



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 01:40 PM UTC

PDB ID : 1Q85 / pdb_00001q85
Title : Cobalamin-dependent methionine synthase (1-566) from *Thermotoga maritima* (Cd²⁺ complex, Se-Met)
Authors : Evans, J.C.; Huddler, D.P.; Hilgers, M.T.; Romanchuk, G.; Matthews, R.G.; Ludwig, M.L.
Deposited on : 2003-08-20
Resolution : 2.00 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

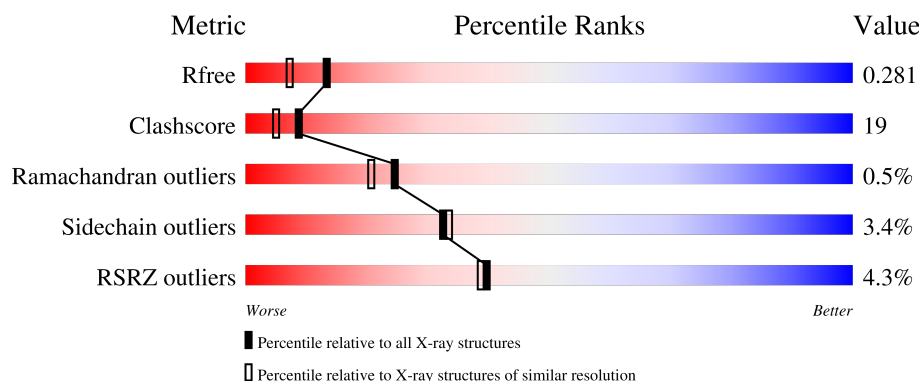
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	566	3% 69% 28% ..
1	B	566	5% 53% 42% ..

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8923 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 5-methyltetrahydrofolate S-homocysteine methyltransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	559	Total	C	N	O	S	Se	0	0	0
			4423	2839	738	833	3	10			
1	B	550	Total	C	N	O	S	Se	0	0	0
			4353	2795	726	819	3	10			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	modified residue	UNP Q9WYA5
A	27	MSE	MET	modified residue	UNP Q9WYA5
A	71	MSE	MET	modified residue	UNP Q9WYA5
A	135	MSE	MET	modified residue	UNP Q9WYA5
A	174	MSE	MET	modified residue	UNP Q9WYA5
A	334	MSE	MET	modified residue	UNP Q9WYA5
A	422	MSE	MET	modified residue	UNP Q9WYA5
A	439	MSE	MET	modified residue	UNP Q9WYA5
A	537	MSE	MET	modified residue	UNP Q9WYA5
A	545	MSE	MET	modified residue	UNP Q9WYA5
B	1	MSE	MET	modified residue	UNP Q9WYA5
B	27	MSE	MET	modified residue	UNP Q9WYA5
B	71	MSE	MET	modified residue	UNP Q9WYA5
B	135	MSE	MET	modified residue	UNP Q9WYA5
B	174	MSE	MET	modified residue	UNP Q9WYA5
B	334	MSE	MET	modified residue	UNP Q9WYA5
B	422	MSE	MET	modified residue	UNP Q9WYA5
B	439	MSE	MET	modified residue	UNP Q9WYA5
B	537	MSE	MET	modified residue	UNP Q9WYA5
B	545	MSE	MET	modified residue	UNP Q9WYA5

- Molecule 2 is CADMIUM ION (CCD ID: CD) (formula: Cd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total 2	Cd 2	0	0
2	B	1	Total 1	Cd 1	0	0

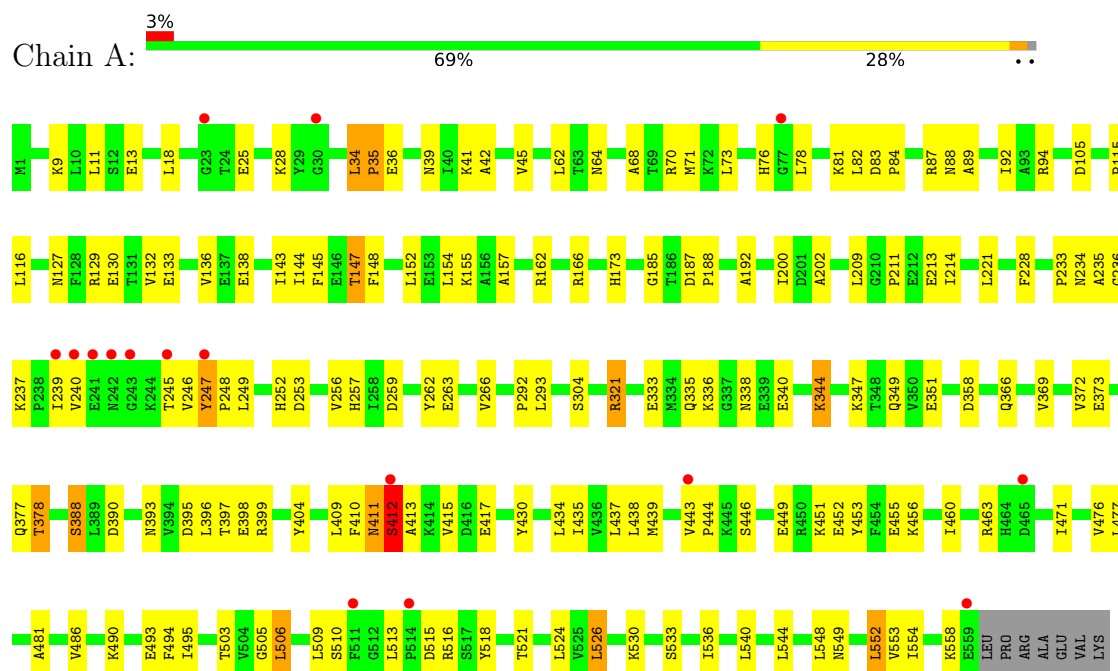
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	105	Total 105	O 105	0	0
3	B	39	Total 39	O 39	0	0

