1:A:321:GLU:HG2	1:A:341:PRO:HB3	2.01	0.42
2:C:8:LYS:O	2:C:11:PHE:HB3	2.19	0.42
1:A:469:VAL:HG13	1:A:522:LEU:HD13	2.00	0.42
1:B:302:LEU:HD12	1:B:323:LEU:HD11	2.00	0.42
1:A:187:ILE:HD12	1:A:187:ILE:HA	1.92	0.42
1:A:551:THR:OG1	1:A:552:LEU:N	2.52	0.42
1:A:245:GLU:HG2	1:A:266:SER:HB3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:229:SER:HA	1:B:250:ILE:HB	2.02	0.42
1:B:434:ASN:ND2	1:B:603:PHE:O	2.53	0.42
1:B:464:PRO:HG2	1:B:467:GLN:HB2	2.00	0.42
1:B:586:PHE:O	1:B:590:GLN:HG2	2.19	0.42
2:C:6:GLU:CG	2:C:7:SER:H	2.32	0.42
1:A:446:GLY:O	1:A:448:ARG:NH1	2.52	0.42
1:B:193:ILE:HA	1:B:194:PRO:HD2	1.88	0.42
1:B:293:ASN:H	1:B:313:SER:HA	1.85	0.42
2:D:6:GLU:N	2:D:6:GLU:OE1	2.53	0.42
1:A:666:TRP:CD2	1:A:668:PRO:HD2	2.55	0.42
1:A:290:VAL:HG23	1:A:311:VAL:HA	2.02	0.42
1:A:188:LEU:HB3	1:A:190:LEU:HD13	2.00	0.41
1:B:453:MET:HA	1:B:454:GLY:HA2	1.51	0.41
1:A:403:LEU:HB2	1:A:404:PRO:HD2	2.01	0.41
1:B:439:MET:CE	1:B:449:VAL:H	2.32	0.41
2:D:74:MET:HA	2:D:75:PRO:C	2.45	0.41
1:A:247:GLU:HA	1:A:268:ASP:HB3	2.01	0.41
1:B:328:LYS:HB3	1:B:349:GLN:HG2	2.01	0.41
1:B:435:MET:HE3	1:B:449:VAL:HG21	2.01	0.41
1:A:409:HIS:O	1:A:446:GLY:HA2	2.21	0.41
1:B:166:GLU:HA	1:B:169:VAL:HG22	2.02	0.41
2:C:6:GLU:HG2	2:C:7:SER:N	2.36	0.41
1:A:469:VAL:HG11	1:A:484:TRP:CE2	2.56	0.41
1:A:657:ARG:NH2	1:A:658:ASP:OD2	2.54	0.41
1:B:366:PRO:HA	1:B:367:PRO:HD3	1.89	0.41
1:B:321:GLU:HA	1:B:341:PRO:HB3	2.03	0.41
1:B:666:TRP:HA	1:B:667:GLY:HA3	1.86	0.41
1:A:152:GLU:OE1	1:A:172:MET:HG2	2.21	0.41
1:A:206:LEU:HD23	1:A:211:LEU:HD21	2.03	0.41
1:B:636:VAL:HG12	1:B:639:ILE:HB	2.03	0.41
1:B:269:LEU:N	1:B:289:SER:O	2.44	0.41
1:B:337:LEU:H	1:B:356:ASN:HB2	1.86	0.40
2:C:18:ALA:HA	2:C:21:LYS:NZ	2.36	0.40
1:B:588:LEU:HD23	1:B:591:ILE:HD12	2.03	0.40
1:A:458:ILE:HD12	1:A:458:ILE:H	1.85	0.40
1:B:426:PRO:HB2	1:B:510:ARG:HH12	1.86	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:357:GLN:NE2	1:A:572:ASP:OD2[4_456]	2.18	0.02

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	595/637 (93%)	539 (91%)	50 (8%)	6 (1%)	12	40
1	B	594/637 (93%)	545 (92%)	45 (8%)	4 (1%)	18	49
2	C	110/117 (94%)	98 (89%)	9 (8%)	3 (3%)	4	22
2	D	110/117 (94%)	106 (96%)	4 (4%)	0	100	100
All	All	1409/1508 (93%)	1288 (91%)	108 (8%)	13 (1%)	14	43

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	408	PRO
1	A	181	THR
1	A	312	GLN
1	A	313	SER
1	B	318	ALA
1	B	198	PRO
2	C	86	VAL
1	A	187	ILE
2	C	-1	GLY
2	C	0	SER
1	A	408	PRO
1	B	234	SER
1	A	692	ILE

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	536/568 (94%)	524 (98%)	12 (2%)	45	66
1	B	535/568 (94%)	526 (98%)	9 (2%)	53	71
2	C	98/102 (96%)	94 (96%)	4 (4%)	27	55
2	D	98/102 (96%)	97 (99%)	1 (1%)	68	76
All	All	1267/1340 (95%)	1241 (98%)	26 (2%)	47	67

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	154	VAL
1	A	185	LEU
1	A	201	ILE
1	A	214	LEU
1	A	294	SER
1	A	311	VAL
1	A	353	VAL
1	A	382	LEU
1	A	461	VAL
1	A	636	VAL
1	A	666	TRP
1	A	751	LEU
1	B	199	GLU
1	B	316	LEU
1	B	321	GLU
1	B	322	THR
1	B	411	ARG
1	B	453	MET
1	B	459	VAL
1	B	476	LEU
1	B	636	VAL
2	C	13	GLU
2	C	24	VAL
2	C	72	LYS

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Mol	Chain	Res	Type
2	C	76	THR
2	D	24	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	209	ASN
1	A	251	ASN
1	A	312	GLN
1	A	391	GLN
1	A	394	GLN
1	A	507	GLN
1	A	508	ASN
1	A	617	GLN
1	A	635	ASN
1	B	208	ASN
1	B	209	ASN
1	B	251	ASN
1	B	508	ASN
1	B	574	HIS
1	B	615	GLN
2	C	43	HIS
2	D	12	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, 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	599/637 (94%)	-0.24	3 (0%) 87 76	65, 106, 166, 191	0
1	B	598/637 (93%)	-0.11	6 (1%) 79 63	72, 123, 178, 201	0
2	C	112/117 (95%)	-0.09	3 (2%) 56 37	81, 121, 161, 168	0
2	D	112/117 (95%)	-0.16	0 100 100	72, 109, 147, 176	0
All	All	1421/1508 (94%)	-0.17	12 (0%) 82 68	65, 114, 170, 201	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	197	ILE	2.8
2	C	26	ASP	2.5
1	B	217	ASN	2.5
1	B	451	PHE	2.4
1	A	206	LEU	2.3
2	C	89	PHE	2.1
2	C	25	VAL	2.1
1	B	214	LEU	2.1
1	A	588	LEU	2.0
1	B	581	ALA	2.0
1	A	575	LEU	2.0
1	B	277	LEU	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 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.