



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:45 PM GMT

PDB ID : 4PUF
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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

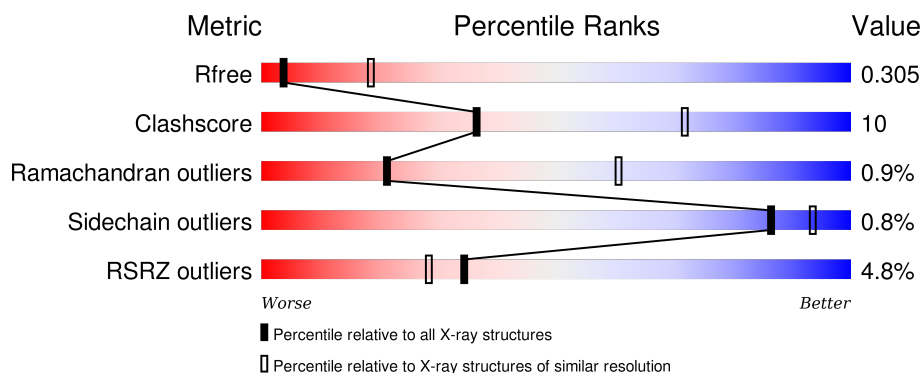
1 Overall quality at a glance

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}	91344	2060 (3.40-3.20)
Clashscore	102246	1058 (3.38-3.22)
Ramachandran outliers	100387	1038 (3.38-3.22)
Sidechain outliers	100360	1037 (3.38-3.22)
RSRZ outliers	91569	2070 (3.40-3.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	<div> <div>3%</div> <div>73%</div> <div>20%</div> <div>6%</div> </div>
1	B	637	<div> <div>5%</div> <div>67%</div> <div>26%</div> <div>6%</div> </div>
2	C	117	<div> <div>6%</div> <div>72%</div> <div>24%</div> <div>.</div> </div>
2	D	117	<div> <div>8%</div> <div>86%</div> <div>9%</div> <div>.</div> </div>

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			

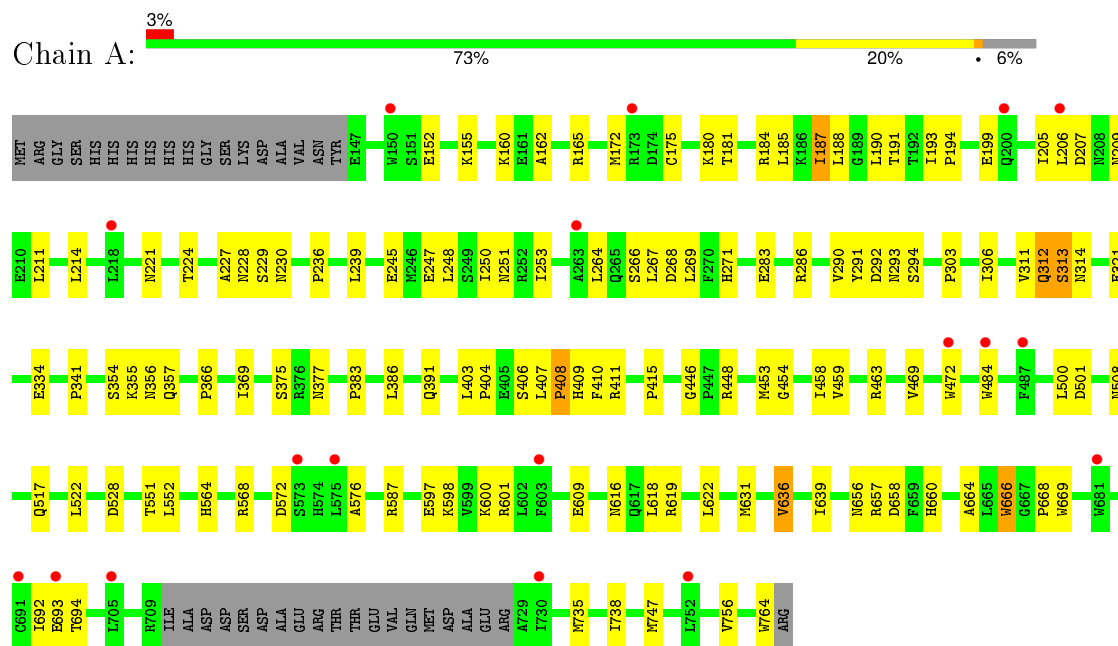
There are 24 discrepancies between the modelled and reference sequences:

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

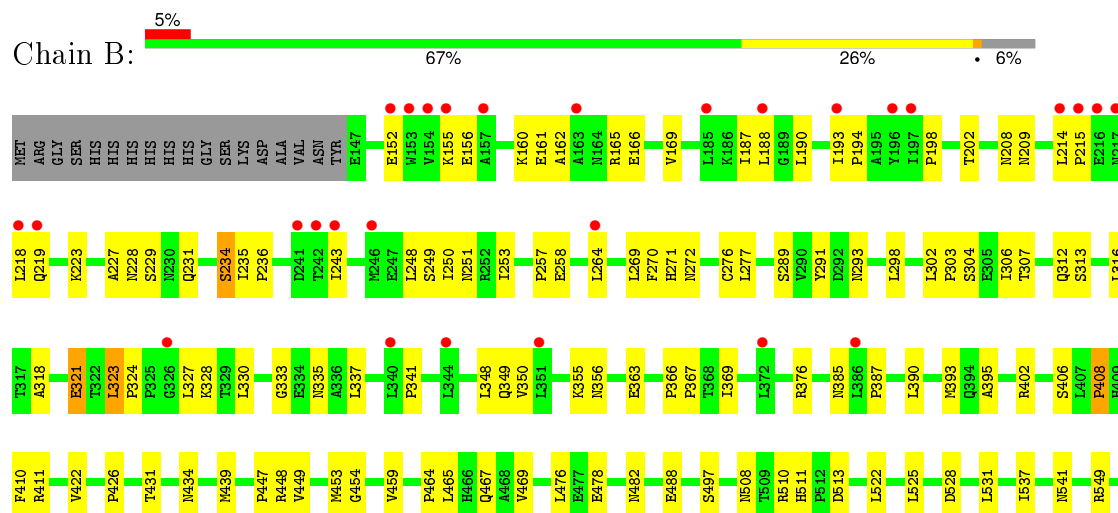
3 Residue-property plots

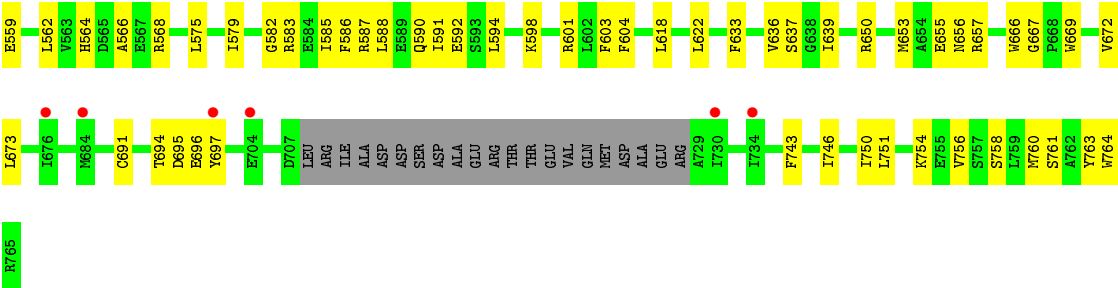
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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

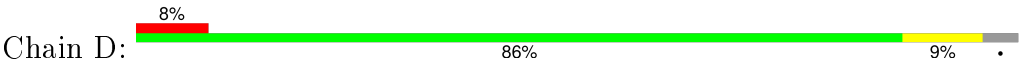




• Molecule 2: Thioredoxin



• 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 0.272 , 0.305	Depositor DCC
R_{free} test set	1698 reflections (5.00%)	DCC
Wilson B-factor (Å ²)	116.0	Xtriage
Anisotropy	0.552	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 54.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	2 of 34022 reflections (0.006%)	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.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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	0.22	0/4908	0.41	0/6667
1	B	0.25	0/4900	0.47	1/6656 (0.0%)
2	C	0.27	0/900	0.46	0/1210
2	D	0.22	0/900	0.34	0/1210
All	All	0.24	0/11608	0.44	1/15743 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	323	LEU	C-N-CD	6.14	141.29	128.40