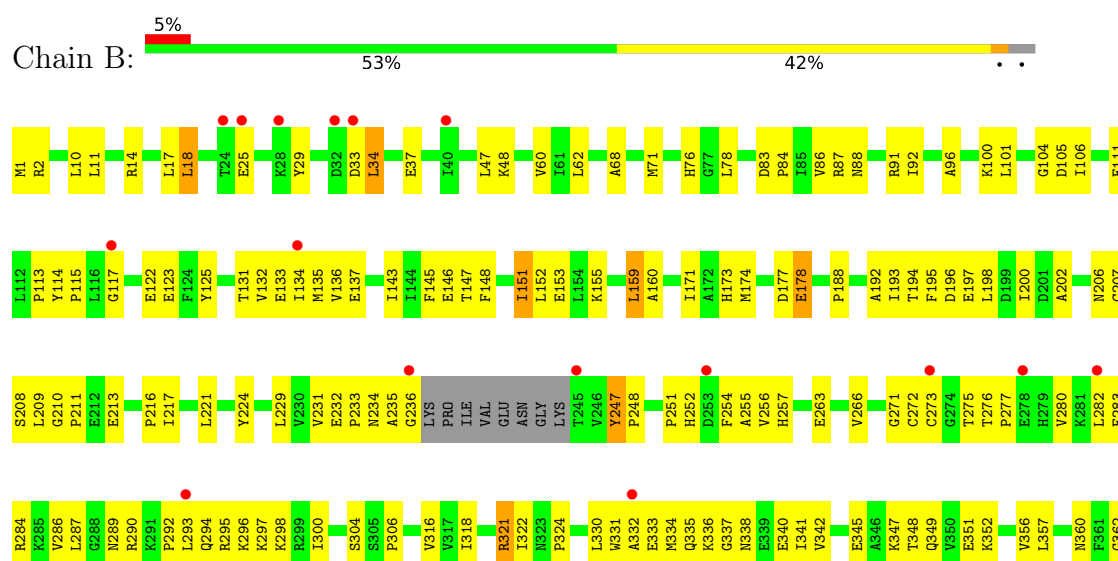
3 Residue-property plots

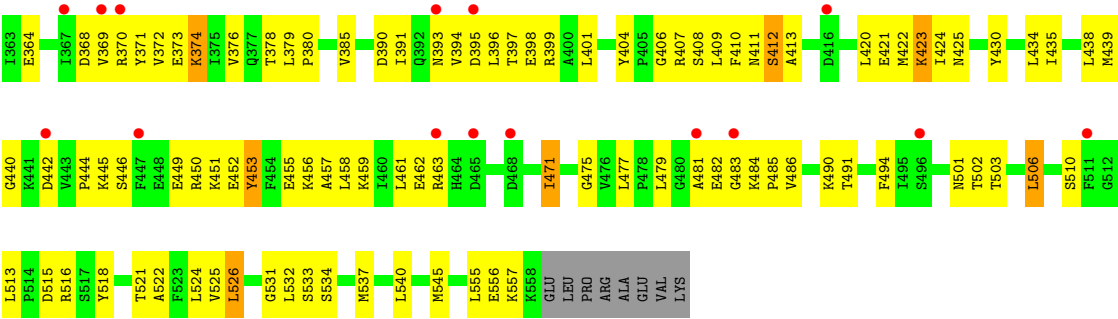
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: 5-methyltetrahydrofolate S-homocysteine methyltransferase



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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	60.30Å 86.17Å 125.88Å 90.00° 100.39° 90.00°	Depositor
Resolution (Å)	20.00 – 2.00 20.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-2.00) 90.8 (20.00-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.48 (at 1.90Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.243 , 0.283 0.242 , 0.281	Depositor DCC
R_{free} test set	7884 reflections (9.01%)	wwPDB-VP
Wilson B-factor (Å ²)	28.8	Xtriage
Anisotropy	0.470	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.45 , 68.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8923	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.68% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

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	0.45	0/4496	0.98	19/6056 (0.3%)
1	B	0.35	0/4424	0.90	9/5958 (0.2%)
All	All	0.40	0/8920	0.94	28/12014 (0.2%)

There are no bond length outliers.

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	321	ARG	N-CA-C	9.79	123.17	111.33
1	A	378	THR	N-CA-C	9.02	120.88	111.14
1	B	247	TYR	CA-C-N	8.68	128.41	119.56
1	B	247	TYR	C-N-CA	8.68	128.41	119.56
1	A	247	TYR	CA-C-N	7.37	127.01	119.56
1	A	247	TYR	C-N-CA	7.37	127.01	119.56
1	A	68	ALA	N-CA-C	6.98	120.38	110.68
1	B	378	THR	N-CA-C	6.94	118.64	111.14
1	A	34	LEU	CA-C-N	6.85	128.41	119.84
1	A	34	LEU	C-N-CA	6.85	128.41	119.84
1	A	304	SER	N-CA-C	6.70	118.95	108.96
1	A	248	PRO	N-CA-C	6.61	122.12	113.86
1	A	234	ASN	N-CA-C	-6.46	101.83	110.55
1	B	210	GLY	CA-C-N	6.39	125.88	119.24
1	B	210	GLY	C-N-CA	6.39	125.88	119.24
1	A	536	ILE	N-CA-C	-6.37	98.49	108.23
1	A	506	LEU	N-CA-C	6.12	118.49	111.02
1	A	116	LEU	N-CA-C	-5.86	104.49	111.69
1	B	304	SER	N-CA-C	5.82	117.63	108.96
1	B	557	LYS	N-CA-C	-5.81	104.86	112.41
1	B	321	ARG	N-CA-C	5.63	118.18	111.71
1	B	68	ALA	N-CA-C	5.52	118.16	109.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	412	SER	N-CA-C	5.40	122.30	110.80
1	A	147	THR	N-CA-C	5.35	119.08	112.24
1	A	388	SER	N-CA-C	-5.32	98.57	107.99
1	A	505	GLY	N-CA-C	-5.32	100.58	113.18
1	A	185	GLY	N-CA-C	5.20	123.23	115.64
1	A	340	GLU	N-CA-C	5.13	117.53	111.33

