



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:36 AM GMT

PDB ID : 2RAL
Title : Crystal Structure Analysis of double cysteine mutant of S.epidermidis adhesin SdrG: Evidence for the Dock,Lock and Latch ligand binding mechanism
Authors : Ponnuraj, K.; Sthanam, N.; Bowden, M.G.; Hook, M.
Deposited on : 2007-09-17
Resolution : 2.80 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

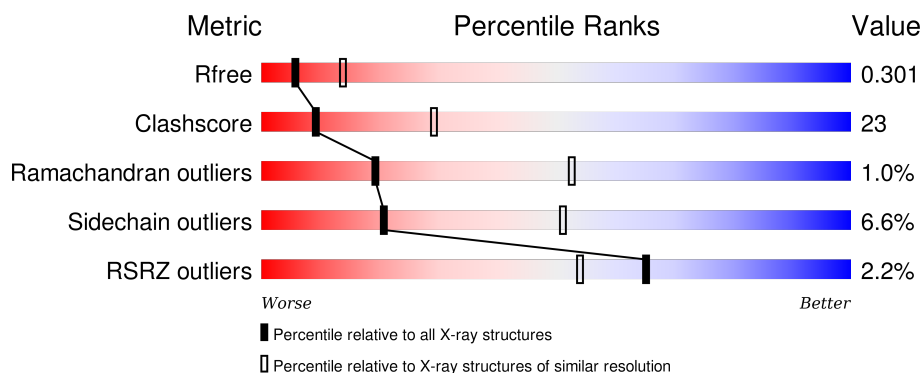
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

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	340	<div> <div></div> <div>60% 32% 6%</div> </div>
1	B	340	<div> <div>3%</div> <div>54% 34% 5% 7%</div> </div>

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 4966 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 Serine-aspartate repeat-containing protein G.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	320	Total	C	N	O	S	0	0	0
			2491	1537	405	544	5			
1	B	315	Total	C	N	O	S	0	0	0
			2475	1530	404	536	5			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	256	MET	-	EXPRESSION TAG	UNP Q9KI13
A	257	GLY	-	EXPRESSION TAG	UNP Q9KI13
A	258	ARG	-	EXPRESSION TAG	UNP Q9KI13
A	259	SER	-	EXPRESSION TAG	UNP Q9KI13
A	260	HIS	-	EXPRESSION TAG	UNP Q9KI13
A	261	HIS	-	EXPRESSION TAG	UNP Q9KI13
A	262	HIS	-	EXPRESSION TAG	UNP Q9KI13
A	263	HIS	-	EXPRESSION TAG	UNP Q9KI13
A	264	HIS	-	EXPRESSION TAG	UNP Q9KI13
A	265	HIS	-	EXPRESSION TAG	UNP Q9KI13
A	266	GLY	-	EXPRESSION TAG	UNP Q9KI13
A	267	SER	-	EXPRESSION TAG	UNP Q9KI13
A	268	LEU	-	EXPRESSION TAG	UNP Q9KI13
A	269	VAL	-	EXPRESSION TAG	UNP Q9KI13
A	270	PRO	-	EXPRESSION TAG	UNP Q9KI13
A	271	ARG	-	EXPRESSION TAG	UNP Q9KI13
A	272	GLY	-	EXPRESSION TAG	UNP Q9KI13
A	273	SER	-	EXPRESSION TAG	UNP Q9KI13
A	381	CYS	GLU	ENGINEERED	UNP Q9KI13
A	595	CYS	PRO	ENGINEERED	UNP Q9KI13
B	256	MET	-	EXPRESSION TAG	UNP Q9KI13
B	257	GLY	-	EXPRESSION TAG	UNP Q9KI13
B	258	ARG	-	EXPRESSION TAG	UNP Q9KI13
B	259	SER	-	EXPRESSION TAG	UNP Q9KI13
B	260	HIS	-	EXPRESSION TAG	UNP Q9KI13