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	84	1
1	B	4807	0	4783	109	0
2	C	880	0	854	21	0
2	D	880	0	854	9	0
All	All	11382	0	11285	216	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 (216) 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: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
1:A:664:ALA:HB2	1:A:747:MET:HE2	1.77	0.64
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: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:386:LEU:HB2	1:A:410:PHE:HE1	1.65	0.61
1:A:286:ARG:HA	1:A:306:ILE:HA	1.82	0.61
1:B:243:ILE:O	1:B:264:LEU:HD12	2.01	0.61
1:A:160:LYS:HZ3	1:A:190:LEU:HA	1.66	0.60
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:478:GLU:O	1:B:482:ASN:ND2	2.36	0.58
1:A:375:SER:O	1:A:377:ASN:ND2	2.37	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:B:234:SER:HA	1:B:253:ILE:HG13	1.85	0.58
1:B:694:THR:HB	1:B:697:TYR:HB2	1.85	0.58
1:A:264:LEU:HD21	1:A:267:LEU:HB2	1.86	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:OE1	2:C:6:GLU:N	2.38	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:A:160:LYS:HZ3	1:A:191:THR:H	1.53	0.56
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
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: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: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:406:SER:C	1:B:408:PRO:HD3	2.29	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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:A:406:SER:C	1:A:408:PRO:HD3	2.30	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:673:LEU:HD11	1:B:750:ILE:HD12	1.92	0.51
1:B:349:GLN:HG3	1:B:350:VAL:HG23	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
2:C:5:ILE:HG21	2:C:57:VAL:HG22	1.93	0.49
1:A:636:VAL:HG12	1:A:639:ILE:HB	1.95	0.49
2:C:23:VAL:HG12	2:C:53:ILE:HB	1.94	0.49
1:A:155:LYS:HE2	1:A:165:ARG:HH21	1.77	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: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:407:LEU:N	1:A:408:PRO:HD3	2.28	0.48
1:A:236:PRO:HG2	1:A:239:LEU:HG	1.95	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:229:SER:H	1:B:250:ILE:HB	1.79	0.47
1:B:227:ALA:HB3	1:B:248:LEU:HD23	1.97	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:277:LEU:HD12	1:B:298:LEU:HD22	1.98	0.46
1:B:439:MET:HE1	1:B:449:VAL:H	1.81	0.46
1:B:465:LEU:HG	1:B:497:SER:HB2	1.97	0.46
1:A:576:ALA:HB2	1:A:756:VAL:HG12	1.98	0.46
1:B:564:HIS:O	1:B:568:ARG:N	2.47	0.46
1:A:453:MET:N	1:A:454:GLY:HA3	2.30	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:B:562:LEU:HD12	1:B:585:ILE:HD11	1.98	0.45
1:A:334:GLU:HG2	1:A:355:LYS:HD2	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
2:C:18:ALA:HA	2:C:21:LYS:HZ2	1.82	0.45
1:B:307:THR:HA	1:B:327:LEU:HA	1.98	0.45
1:A:735:MET:O	1:A:738:ILE:HG13	2.17	0.45
1:A:564:HIS:HB3	1:A:568:ARG:NH1	2.32	0.44
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:694:THR:HG22	1:B:696:GLU:H	1.83	0.44
1:B:327:LEU:HD21	1:B:330:LEU:HB2	2.00	0.44
1:A:184:ARG:HH21	2:D:61:ASP:CG	2.20	0.44
2:C:3:LYS:HB3	2:C:55:LEU:HD23	2.00	0.44
1:B:549:ARG:NH2	1:B:592:GLU:OE2	2.51	0.44
1:B:161:GLU:OE1	1:B:162:ALA:N	2.50	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
1:A:248:LEU:HD22	1:A:253:ILE:HD11	1.99	0.43
2:D:46:SER:HA	2:D:54:PHE:CE2	2.53	0.43
1:B:585:ILE:HD12	1:B:666:TRP:CZ2	2.53	0.43
1:B:152:GLU:O	1:B:156:GLU:N	2.51	0.43
1:B:469:VAL:HG13	1:B:522:LEU:HD23	1.99	0.43
1:B:258:GLU:HG3	1:B:258:GLU:O	2.18	0.43
2:C:80:PHE:CE1	2:C:85:LYS:HA	2.52	0.43
1:B:249:SER:HB3	1:B:270:PHE:HD2	1.84	0.43
1:B:582:GLY:HA2	1:B:666:TRP:CH2	2.54	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:A:366:PRO:HB2	1:A:369:ILE:HG13	2.00	0.43
1:B:508:ASN:HB3	1:B:633:PHE:CD1	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:271:HIS:HE1	2:C:-2:HIS:HB2	1.84	0.42
2:D:38:ILE:HD13	2:D:77:PHE:CZ	2.54	0.42
1:B:395:ALA:HB3	1:B:422:VAL:HG12	2.01	0.42
2:C:8:LYS:O	2:C:11:PHE:HB3	2.19	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
1:A:160:LYS:NZ	1:A:190:LEU:HA	2.33	0.42
1:B:439:MET:HE1	1:B:448:ARG:HA	2.01	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:B:229:SER:HA	1:B:250:ILE:HB	2.02	0.42
2:C:6:GLU:CG	2:C:7:SER:H	2.32	0.42
1:A:245:GLU:HG2	1:A:266:SER:HB3	2.01	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
1:B:434:ASN:ND2	1:B:603:PHE:O	2.53	0.42
2:D:6:GLU:OE1	2:D:6:GLU:N	2.53	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
1:A:446:GLY:O	1:A:448:ARG:NH1	2.52	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
1:B:328:LYS:HB3	1:B:349:GLN:HG2	2.01	0.41
1:A:247:GLU:HA	1:A:268:ASP:HB3	2.01	0.41
2:C:6:GLU:HG2	2:C:7:SER:N	2.36	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
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:666:TRP:HA	1:B:667:GLY:HA3	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:321:GLU:HA	1:B:341:PRO:HB3	2.03	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:A:152:GLU:OE1	1:A:172:MET:HG2	2.21	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:H	1:A:458:ILE:HD12	1.85	0.40
1:B:695:ASP:N	1:B:695:ASP:OD1	2.55	0.40
1:B:426:PRO:HB2	1:B:510:ARG:HH12	1.86	0.40
2:D:73:CYS:O	2:D:76:THR:OG1	2.24	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

