



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 05:21 PM UTC

PDB ID : 2OND / pdb_00002ond
Title : Crystal Structure of the HAT-C domain of murine CstF-77
Authors : Bai, Y.; Auperin, T.C.; Chou, C.-Y.; Chang, G.-G.; Manley, J.L.; Tong, L.
Deposited on : 2007-01-23
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

<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

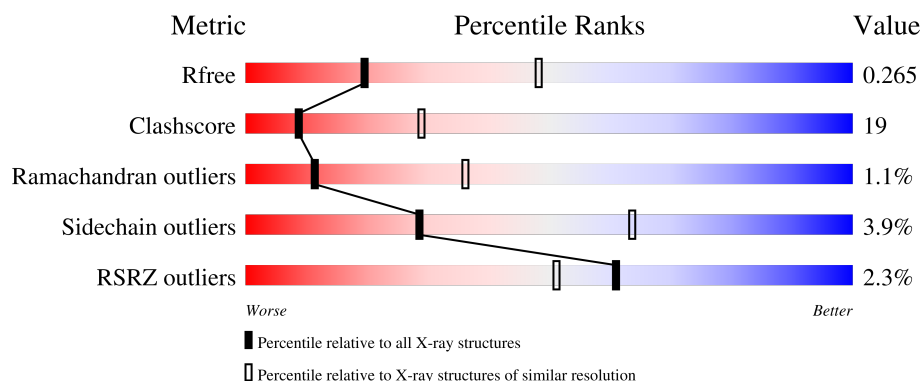
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (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 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	308	
1	B	308	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 5099 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 Cleavage stimulation factor 77 kDa subunit.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	308	Total	C	N	O	S	Se	0	0	0
			2549	1641	424	472	3	9			
1	B	308	Total	C	N	O	S	Se	0	0	0
			2550	1641	424	473	3	9			

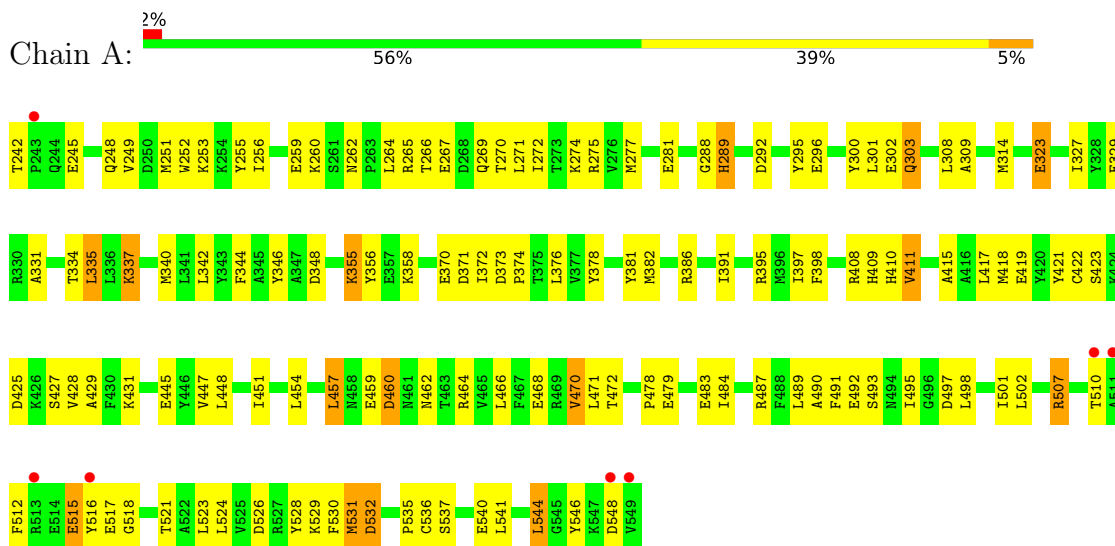
There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	251	MSE	MET	modified residue	UNP Q99LI7
A	277	MSE	MET	modified residue	UNP Q99LI7
A	314	MSE	MET	modified residue	UNP Q99LI7
A	340	MSE	MET	modified residue	UNP Q99LI7
A	354	MSE	MET	modified residue	UNP Q99LI7
A	382	MSE	MET	modified residue	UNP Q99LI7
A	396	MSE	MET	modified residue	UNP Q99LI7
A	418	MSE	MET	modified residue	UNP Q99LI7
A	531	MSE	MET	modified residue	UNP Q99LI7
B	251	MSE	MET	modified residue	UNP Q99LI7
B	277	MSE	MET	modified residue	UNP Q99LI7
B	314	MSE	MET	modified residue	UNP Q99LI7
B	340	MSE	MET	modified residue	UNP Q99LI7
B	354	MSE	MET	modified residue	UNP Q99LI7
B	382	MSE	MET	modified residue	UNP Q99LI7
B	396	MSE	MET	modified residue	UNP Q99LI7
B	418	MSE	MET	modified residue	UNP Q99LI7
B	531	MSE	MET	modified residue	UNP Q99LI7