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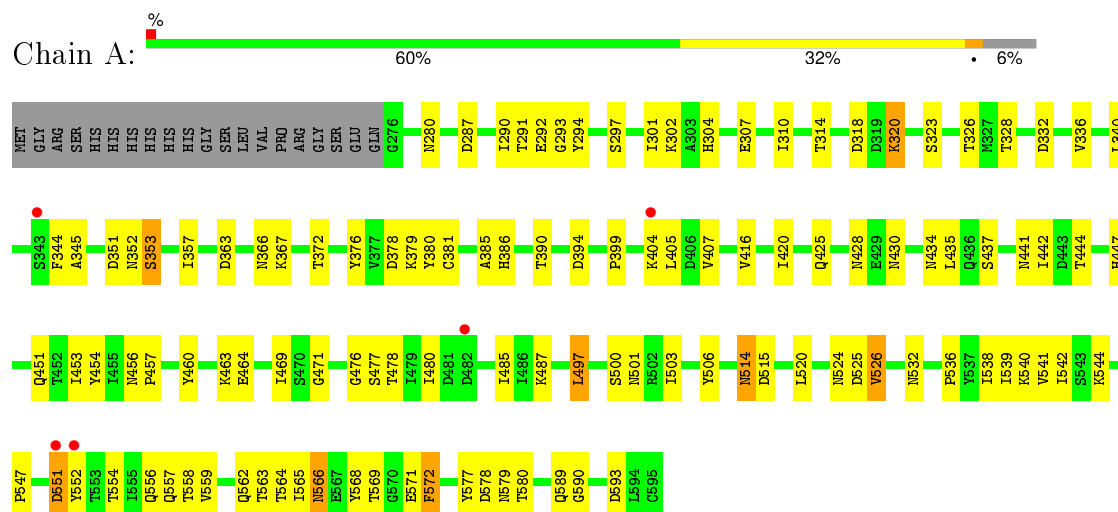
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Chain	Residue	Modelled	Actual	Comment	Reference
B	261	HIS	-	EXPRESSION TAG	UNP Q9KI13
B	262	HIS	-	EXPRESSION TAG	UNP Q9KI13
B	263	HIS	-	EXPRESSION TAG	UNP Q9KI13
B	264	HIS	-	EXPRESSION TAG	UNP Q9KI13
B	265	HIS	-	EXPRESSION TAG	UNP Q9KI13
B	266	GLY	-	EXPRESSION TAG	UNP Q9KI13
B	267	SER	-	EXPRESSION TAG	UNP Q9KI13
B	268	LEU	-	EXPRESSION TAG	UNP Q9KI13
B	269	VAL	-	EXPRESSION TAG	UNP Q9KI13
B	270	PRO	-	EXPRESSION TAG	UNP Q9KI13
B	271	ARG	-	EXPRESSION TAG	UNP Q9KI13
B	272	GLY	-	EXPRESSION TAG	UNP Q9KI13
B	273	SER	-	EXPRESSION TAG	UNP Q9KI13
B	381	CYS	GLU	ENGINEERED	UNP Q9KI13
B	595	CYS	PRO	ENGINEERED	UNP Q9KI13