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	4423	0	4510	118	0
1	B	4353	0	4436	217	0
2	A	2	0	0	0	0
2	B	1	0	0	0	0
3	A	105	0	0	10	0
3	B	39	0	0	10	0
All	All	8923	0	8946	332	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 19.

All (332) 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:106:ILE:HD11	1:B:135:MSE:HE1	1.45	0.96
1:B:34:LEU:HD22	1:B:34:LEU:H	1.32	0.95
1:B:131:THR:HG22	1:B:135:MSE:HE2	1.48	0.94
1:A:515:ASP:HB3	1:A:518:TYR:HD1	1.34	0.91
1:B:482:GLU:HB3	1:B:484:LYS:HE2	1.53	0.89
1:B:131:THR:HG22	1:B:135:MSE:CE	2.04	0.87
1:B:459:LYS:HB3	1:B:463:ARG:HH12	1.38	0.86
1:B:435:ILE:HD13	1:B:471:ILE:HG23	1.56	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:273:CYS:SG	3:B:634:HOH:O	2.27	0.84
1:B:47:LEU:HD12	1:B:96:ALA:HB2	1.61	0.82
1:A:411:ASN:C	1:A:411:ASN:HD22	1.87	0.81
1:B:372:VAL:HG21	1:B:396:LEU:HD21	1.64	0.80
1:A:39:ASN:ND2	1:A:64:ASN:HD22	1.80	0.80
1:A:515:ASP:HB3	1:A:518:TYR:CD1	2.17	0.79
1:B:458:LEU:O	1:B:462:GLU:HG2	1.83	0.77
1:B:409:LEU:HD21	1:B:435:ILE:HD11	1.66	0.77
1:B:271:GLY:HA3	1:B:275:THR:HG21	1.68	0.76
1:A:39:ASN:HD21	1:A:64:ASN:HD22	1.34	0.75
1:A:411:ASN:HD22	1:A:412:SER:N	1.84	0.75
1:A:247:TYR:HD1	3:A:690:HOH:O	1.68	0.75
1:B:435:ILE:HD13	1:B:471:ILE:CG2	2.16	0.75
1:B:522:ALA:O	1:B:526:LEU:HD22	1.86	0.75
1:B:83:ASP:HB3	1:B:84:PRO:HD3	1.70	0.73
1:B:333:GLU:HB3	1:B:338:ASN:HB3	1.71	0.72
1:B:449:GLU:O	1:B:452:GLU:HG2	1.90	0.72
1:B:287:LEU:HD12	1:B:290:ARG:HE	1.54	0.71
1:B:151:ILE:HD11	1:B:306:PRO:HG3	1.71	0.71
1:A:89:ALA:HA	3:A:616:HOH:O	1.90	0.71
1:B:234:ASN:HD22	1:B:272:CYS:HB3	1.57	0.70
1:A:73:LEU:HD12	1:A:82:LEU:HD23	1.72	0.70
1:A:493:GLU:HB2	1:A:530:LYS:HG3	1.73	0.70
1:B:60:VAL:HA	3:B:606:HOH:O	1.93	0.69
1:B:159:LEU:HD23	3:B:639:HOH:O	1.92	0.69
1:B:442:ASP:O	1:B:444:PRO:HD3	1.93	0.68
1:B:284:ARG:HB2	3:B:619:HOH:O	1.93	0.67
1:A:94:ARG:HG3	1:A:94:ARG:HH11	1.59	0.67
1:B:234:ASN:ND2	1:B:272:CYS:HB3	2.10	0.67
1:A:235:ALA:C	1:A:249:LEU:HB2	2.21	0.66
1:B:132:VAL:HA	1:B:135:MSE:HE3	1.77	0.66
1:B:2:ARG:NH2	1:B:101:LEU:HD23	2.11	0.66
1:B:330:LEU:HG	1:B:334:MSE:HE3	1.78	0.65
1:A:413:ALA:HB2	1:A:434:LEU:HD11	1.78	0.65
1:B:34:LEU:HD23	1:B:37:GLU:OE2	1.97	0.65
1:B:114:TYR:CD1	1:B:115:PRO:HA	2.32	0.64
1:B:515:ASP:HB3	1:B:518:TYR:HD1	1.62	0.64
1:A:347:LYS:O	1:A:351:GLU:HG3	1.98	0.64
1:B:151:ILE:HD12	1:B:152:LEU:H	1.63	0.64
1:B:178:GLU:H	1:B:178:GLU:CD	2.01	0.64
1:B:11:LEU:O	1:B:292:PRO:HG2	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:132:VAL:HG13	1:B:143:ILE:CD1	2.28	0.63
1:B:208:SER:O	1:B:209:LEU:HD23	1.99	0.62
1:A:192:ALA:HB2	1:A:221:LEU:HD12	1.81	0.62
1:B:412:SER:H	1:B:435:ILE:HB	1.65	0.62
1:B:229:LEU:HG	1:B:295:ARG:NH2	2.15	0.62
1:A:372:VAL:HG21	1:A:396:LEU:HD21	1.82	0.61
1:B:151:ILE:HD12	1:B:152:LEU:N	2.14	0.61
1:A:411:ASN:C	1:A:411:ASN:ND2	2.55	0.61
1:B:445:LYS:HD2	1:B:445:LYS:N	2.15	0.61
1:B:47:LEU:CD1	1:B:96:ALA:HB2	2.30	0.60
1:B:396:LEU:HD13	1:B:396:LEU:O	2.01	0.60
1:A:252:HIS:O	1:A:256:VAL:HG13	2.01	0.60
1:B:235:ALA:HB3	1:B:247:TYR:HB3	1.83	0.60
1:A:443:VAL:HG11	1:A:477:LEU:HD21	1.84	0.60
1:B:71:MSE:HE3	1:B:123:GLU:HG2	1.84	0.59
1:B:193:ILE:HD11	1:B:300:ILE:HD11	1.83	0.59
1:B:486:VAL:O	1:B:490:LYS:HG3	2.03	0.59
1:B:368:ASP:O	1:B:371:TYR:HB3	2.02	0.59
1:A:321:ARG:NE	1:A:540:LEU:HD22	2.17	0.59
1:A:88:ASN:O	1:A:92:ILE:HG13	2.02	0.59
1:B:410:PHE:HE2	1:B:423:LYS:HD3	1.67	0.59
1:B:193:ILE:CD1	1:B:300:ILE:HD11	2.33	0.58
1:B:148:PHE:HE2	1:B:153:GLU:HG2	1.69	0.58
1:B:502:THR:H	1:B:533:SER:HB2	1.68	0.58
1:B:510:SER:OG	1:B:516:ARG:HB2	2.03	0.58
1:B:251:PRO:CB	1:B:282:LEU:HD12	2.32	0.57
1:B:372:VAL:O	1:B:376:VAL:HG23	2.04	0.57
1:B:411:ASN:HB3	3:B:625:HOH:O	2.04	0.57
1:B:76:HIS:O	1:B:78:LEU:HD22	2.04	0.57