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	595/637 (93%)	539 (91%)	50 (8%)	6 (1%)	19	58
1	B	594/637 (93%)	545 (92%)	45 (8%)	4 (1%)	26	66
2	C	110/117 (94%)	98 (89%)	9 (8%)	3 (3%)	6	35
2	D	110/117 (94%)	106 (96%)	4 (4%)	0	100	100
All	All	1409/1508 (93%)	1288 (91%)	108 (8%)	13 (1%)	21	60

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%)	531 (99%)	5 (1%)	84	92
1	B	535/568 (94%)	531 (99%)	4 (1%)	88	94
2	C	98/102 (96%)	97 (99%)	1 (1%)	82	91
2	D	98/102 (96%)	98 (100%)	0	100	100
All	All	1267/1340 (95%)	1257 (99%)	10 (1%)	86	93

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	214	LEU
1	A	294	SER
1	A	636	VAL
1	A	660	HIS
1	A	666	TRP
1	B	316	LEU
1	B	321	GLU
1	B	411	ARG
1	B	476	LEU

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Mol	Chain	Res	Type
2	C	76	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	209	ASN
1	A	251	ASN
1	A	391	GLN
1	B	208	ASN
1	B	251	ASN
1	B	508	ASN
2	C	43	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 carbohydrates 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	599/637 (94%)	0.22	18 (3%) 54 47	65, 106, 166, 191	0
1	B	598/637 (93%)	0.37	34 (5%) 27 22	72, 123, 178, 201	0
2	C	112/117 (95%)	0.53	7 (6%) 23 19	81, 121, 161, 168	0
2	D	112/117 (95%)	0.64	9 (8%) 15 12	72, 109, 147, 176	0
All	All	1421/1508 (94%)	0.34	68 (4%) 34 28	65, 114, 170, 201	0

All (68) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	217	ASN	10.1
1	B	196	TYR	5.9
1	B	219	GLN	5.9
1	B	242	THR	5.0
1	A	693	GLU	4.6
1	B	684	MET	4.6
1	B	730	ILE	4.5
2	D	81	LYS	4.2
1	B	326	GLY	3.7
1	A	575	LEU	3.5
1	B	197	ILE	3.5
2	D	20	ASP	3.4
1	B	241	ASP	3.4
1	A	173	ARG	3.3
1	A	681	TRP	3.2
1	B	154	VAL	3.2
2	C	13	GLU	3.1
1	B	216	GLU	3.1
1	B	163	ALA	3.1
1	B	676	ILE	3.0
1	B	704	GLU	3.0

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Mol	Chain	Res	Type	RSRZ
2	C	-1	GLY	2.8
1	A	573	SER	2.8
1	B	243	ILE	2.8
2	C	55	LEU	2.8
2	D	21	LYS	2.8
2	C	79	PHE	2.7
2	D	17	ALA	2.7
1	A	150	TRP	2.7
1	B	218	LEU	2.7
1	A	218	LEU	2.7
1	A	730	ILE	2.6
1	B	734	ILE	2.6
1	B	372	LEU	2.6
1	B	193	ILE	2.6
1	B	386	LEU	2.5
1	A	691	CYS	2.5
2	D	80	PHE	2.5
1	B	215	PRO	2.4
1	B	214	LEU	2.4
1	B	351	LEU	2.4
1	B	246	MET	2.4
1	A	484	TRP	2.4
2	C	19	GLY	2.3
1	A	200	GLN	2.3
1	B	152	GLU	2.3
1	B	155	LYS	2.3
1	B	697	TYR	2.2
2	C	26	ASP	2.2
1	B	157	ALA	2.2
1	A	487	PHE	2.2
1	B	264	LEU	2.2
1	B	344	LEU	2.2
1	A	603	PHE	2.2
1	A	206	LEU	2.2
1	B	340	LEU	2.2
2	D	54	PHE	2.2
1	B	188	LEU	2.2
2	D	79	PHE	2.2
1	A	752	LEU	2.1
1	B	185	LEU	2.1
2	C	15	LEU	2.1
1	A	472	TRP	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	153	TRP	2.1
2	D	4	GLN	2.1
1	A	705	LEU	2.0
2	D	57	VAL	2.0
1	A	263	ALA	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 carbohydrates 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.