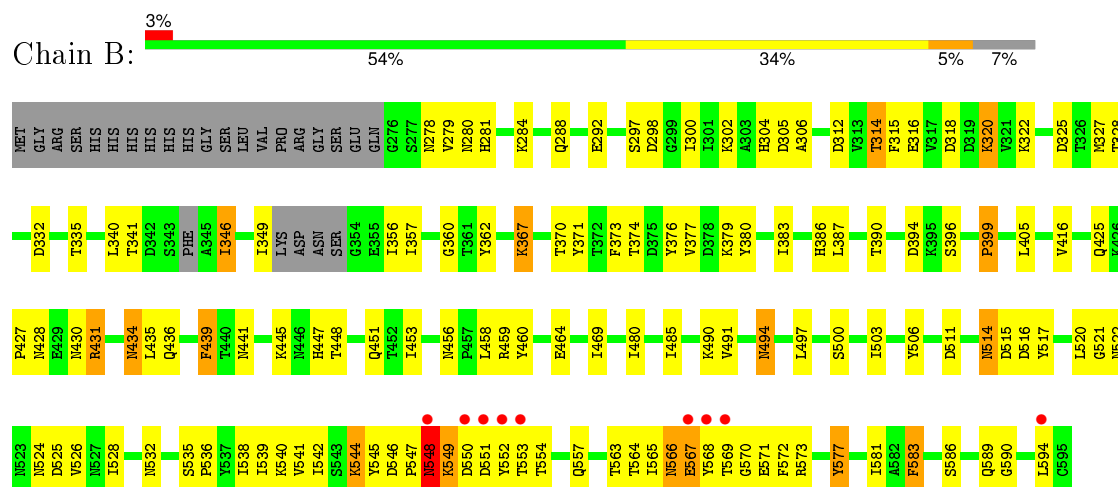
3 Residue-property plots

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 ($\text{RSRZ} > 2$). 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: Serine-aspartate repeat-containing protein G



• Molecule 1: Serine-aspartate repeat-containing protein G



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	61.51Å 94.17Å 129.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.99 – 2.80 32.38 – 2.79	Depositor EDS
% Data completeness (in resolution range)	97.3 (19.99-2.80) 96.7 (32.38-2.79)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.57 (at 2.81Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.242 , 0.303 0.241 , 0.301	Depositor DCC
R_{free} test set	744 reflections (4.00%)	DCC
Wilson B-factor (Å ²)	32.7	Xtriage
Anisotropy	0.220	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 29.5	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	0 of 19157 reflections	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	4966	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.17% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.44	0/2529	0.70	0/3444
1	B	0.44	0/2511	0.68	1/3414 (0.0%)
All	All	0.44	0/5040	0.69	1/6858 (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	405	LEU	CA-CB-CG	6.03	129.16	115.30

There are no chirality outliers.

There are no planarity outliers.

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	2491	0	2342	101	0
1	B	2475	0	2353	122	0
All	All	4966	0	4695	221	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 23.