1:B:200:ILE:C	1:B:200:ILE:HD12	2.30	0.56
1:B:321:ARG:HD2	1:B:540:LEU:HD13	1.86	0.56
1:B:372:VAL:HG21	1:B:396:LEU:CD2	2.36	0.56
1:B:390:ASP:HA	1:B:411:ASN:HB3	1.88	0.56
1:B:413:ALA:HB2	1:B:434:LEU:HD11	1.87	0.56
1:B:293:LEU:HD23	1:B:294:GLN:O	2.06	0.56
1:B:444:PRO:HB3	1:B:449:GLU:HB3	1.87	0.56
1:B:391:ILE:HD13	1:B:396:LEU:HD12	1.88	0.56
1:A:344:LYS:HB2	1:A:344:LYS:NZ	2.21	0.56
1:B:321:ARG:NE	1:B:540:LEU:HD22	2.21	0.56
1:B:100:LYS:HB3	3:B:606:HOH:O	2.07	0.55
1:B:368:ASP:OD2	1:B:370:ARG:HG2	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:ARG:HH11	1:A:129:ARG:HG2	1.72	0.55
1:B:479:LEU:HD13	1:B:485:PRO:HG2	1.87	0.55
1:A:76:HIS:HB2	1:A:78:LEU:HD23	1.87	0.54
1:A:554:ILE:HG23	1:B:485:PRO:HB2	1.89	0.54
1:B:145:PHE:O	1:B:173:HIS:HB2	2.08	0.54
1:A:393:ASN:ND2	1:A:395:ASP:H	2.05	0.54
1:B:251:PRO:HB3	1:B:282:LEU:HD12	1.88	0.54
1:A:81:LYS:C	1:A:84:PRO:HD2	2.32	0.54
1:A:240:VAL:HG22	1:A:245:THR:HG22	1.89	0.54
1:A:71:MSE:H	1:A:127:ASN:HD21	1.56	0.54
1:A:213:GLU:HG3	3:A:638:HOH:O	2.08	0.53
1:B:484:LYS:N	1:B:485:PRO:HD3	2.22	0.53
1:B:450:ARG:HD2	1:B:491:THR:OG1	2.09	0.53
1:B:451:LYS:HE3	1:B:494:PHE:CZ	2.42	0.53
1:A:411:ASN:O	1:A:412:SER:HB2	2.08	0.53
1:B:484:LYS:N	1:B:484:LYS:HD2	2.24	0.53
1:A:132:VAL:HG13	1:A:143:ILE:CD1	2.39	0.53
1:A:446:SER:OG	1:A:449:GLU:HG3	2.09	0.53
1:A:94:ARG:HG3	1:A:94:ARG:NH1	2.24	0.52
1:B:171:ILE:HG12	1:B:202:ALA:HB3	1.91	0.52
1:A:526:LEU:HD13	1:B:525:VAL:HG12	1.91	0.52
1:B:510:SER:CB	1:B:516:ARG:HB2	2.39	0.52
1:B:17:LEU:HD11	1:B:283:PHE:HB3	1.91	0.52
1:B:25:GLU:OE1	1:B:277:PRO:HD3	2.09	0.52
1:A:71:MSE:H	1:A:127:ASN:ND2	2.07	0.52
1:B:457:ALA:O	1:B:461:LEU:HD13	2.09	0.52
1:A:83:ASP:HB3	1:A:84:PRO:HD3	1.91	0.52
1:B:115:PRO:C	1:B:117:GLY:H	2.16	0.52
1:B:336:LYS:C	1:B:336:LYS:HD2	2.35	0.52
1:B:397:THR:O	1:B:401:LEU:HD23	2.10	0.52
1:B:177:ASP:HB2	1:B:178:GLU:OE1	2.11	0.51
1:A:144:ILE:HD11	1:A:173:HIS:CE1	2.46	0.51
1:B:537:MSE:HE2	1:B:545:MSE:SE	2.61	0.51
1:B:125:TYR:CE1	1:B:160:ALA:HA	2.45	0.51
1:A:202:ALA:HB2	1:A:228:PHE:HB2	1.92	0.51
1:B:341:ILE:HG23	1:B:342:VAL:N	2.25	0.51
1:B:360:ASN:ND2	1:B:362:GLY:H	2.09	0.50
1:B:76:HIS:HB2	1:B:78:LEU:HD23	1.93	0.50
1:B:446:SER:HB3	1:B:449:GLU:HG3	1.93	0.50
1:B:131:THR:O	1:B:135:MSE:HE3	2.12	0.50
1:B:113:PRO:HB3	1:B:152:LEU:HD23	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:336:LYS:HE3	1:B:338:ASN:N	2.27	0.50
1:A:9:LYS:HE2	1:A:13:GLU:OE2	2.11	0.50
1:B:379:LEU:HB2	1:B:380:PRO:HD3	1.93	0.50
1:B:521:THR:O	1:B:525:VAL:HG23	2.12	0.50
1:B:555:LEU:O	1:B:556:GLU:HG2	2.12	0.50
1:B:197:GLU:OE1	1:B:407:ARG:HD2	2.12	0.50
1:A:129:ARG:HG2	1:A:129:ARG:NH1	2.26	0.50
1:B:393:ASN:HD22	1:B:393:ASN:C	2.20	0.50
1:A:81:LYS:O	1:A:84:PRO:HD2	2.12	0.49
1:B:421:GLU:HG2	1:B:425:ASN:ND2	2.27	0.49
1:B:87:ARG:O	1:B:91:ARG:HG3	2.13	0.49
1:B:347:LYS:O	1:B:351:GLU:HG3	2.12	0.49
1:B:398:GLU:HG3	1:B:430:TYR:CE2	2.47	0.49
1:B:477:LEU:HB2	1:B:483:GLY:HA2	1.95	0.49
1:A:549:ASN:O	1:A:553:VAL:HG23	2.13	0.49
1:B:463:ARG:HG3	1:B:463:ARG:HH11	1.77	0.49
1:B:421:GLU:HG2	1:B:425:ASN:HD21	1.77	0.49
1:A:214:ILE:HD12	1:A:233:PRO:HB3	1.95	0.49
1:B:276:THR:HB	1:B:277:PRO:CD	2.43	0.49
1:B:369:VAL:O	1:B:372:VAL:HG22	2.13	0.49
1:A:145:PHE:CE2	1:A:157:ALA:HB1	2.48	0.48
1:A:552:LEU:HB3	1:A:558:LYS:HG2	1.95	0.48
1:B:446:SER:HB3	1:B:449:GLU:CG	2.43	0.48
1:A:369:VAL:O	1:A:373:GLU:HG3	2.14	0.48
1:A:444:PRO:HG3	1:A:453:TYR:CE1	2.48	0.48
1:B:213:GLU:O	1:B:216:PRO:HD2	2.14	0.48
1:B:393:ASN:ND2	1:B:395:ASP:H	2.12	0.48
1:A:87:ARG:HA	1:A:138:GLU:HG3	1.95	0.48
1:A:398:GLU:HG3	1:A:430:TYR:CE2	2.48	0.48
1:B:287:LEU:HD12	1:B:290:ARG:HH11	1.79	0.48
1:A:471:ILE:HD11	1:A:503:THR:HB	1.94	0.48
1:B:506:LEU:HD11	1:B:524:LEU:HB2	1.96	0.48
1:B:132:VAL:O	1:B:136:VAL:HG23	2.14	0.48