All (221) 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:304:HIS:HD2	1:B:425:GLN:H	1.11	0.94
1:A:304:HIS:HD2	1:A:425:GLN:H	0.94	0.94
1:A:304:HIS:CD2	1:A:425:GLN:H	1.85	0.92
1:A:480:ILE:H	1:A:524:ASN:HD21	1.15	0.91
1:B:456:ASN:HD21	1:B:460:TYR:H	1.20	0.89
1:B:346:ILE:HD11	1:B:362:TYR:HB2	1.57	0.87
1:B:314:THR:HG23	1:B:386:HIS:HB2	1.57	0.87
1:A:480:ILE:H	1:A:524:ASN:ND2	1.77	0.83
1:B:544:LYS:H	1:B:544:LYS:HD2	1.40	0.83
1:B:520:LEU:HA	1:B:526:VAL:HG23	1.59	0.82
1:A:332:ASP:OD1	1:A:407:VAL:HA	1.79	0.82
1:A:464:GLU:H	1:A:532:ASN:HD22	1.28	0.81
1:B:280:ASN:HD21	1:B:328:THR:H	1.30	0.79
1:B:304:HIS:CD2	1:B:425:GLN:H	1.98	0.79
1:A:464:GLU:H	1:A:532:ASN:ND2	1.82	0.76
1:B:278:ASN:HD21	1:B:280:ASN:HB2	1.49	0.76
1:A:503:ILE:HG13	1:A:538:ILE:HD13	1.66	0.76
1:A:451:GLN:NE2	1:A:557:GLN:HE21	1.85	0.75
1:A:447:HIS:O	1:A:544:LYS:HG3	1.87	0.73
1:A:318:ASP:OD2	1:A:320:LYS:HB2	1.88	0.73
1:B:346:ILE:HG21	1:B:360:GLY:O	1.90	0.71
1:B:494:ASN:HD22	1:B:494:ASN:N	1.88	0.71
1:B:451:GLN:NE2	1:B:557:GLN:HE21	1.90	0.70
1:A:304:HIS:HD2	1:A:425:GLN:N	1.79	0.69
1:A:566:ASN:HD22	1:A:569:THR:H	1.41	0.69
1:B:567:GLU:H	1:B:567:GLU:CD	1.96	0.69
1:A:340:LEU:HB3	1:A:442:ILE:HG22	1.76	0.68
1:B:552:TYR:HB2	1:B:583:PHE:CE2	2.29	0.68
1:B:544:LYS:CD	1:B:544:LYS:H	2.08	0.66
1:B:314:THR:OG1	1:B:386:HIS:HD2	1.79	0.65
1:B:494:ASN:HD22	1:B:494:ASN:H	1.44	0.65
1:B:304:HIS:HD2	1:B:425:GLN:N	1.89	0.65
1:A:314:THR:OG1	1:A:386:HIS:HD2	1.78	0.65
1:B:514:ASN:HD22	1:B:515:ASP:N	1.95	0.65
1:B:494:ASN:H	1:B:494:ASN:ND2	1.96	0.64
1:A:564:THR:HA	1:A:572:PHE:HB3	1.80	0.64
1:A:425:GLN:HE21	1:A:501:ASN:ND2	1.95	0.63
1:B:552:TYR:HB2	1:B:583:PHE:HE2	1.62	0.63
1:B:304:HIS:CD2	1:B:425:GLN:HG2	2.33	0.63
1:B:480:ILE:H	1:B:524:ASN:HD21	1.46	0.62
1:B:520:LEU:CA	1:B:526:VAL:HG23	2.29	0.62
1:B:567:GLU:HG2	1:B:568:TYR:CD1	2.35	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:594:LEU:N	1:B:594:LEU:HD12	2.14	0.62
1:B:434:ASN:HD21	1:B:456:ASN:HD22	1.45	0.62
1:A:280:ASN:HD21	1:A:328:THR:H	1.47	0.62
1:A:480:ILE:N	1:A:524:ASN:HD21	1.95	0.61
1:A:476:GLY:HA3	1:A:554:THR:O	2.01	0.61
1:A:563:THR:O	1:A:572:PHE:CB	2.49	0.61
1:B:434:ASN:HD21	1:B:456:ASN:ND2	1.99	0.59
1:A:544:LYS:NZ	1:A:544:LYS:HB2	2.17	0.59
1:A:471:GLY:O	1:A:477:SER:HB2	2.02	0.59
1:B:565:ILE:HD12	1:B:566:ASN:N	2.17	0.59
1:B:544:LYS:CD	1:B:544:LYS:N	2.66	0.59
1:B:480:ILE:H	1:B:524:ASN:ND2	2.01	0.59
1:A:456:ASN:HD21	1:A:460:TYR:H	1.51	0.59
1:A:563:THR:O	1:A:572:PHE:HB2	2.03	0.59
1:B:548:ASN:ND2	1:B:548:ASN:H	2.01	0.58
1:A:379:LYS:HE2	1:A:380:TYR:CE1	2.38	0.58
1:B:503:ILE:HG13	1:B:538:ILE:HD13	1.84	0.58
1:B:594:LEU:CD1	1:B:594:LEU:H	2.17	0.57
1:B:514:ASN:HD22	1:B:514:ASN:C	2.07	0.57
1:A:351:ASP:O	1:A:353:SER:N	2.38	0.57
1:B:428:ASN:HB2	1:B:577:TYR:HD1	1.70	0.57
1:B:566:ASN:O	1:B:570:GLY:N	2.38	0.56
1:A:363:ASP:HB3	1:A:366:ASN:OD1	2.05	0.56
1:A:554:THR:CG2	1:A:580:THR:HG23	2.35	0.56
1:A:351:ASP:OD2	1:A:353:SER:HB3	2.05	0.56
1:B:377:VAL:HA	1:B:383:ILE:CD1	2.35	0.56
1:A:463:LYS:HA	1:A:532:ASN:HD22	1.71	0.55
1:B:292:GLU:OE1	1:B:297:SER:O	2.23	0.55
1:A:457:PRO:HA	1:A:536:PRO:HB3	1.89	0.55
1:A:554:THR:HG21	1:A:580:THR:HG23	1.88	0.55
1:B:288:GLN:HA	1:B:312:ASP:O	2.06	0.55
1:A:425:GLN:HE21	1:A:501:ASN:HD22	1.52	0.55
1:B:428:ASN:HB2	1:B:577:TYR:CD1	2.42	0.54
1:A:566:ASN:ND2	1:A:568:TYR:H	2.05	0.54
1:A:520:LEU:HA	1:A:526:VAL:HG23	1.88	0.54
1:B:322:LYS:HG2	1:B:325:ASP:OD1	2.08	0.54
1:B:594:LEU:N	1:B:594:LEU:CD1	2.71	0.54
1:A:520:LEU:HD23	1:A:526:VAL:HG23	1.90	0.54
1:B:453:ILE:HB	1:B:539:ILE:HB	1.90	0.54
1:A:290:ILE:CD1	1:A:416:VAL:HG13	2.38	0.53
1:B:548:ASN:O	1:B:550:ASP:N	2.41	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:539:ILE:HG22	1:A:541:VAL:HG23	1.90	0.53
1:A:326:THR:HG22	1:A:372:THR:HG23	1.91	0.53
1:A:551:ASP:HB3	1:A:552:TYR:CD1	2.44	0.53
1:B:464:GLU:H	1:B:532:ASN:ND2	2.07	0.52
1:B:447:HIS:O	1:B:544:LYS:HB2	2.08	0.52
1:B:517:TYR:O	1:B:528:ILE:HA	2.09	0.52
1:B:386:HIS:C	1:B:387:LEU:HD23	2.30	0.52
1:B:399:PRO:HA	1:B:500:SER:HB3	1.91	0.52
1:A:566:ASN:ND2	1:A:569:THR:H	2.05	0.52
1:B:427:PRO:HA	1:B:436:GLN:HB3	1.90	0.52
1:B:451:GLN:HE21	1:B:557:GLN:HE21	1.58	0.52
1:B:373:PHE:HB3	1:B:377:VAL:HG11	1.92	0.52
1:B:469:ILE:HD13	1:B:541:VAL:HG21	1.92	0.52
1:B:376:TYR:O	1:B:380:TYR:HD1	1.92	0.52
1:A:525:ASP:OD1	1:A:525:ASP:N	2.43	0.52
1:A:351:ASP:C	1:A:353:SER:H	2.13	0.51
1:A:453:ILE:HG12	1:A:559:VAL:HG21	1.93	0.51
1:A:345:ALA:HB2	1:B:522:ASN:HA	1.93	0.51
1:A:404:LYS:O	1:A:405:LEU:HD23	2.11	0.51
1:A:478:THR:HG23	1:A:556:GLN:O	2.11	0.51
1:B:445:LYS:HG2	1:B:445:LYS:O	2.11	0.51
1:A:566:ASN:HD22	1:A:568:TYR:H	1.58	0.50
1:A:292:GLU:OE1	1:A:297:SER:O	2.28	0.50
1:A:453:ILE:HB	1:A:539:ILE:HB	1.93	0.50
1:A:404:LYS:N	1:A:404:LYS:HD2	2.26	0.50
1:B:340:LEU:O	1:B:441:ASN:HA	2.11	0.50
1:B:548:ASN:ND2	1:B:548:ASN:N	2.57	0.50
1:B:547:PRO:O	1:B:549:LYS:N	2.45	0.50
1:B:297:SER:HB2	1:B:302:LYS:HE2	1.93	0.49