1:A:155:LYS:NZ	3:A:685:HOH:O	2.46	0.48
1:B:231:VAL:HG22	1:B:233:PRO:HD3	1.95	0.48
1:B:207:CYS:SG	1:B:208:SER:N	2.84	0.47
1:B:348:THR:O	1:B:352:LYS:HG2	2.14	0.47
1:B:510:SER:HB2	1:B:516:ARG:HB2	1.97	0.47
1:B:211:PRO:HG3	1:B:254:PHE:CE1	2.49	0.47
1:A:36:GLU:H	1:A:36:GLU:CD	2.22	0.47
1:A:548:LEU:O	1:A:552:LEU:HD22	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:115:PRO:C	1:B:117:GLY:N	2.70	0.47
1:B:231:VAL:HG12	1:B:266:VAL:HG21	1.97	0.47
1:A:129:ARG:O	1:A:133:GLU:HG3	2.15	0.47
1:B:29:TYR:OH	1:B:48:LYS:HE2	2.14	0.47
1:A:397:THR:HG23	1:A:410:PHE:CE1	2.49	0.47
1:B:555:LEU:C	1:B:556:GLU:HG2	2.40	0.47
1:A:321:ARG:HB2	1:A:349:GLN:CD	2.40	0.46
1:A:166:ARG:HG2	1:A:166:ARG:HH11	1.80	0.46
1:B:133:GLU:O	1:B:137:GLU:HG3	2.14	0.46
1:B:235:ALA:HB3	1:B:247:TYR:CB	2.45	0.46
1:B:321:ARG:HG2	1:B:321:ARG:HH11	1.81	0.46
1:B:391:ILE:HD13	1:B:396:LEU:CD1	2.45	0.46
1:A:36:GLU:O	1:A:39:ASN:HB2	2.16	0.46
1:B:252:HIS:O	1:B:256:VAL:HG13	2.14	0.46
1:A:333:GLU:OE1	1:A:336:LYS:HD2	2.16	0.46
1:B:445:LYS:HD2	1:B:445:LYS:H	1.80	0.46
1:A:249:LEU:HD22	1:A:253:ASP:HB2	1.98	0.46
1:A:259:ASP:O	1:A:263:GLU:HG3	2.16	0.46
1:A:415:VAL:O	1:A:415:VAL:HG12	2.16	0.46
1:A:451:LYS:HE3	1:A:494:PHE:CZ	2.51	0.46
1:B:122:GLU:OE1	1:B:122:GLU:HA	2.16	0.46
1:B:296:LYS:O	1:B:297:LYS:HB3	2.16	0.46
1:A:73:LEU:CD1	1:A:82:LEU:HD23	2.45	0.46
1:B:211:PRO:HB2	1:B:257:HIS:CD2	2.51	0.46
1:B:287:LEU:CD1	1:B:290:ARG:HH11	2.28	0.46
1:B:331:TRP:O	1:B:335:GLN:HG2	2.16	0.46
1:B:444:PRO:HG2	1:B:453:TYR:CE1	2.50	0.46
1:A:321:ARG:CZ	1:A:540:LEU:HD22	2.46	0.46
1:B:324:PRO:HG3	1:B:334:MSE:SE	2.65	0.46
1:A:76:HIS:O	1:A:78:LEU:HD22	2.16	0.45
1:A:510:SER:HA	1:A:513:LEU:HD12	1.97	0.45
1:B:18:LEU:HD13	1:B:232:GLU:CD	2.41	0.45
1:B:62:LEU:HB3	1:B:105:ASP:HB2	1.98	0.45
1:B:83:ASP:CG	1:B:87:ARG:HE	2.25	0.45
1:B:151:ILE:HG22	1:B:195:PHE:HE1	1.82	0.45
1:A:438:LEU:HD11	1:A:495:ILE:HD11	1.98	0.45
1:B:263:GLU:O	1:B:293:LEU:HD13	2.16	0.45
1:A:41:LYS:HG3	1:A:42:ALA:N	2.31	0.45
1:A:209:LEU:HB3	1:A:213:GLU:HB2	1.99	0.45
1:B:398:GLU:HG3	1:B:430:TYR:HE2	1.81	0.45
1:B:486:VAL:HG12	1:B:490:LYS:HE3	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:444:PRO:HG3	1:A:453:TYR:HE1	1.81	0.45
1:A:417:GLU:HG2	1:A:460:ILE:CD1	2.47	0.45
1:A:526:LEU:HD13	1:B:525:VAL:CG1	2.46	0.45
1:B:321:ARG:HB2	1:B:349:GLN:NE2	2.31	0.45
1:B:1:MSE:SE	3:B:614:HOH:O	2.84	0.45
1:B:83:ASP:OD1	1:B:87:ARG:NE	2.43	0.45
1:B:132:VAL:HG13	1:B:143:ILE:HD11	1.99	0.45
1:B:200:ILE:HD12	1:B:200:ILE:O	2.17	0.45
1:B:209:LEU:HB3	1:B:213:GLU:HB2	1.97	0.45
1:A:148:PHE:CG	1:A:154:LEU:HD13	2.52	0.45
1:B:148:PHE:CE2	1:B:153:GLU:HG2	2.51	0.45
1:B:446:SER:HB3	1:B:449:GLU:CD	2.42	0.45
1:A:187:ASP:HB2	1:A:188:PRO:CD	2.47	0.44
1:B:318:ILE:HG12	1:B:356:VAL:HB	1.99	0.44
1:B:207:CYS:HA	1:B:234:ASN:ND2	2.32	0.44
1:B:336:LYS:HE3	1:B:338:ASN:CA	2.47	0.44
1:B:88:ASN:O	1:B:92:ILE:HG13	2.16	0.44
1:B:192:ALA:O	1:B:196:ASP:HB2	2.18	0.44
1:B:471:ILE:HG13	1:B:501:ASN:C	2.43	0.44
1:B:502:THR:O	1:B:533:SER:HB2	2.18	0.44
1:A:42:ALA:HB1	1:A:45:VAL:HG23	1.98	0.44
1:A:393:ASN:OD1	1:A:395:ASP:HB3	2.17	0.44
1:A:404:TYR:HB3	3:A:685:HOH:O	2.18	0.44
1:A:415:VAL:CG1	1:A:456:LYS:HD3	2.47	0.44
1:A:510:SER:OG	1:A:516:ARG:HB2	2.17	0.44
1:B:146:GLU:OE2	1:B:232:GLU:OE2	2.36	0.44
1:A:62:LEU:HB3	1:A:105:ASP:HB2	2.00	0.44
1:A:70:ARG:CD	1:A:130:GLU:OE2	2.66	0.44
1:A:25:GLU:O	1:A:28:LYS:HB3	2.18	0.43
1:A:39:ASN:ND2	3:A:616:HOH:O	2.50	0.43
1:A:435:ILE:HA	1:A:471:ILE:HG23	2.00	0.43
1:A:476:VAL:HG11	1:A:509:LEU:HB2	2.00	0.43
1:B:364:GLU:OE1	1:B:391:ILE:HG23	2.17	0.43
1:B:479:LEU:HB3	1:B:513:LEU:HD11	2.00	0.43
1:B:369:VAL:C	1:B:371:TYR:H	2.26	0.43
1:B:321:ARG:CD	1:B:540:LEU:HD22	2.49	0.43
1:B:330:LEU:CG	1:B:334:MSE:HE3	2.45	0.43
1:B:321:ARG:HG2	1:B:321:ARG:NH1	2.34	0.43