1:A:514:ASN:HD22	1:A:514:ASN:C	2.14	0.49
1:B:292:GLU:OE2	1:B:298:ASP:HA	2.11	0.49
1:A:291:THR:O	1:A:310:ILE:N	2.42	0.49
1:B:357:ILE:CD1	1:B:590:GLY:HA3	2.42	0.49
1:A:454:TYR:CZ	1:A:497:LEU:HD22	2.47	0.49
1:A:464:GLU:N	1:A:532:ASN:HD22	2.03	0.49
1:A:551:ASP:O	1:A:552:TYR:C	2.50	0.49
1:A:463:LYS:HA	1:A:532:ASN:ND2	2.29	0.48
1:B:428:ASN:O	1:B:434:ASN:HA	2.13	0.48
1:A:297:SER:OG	1:A:297:SER:O	2.25	0.48
1:A:293:GLY:C	1:A:294:TYR:HD2	2.17	0.48
1:B:284:LYS:HD2	1:B:316:GLU:OE1	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:434:ASN:N	1:B:434:ASN:HD22	2.10	0.48
1:B:566:ASN:HB3	1:B:571:GLU:HB3	1.95	0.48
1:A:565:ILE:HD12	1:A:565:ILE:C	2.34	0.48
1:B:521:GLY:H	1:B:526:VAL:HA	1.79	0.48
1:B:390:THR:HG22	1:B:586:SER:OG	2.13	0.48
1:B:506:TYR:CE2	1:B:540:LYS:HD3	2.49	0.47
1:A:566:ASN:HD22	1:A:568:TYR:N	2.12	0.47
1:A:435:LEU:N	1:A:435:LEU:HD23	2.28	0.47
1:A:351:ASP:C	1:A:353:SER:N	2.67	0.47
1:B:485:ILE:HB	1:B:542:ILE:HB	1.97	0.47
1:B:546:ASP:HB3	1:B:549:LYS:HG3	1.96	0.47
1:B:305:ASP:O	1:B:306:ALA:HB3	2.15	0.47
1:B:514:ASN:ND2	1:B:514:ASN:C	2.68	0.47
1:B:458:LEU:O	1:B:459:ARG:HB2	2.14	0.47
1:A:506:TYR:CZ	1:A:540:LYS:HD3	2.49	0.47
1:B:332:ASP:HB3	1:B:335:THR:HG23	1.97	0.47
1:A:357:ILE:HG12	1:A:376:TYR:CE2	2.50	0.47
1:A:442:ILE:CD1	1:A:444:THR:HG22	2.44	0.46
1:B:583:PHE:C	1:B:583:PHE:CD1	2.88	0.46
1:B:304:HIS:HB3	1:B:425:GLN:OE1	2.16	0.46
1:A:476:GLY:O	1:A:477:SER:HB3	2.16	0.46
1:B:427:PRO:HG3	1:B:458:LEU:HD21	1.97	0.46
1:B:374:THR:O	1:B:377:VAL:HG22	2.16	0.46
1:A:323:SER:HB3	1:A:378:ASP:OD2	2.15	0.46
1:B:566:ASN:OD1	1:B:569:THR:CG2	2.64	0.46
1:A:291:THR:HB	1:A:310:ILE:HB	1.98	0.45
1:A:425:GLN:NE2	1:A:501:ASN:ND2	2.63	0.45
1:B:360:GLY:HA2	1:B:370:THR:O	2.17	0.45
1:B:380:TYR:CD1	1:B:380:TYR:N	2.84	0.45
1:B:318:ASP:OD2	1:B:320:LYS:HE3	2.17	0.45
1:B:564:THR:HA	1:B:572:PHE:HA	1.98	0.45
1:A:558:THR:HG23	1:A:578:ASP:HB3	1.97	0.45
1:A:301:ILE:HG12	1:A:420:ILE:HD12	1.98	0.45
1:B:544:LYS:HE3	1:B:544:LYS:N	2.31	0.45
1:A:293:GLY:O	1:A:294:TYR:HD2	2.00	0.44
1:A:469:ILE:HD13	1:A:541:VAL:HG21	1.99	0.44
1:A:506:TYR:CE2	1:A:540:LYS:HD3	2.53	0.44
1:B:394:ASP:OD2	1:B:396:SER:HB3	2.17	0.44
1:A:336:VAL:HG13	1:A:394:ASP:HB2	2.00	0.44
1:B:544:LYS:CE	1:B:544:LYS:N	2.81	0.44
1:B:439:PHE:CD1	1:B:581:ILE:HG12	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:566:ASN:HD22	1:B:566:ASN:C	2.20	0.44
1:B:550:ASP:CG	1:B:551:ASP:N	2.72	0.43
1:B:551:ASP:HB3	1:B:554:THR:OG1	2.19	0.43
1:B:431:ARG:HG3	1:B:573:ARG:HG3	2.00	0.43
1:A:340:LEU:O	1:A:441:ASN:HA	2.17	0.43
1:A:566:ASN:ND2	1:A:569:THR:N	2.66	0.43
1:B:292:GLU:OE1	1:B:300:ILE:O	2.35	0.43
1:B:547:PRO:C	1:B:549:LYS:N	2.70	0.43
1:B:280:ASN:HD21	1:B:328:THR:N	2.07	0.43
1:B:349:ILE:HG13	1:B:371:TYR:CE2	2.53	0.43
1:B:514:ASN:ND2	1:B:516:ASP:H	2.17	0.43
1:B:539:ILE:HG22	1:B:541:VAL:HG23	2.00	0.43
1:B:315:PHE:N	1:B:315:PHE:CD2	2.86	0.43
1:A:301:ILE:HA	1:A:307:GLU:OE2	2.19	0.42
1:A:301:ILE:CG1	1:A:420:ILE:HD12	2.48	0.42
1:B:315:PHE:CE1	1:B:327:MET:SD	3.12	0.42
1:A:437:SER:OG	1:A:579:ASN:ND2	2.52	0.42
1:A:469:ILE:HB	1:A:526:VAL:HG12	2.02	0.42
1:A:566:ASN:HB3	1:A:569:THR:OG1	2.20	0.42
1:A:399:PRO:HA	1:A:500:SER:HB3	2.00	0.42
1:A:487:LYS:HA	1:A:487:LYS:HD3	1.84	0.42
1:A:562:GLN:HG2	1:A:563:THR:N	2.35	0.42
1:B:548:ASN:HD22	1:B:548:ASN:N	2.17	0.42
1:A:471:GLY:O	1:A:477:SER:CB	2.67	0.42
1:B:279:VAL:HG13	1:B:325:ASP:HB3	2.01	0.42
1:B:340:LEU:HD13	1:B:581:ILE:HD12	2.02	0.42
1:A:464:GLU:N	1:A:532:ASN:ND2	2.58	0.42
1:B:547:PRO:C	1:B:549:LYS:H	2.21	0.42
1:B:563:THR:O	1:B:572:PHE:HB2	2.20	0.41
1:A:302:LYS:N	1:A:307:GLU:OE2	2.42	0.41
1:B:583:PHE:C	1:B:583:PHE:HD1	2.24	0.41
1:B:356:ILE:HG22	1:B:374:THR:HG21	2.01	0.41
1:A:435:LEU:N	1:A:435:LEU:CD2	2.83	0.41
1:A:385:ALA:HA	1:A:590:GLY:HA2	2.01	0.41
1:A:345:ALA:HA	1:B:522:ASN:HB2	2.02	0.41
1:A:381:CYS:N	1:A:593:ASP:O	2.37	0.41
1:B:520:LEU:N	1:B:526:VAL:HG23	2.36	0.41
1:B:356:ILE:HG22	1:B:356:ILE:O	2.20	0.41
1:B:459:ARG:HG2	1:B:536:PRO:HD3	2.02	0.41
1:A:290:ILE:HD12	1:A:416:VAL:HG13	2.03	0.41
1:B:525:ASP:OD1	1:B:525:ASP:N	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:363:ASP:OD2	1:A:366:ASN:ND2	2.53	0.41
1:B:545:TYR:HB2	1:B:549:LYS:HD2	2.02	0.41
1:B:367:LYS:N	1:B:367:LYS:HD2	2.36	0.41
1:A:544:LYS:HZ2	1:A:544:LYS:HB2	1.85	0.41
1:B:373:PHE:HB3	1:B:377:VAL:CG1	2.51	0.41
1:A:569:THR:OG1	1:A:571:GLU:HB3	2.20	0.40
1:B:567:GLU:CD	1:B:567:GLU:N	2.69	0.40
1:A:357:ILE:HG12	1:A:376:TYR:CD2	2.56	0.40
1:B:566:ASN:OD1	1:B:569:THR:HG23	2.22	0.40
1:B:490:LYS:HE3	1:B:535:SER:CB	2.52	0.40
1:B:491:VAL:HG21	1:B:497:LEU:HD21	2.02	0.40
1:A:485:ILE:HB	1:A:542:ILE:HB	2.03	0.40
1:B:448:THR:HA	1:B:544:LYS:HA	2.03	0.40
1:A:428:ASN:O	1:A:434:ASN:HA	2.22	0.40
1:B:511:ASP:OD1	1:B:511:ASP:C	2.60	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	318/340 (94%)	290 (91%)	26 (8%)	2 (1%)	30	65
1	B	309/340 (91%)	276 (89%)	29 (9%)	4 (1%)	15	44
All	All	627/680 (92%)	566 (90%)	55 (9%)	6 (1%)	19	52