1:B:345:GLU:HA	1:B:345:GLU:OE1	2.18	0.43
1:B:357:LEU:HG	1:B:385:VAL:HG21	1.99	0.43
1:B:393:ASN:C	1:B:393:ASN:ND2	2.75	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:510:SER:HB2	1:B:516:ARG:CB	2.48	0.43
1:B:280:VAL:HA	1:B:283:PHE:CD2	2.53	0.43
1:B:374:LYS:CB	1:B:374:LYS:NZ	2.81	0.43
1:B:393:ASN:HD22	1:B:394:VAL:N	2.16	0.43
1:B:479:LEU:C	1:B:481:ALA:H	2.26	0.43
1:B:255:ALA:HB1	1:B:286:VAL:HG21	1.99	0.43
1:B:296:LYS:HB3	1:B:297:LYS:H	1.55	0.43
1:B:456:LYS:HA	1:B:459:LYS:NZ	2.32	0.43
1:B:276:THR:HB	1:B:277:PRO:HD2	2.01	0.43
1:A:34:LEU:HD12	1:A:34:LEU:O	2.18	0.43
1:A:409:LEU:HD11	1:A:435:ILE:HG13	2.01	0.43
1:B:145:PHE:HB3	1:B:148:PHE:CE1	2.54	0.43
1:B:462:GLU:OE2	1:B:462:GLU:HA	2.19	0.43
1:A:115:PRO:HG3	1:A:378:THR:HG23	2.01	0.43
1:A:152:LEU:CD2	1:A:377:GLN:HG3	2.49	0.43
1:B:420:LEU:O	1:B:424:ILE:HG13	2.18	0.43
1:A:132:VAL:O	1:A:136:VAL:HG23	2.18	0.42
1:A:162:ARG:HD2	1:A:166:ARG:HA	2.00	0.42
1:A:292:PRO:HG3	3:A:606:HOH:O	2.19	0.42
1:B:34:LEU:HD23	1:B:37:GLU:CD	2.44	0.42
1:B:151:ILE:HG23	1:B:174:MSE:HE1	2.00	0.42
1:B:336:LYS:HG3	1:B:337:GLY:N	2.34	0.42
1:B:14:ARG:HH21	1:B:289:ASN:ND2	2.16	0.42
1:B:332:ALA:O	1:B:335:GLN:HB2	2.19	0.42
1:A:132:VAL:HG13	1:A:143:ILE:HD12	2.01	0.42
1:A:390:ASP:HA	1:A:411:ASN:HB3	2.01	0.42
1:A:437:LEU:HG	1:A:439:MSE:HG2	2.00	0.42
1:B:206:ASN:HB2	1:B:232:GLU:HB2	2.00	0.42
1:A:451:LYS:O	1:A:455:GLU:HG3	2.20	0.42
1:B:316:VAL:O	1:B:534:SER:HB2	2.20	0.42
1:B:404:TYR:CD1	1:B:408:SER:HB3	2.54	0.42
1:A:333:GLU:HB3	1:A:338:ASN:HB3	2.01	0.42
1:B:188:PRO:HG3	1:B:217:ILE:HG23	2.02	0.42
1:A:70:ARG:HD3	1:A:130:GLU:OE2	2.19	0.42
1:A:213:GLU:CD	1:A:237:LYS:HE3	2.44	0.42
1:B:247:TYR:HA	1:B:248:PRO:HD2	1.80	0.42
1:A:443:VAL:HA	1:A:444:PRO:HD3	1.90	0.42
1:A:486:VAL:O	1:A:490:LYS:HG3	2.20	0.42
1:B:224:TYR:CZ	1:B:298:LYS:HD2	2.55	0.42
1:A:11:LEU:O	1:A:292:PRO:HG2	2.19	0.42
1:A:34:LEU:HA	1:A:35:PRO:HD2	1.76	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:88:ASN:ND2	3:A:672:HOH:O	2.53	0.42
1:B:515:ASP:HB3	1:B:518:TYR:CD1	2.49	0.42
1:A:358:ASP:HA	1:A:388:SER:HB3	2.02	0.42
1:B:322:ILE:O	1:B:322:ILE:HG22	2.19	0.42
1:B:404:TYR:CE1	1:B:408:SER:HB3	2.55	0.42
1:B:86:VAL:HB	1:B:134:ILE:HG21	2.01	0.41
1:A:209:LEU:HB2	1:A:214:ILE:HG13	2.02	0.41
1:B:234:ASN:C	1:B:236:GLY:H	2.27	0.41
1:B:251:PRO:HB2	1:B:282:LEU:HD12	2.01	0.41
1:B:34:LEU:H	1:B:34:LEU:CD2	2.11	0.41
1:B:159:LEU:HG	3:B:640:HOH:O	2.20	0.41
1:A:41:LYS:HE3	1:A:41:LYS:HB2	1.87	0.41
1:A:162:ARG:HA	1:A:162:ARG:HD3	1.72	0.41
1:A:521:THR:HA	1:A:544:LEU:CD1	2.51	0.41
1:B:104:GLY:HA3	1:B:135:MSE:SE	2.71	0.41
1:B:188:PRO:CG	1:B:217:ILE:HG23	2.50	0.41
1:A:395:ASP:O	1:A:399:ARG:HD3	2.20	0.41
1:A:239:ILE:O	1:A:246:VAL:N	2.48	0.41
1:A:262:TYR:HA	1:A:266:VAL:HG12	2.01	0.41
1:B:373:GLU:OE1	1:B:399:ARG:HD2	2.19	0.41
1:A:39:ASN:CG	3:A:616:HOH:O	2.64	0.41
1:B:287:LEU:HD12	1:B:290:ARG:NE	2.29	0.41
1:A:443:VAL:HG11	1:A:477:LEU:CD2	2.50	0.41
1:B:106:ILE:CD1	1:B:135:MSE:HE1	2.31	0.41
1:B:151:ILE:HG13	3:B:617:HOH:O	2.20	0.41
1:B:194:THR:HG21	1:B:406:GLY:HA2	2.02	0.41
1:B:422:MSE:HE3	1:B:423:LYS:NZ	2.36	0.41
1:B:439:MSE:HG2	1:B:440:GLY:N	2.35	0.41
1:B:458:LEU:O	1:B:458:LEU:HD23	2.21	0.41
1:B:463:ARG:HG3	1:B:463:ARG:NH1	2.36	0.41
1:A:236:GLY:O	3:A:690:HOH:O	2.22	0.40
1:B:531:GLY:O	1:B:532:LEU:C	2.64	0.40
1:B:435:ILE:HD12	1:B:503:THR:HG21	2.02	0.40
1:A:211:PRO:HB2	1:A:257:HIS:CD2	2.56	0.40
1:B:88:ASN:N	1:B:88:ASN:HD22	2.18	0.40
1:B:155:LYS:HE2	1:B:404:TYR:O	2.21	0.40
1:B:111:GLU:HB2	3:B:608:HOH:O	2.21	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	557/566 (98%)	533 (96%)	21 (4%)	3 (0%)	24	21
1	B	546/566 (96%)	505 (92%)	38 (7%)	3 (0%)	24	21
All	All	1103/1132 (97%)	1038 (94%)	59 (5%)	6 (0%)	24	21