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	352	ASN
1	B	549	LYS
1	B	548	ASN

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Mol	Chain	Res	Type
1	B	399	PRO
1	A	515	ASP
1	B	346	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	286/310 (92%)	270 (94%)	16 (6%)	26	59
1	B	286/310 (92%)	264 (92%)	22 (8%)	16	41
All	All	572/620 (92%)	534 (93%)	38 (7%)	21	51

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	287	ASP
1	A	320	LYS
1	A	344	PHE
1	A	353	SER
1	A	367	LYS
1	A	390	THR
1	A	430	ASN
1	A	497	LEU
1	A	514	ASN
1	A	526	VAL
1	A	547	PRO
1	A	551	ASP
1	A	566	ASN
1	A	572	PHE
1	A	577	TYR
1	A	589	GLN
1	B	281	HIS
1	B	314	THR
1	B	320	LYS
1	B	341	THR

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Mol	Chain	Res	Type
1	B	367	LYS
1	B	379	LYS
1	B	416	VAL
1	B	430	ASN
1	B	431	ARG
1	B	434	ASN
1	B	435	LEU
1	B	439	PHE
1	B	494	ASN
1	B	514	ASN
1	B	544	LYS
1	B	548	ASN
1	B	553	THR
1	B	566	ASN
1	B	567	GLU
1	B	577	TYR
1	B	583	PHE
1	B	589	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (43) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	278	ASN
1	A	280	ASN
1	A	304	HIS
1	A	368	GLN
1	A	386	HIS
1	A	400	ASN
1	A	402	ASN
1	A	430	ASN
1	A	436	GLN
1	A	447	HIS
1	A	451	GLN
1	A	456	ASN
1	A	472	ASN
1	A	496	ASN
1	A	501	ASN
1	A	514	ASN
1	A	524	ASN
1	A	532	ASN
1	A	548	ASN
1	A	566	ASN

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Mol	Chain	Res	Type
1	A	579	ASN
1	B	278	ASN
1	B	280	ASN
1	B	281	HIS
1	B	304	HIS
1	B	334	ASN
1	B	386	HIS
1	B	400	ASN
1	B	402	ASN
1	B	430	ASN
1	B	436	GLN
1	B	451	GLN
1	B	456	ASN
1	B	466	ASN
1	B	472	ASN
1	B	494	ASN
1	B	501	ASN
1	B	514	ASN
1	B	524	ASN
1	B	529	ASN
1	B	532	ASN
1	B	579	ASN
1	B	589	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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	320/340 (94%)	-0.33	5 (1%) 74 66	14, 23, 33, 43	0
1	B	315/340 (92%)	-0.11	9 (2%) 55 43	18, 27, 37, 43	0
All	All	635/680 (93%)	-0.22	14 (2%) 65 54	14, 25, 36, 43	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	553	THR	5.8
1	B	552	TYR	5.5
1	B	550	ASP	3.6
1	B	551	ASP	3.5
1	B	568	TYR	3.3
1	B	594	LEU	3.1
1	A	551	ASP	3.1
1	A	552	TYR	2.8
1	B	567	GLU	2.6
1	A	343	SER	2.5
1	B	569	THR	2.4
1	A	482	ASP	2.4
1	B	548	ASN	2.3
1	A	404	LYS	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no 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.