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	481	ALA
1	B	412	SER
1	A	412	SER
1	B	438	LEU
1	B	475	GLY
1	A	35	PRO

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	488/484 (101%)	473 (97%)	15 (3%)	35	37
1	B	480/484 (99%)	462 (96%)	18 (4%)	29	29
All	All	968/968 (100%)	935 (97%)	33 (3%)	32	33

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	LEU
1	A	147	THR
1	A	200	ILE
1	A	293	LEU
1	A	335	GLN
1	A	344	LYS
1	A	366	GLN
1	A	411	ASN
1	A	452	GLU
1	A	463	ARG
1	A	506	LEU
1	A	524	LEU
1	A	526	LEU
1	A	533	SER
1	A	552	LEU
1	B	10	LEU
1	B	18	LEU
1	B	33	ASP
1	B	34	LEU
1	B	147	THR
1	B	151	ILE
1	B	159	LEU
1	B	178	GLU
1	B	198	LEU
1	B	221	LEU
1	B	340	GLU
1	B	374	LYS
1	B	423	LYS
1	B	453	TYR
1	B	455	GLU
1	B	471	ILE
1	B	506	LEU
1	B	526	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	ASN
1	A	127	ASN
1	A	257	HIS
1	A	335	GLN
1	A	392	GLN
1	A	411	ASN

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Mol	Chain	Res	Type
1	A	508	ASN
1	A	549	ASN
1	B	64	ASN
1	B	76	HIS
1	B	88	ASN
1	B	219	GLN
1	B	234	ASN
1	B	257	HIS
1	B	289	ASN
1	B	360	ASN
1	B	393	ASN
1	B	425	ASN
1	B	501	ASN
1	B	520	ASN

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 3 ligands modelled in this entry, 3 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

There are no such residues in this entry.

5.8 Polymer linkage issues

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	549/566 (96%)	0.01	16 (2%) 53 52	17, 34, 57, 79	0
1	B	540/566 (95%)	0.59	31 (5%) 29 28	32, 50, 74, 79	0
All	All	1089/1132 (96%)	0.29	47 (4%) 40 39	17, 43, 71, 79	0

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	483	GLY	4.5
1	A	239	ILE	3.2
1	B	511	PHE	3.1
1	B	273	CYS	3.0
1	A	412	SER	3.0
1	A	77	GLY	2.8
1	B	245	THR	2.7
1	B	40	ILE	2.7
1	B	465	ASP	2.7
1	A	243	GLY	2.7
1	A	511	PHE	2.7
1	B	24	THR	2.6
1	A	240	VAL	2.6
1	B	25	GLU	2.6
1	B	468	ASP	2.6
1	B	282	LEU	2.6
1	B	253	ASP	2.5
1	B	496	SER	2.5
1	B	33	ASP	2.5
1	B	395	ASP	2.5
1	A	245	THR	2.4
1	A	443	VAL	2.4
1	B	442	ASP	2.4
1	A	514	PRO	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	278	GLU	2.3
1	A	242	ASN	2.3
1	A	23	GLY	2.3
1	B	463	ARG	2.3
1	B	117	GLY	2.3
1	B	370	ARG	2.2
1	B	369	VAL	2.2
1	B	293	LEU	2.2
1	A	241	GLU	2.1
1	A	559	GLU	2.1
1	A	30	GLY	2.1
1	B	332	ALA	2.1
1	B	367	ILE	2.1
1	B	28	LYS	2.1
1	B	481	ALA	2.1
1	B	416	ASP	2.1
1	B	236	GLY	2.1
1	A	465	ASP	2.0
1	B	32	ASP	2.0
1	A	247	TYR	2.0
1	B	134	ILE	2.0
1	B	447	PHE	2.0
1	B	393	ASN	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](#)

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.

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	CD	B	602	1/1	0.91	0.05	58,58,58,58	0
2	CD	A	603	1/1	0.95	0.05	45,45,45,45	0
2	CD	A	601	1/1	0.95	0.05	41,41,41,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.