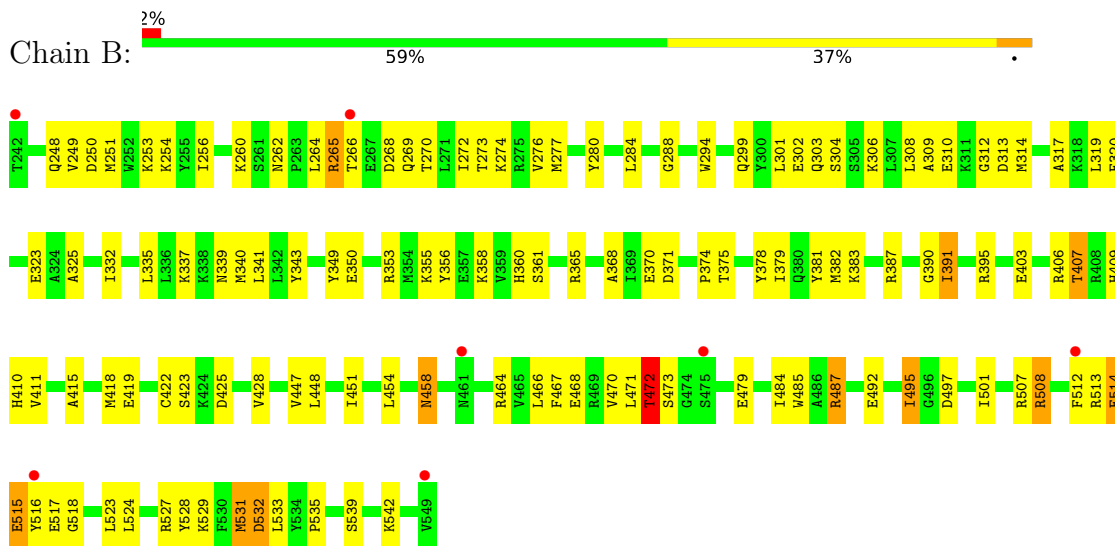
3 Residue-property plots

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: Cleavage stimulation factor 77 kDa subunit



• Molecule 1: Cleavage stimulation factor 77 kDa subunit



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	155.40Å 155.40Å 161.60Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.84 – 2.80 29.84 – 2.80	Depositor EDS
% Data completeness (in resolution range)	97.2 (29.84-2.80) 97.3 (29.84-2.80)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.71 (at 2.80Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.224 , 0.262 0.231 , 0.265	Depositor DCC
R_{free} test set	3946 reflections (7.38%)	wwPDB-VP
Wilson B-factor (Å ²)	60.9	Xtriage
Anisotropy	0.039	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 47.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5099	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.56% 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

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	0.47	0/2597	0.97	13/3484 (0.4%)
1	B	0.44	0/2598	0.94	12/3484 (0.3%)
All	All	0.45	0/5195	0.96	25/6968 (0.4%)

There are no bond length outliers.

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	370	GLU	N-CA-C	7.42	120.31	111.33
1	B	495	ILE	N-CA-C	7.36	119.96	112.90
1	A	373	ASP	CA-C-N	6.84	126.52	119.82
1	A	373	ASP	C-N-CA	6.84	126.52	119.82
1	A	518	GLY	N-CA-C	-6.62	106.10	115.43
1	B	518	GLY	N-CA-C	-6.44	106.02	115.30
1	B	390	GLY	N-CA-C	6.42	118.05	111.95
1	B	472	THR	N-CA-C	6.33	121.24	112.45
1	B	370	GLU	N-CA-C	6.25	117.79	110.97
1	B	356	TYR	N-CA-C	6.16	117.99	111.28
1	B	458	ASN	N-CA-C	5.85	119.64	111.74
1	A	544	LEU	N-CA-C	-5.84	106.13	113.72
1	A	371	ASP	N-CA-C	5.76	118.69	110.68
1	A	409	HIS	N-CA-C	5.74	118.31	111.71
1	B	409	HIS	N-CA-C	5.66	119.00	111.75
1	A	337	LYS	N-CA-C	5.59	118.90	111.75
1	B	340	MSE	N-CA-C	5.55	117.77	111.11
1	B	371	ASP	N-CA-C	5.50	118.13	108.56
1	A	356	TYR	N-CA-C	5.47	117.25	111.28
1	B	487	ARG	N-CA-C	-5.45	105.24	111.07
1	A	335	LEU	N-CA-C	5.42	118.11	111.82
1	A	457	LEU	N-CA-C	-5.32	106.63	113.01
1	B	473	SER	N-CA-C	-5.22	103.25	110.35
1	A	272	ILE	N-CA-C	-5.14	105.69	110.53
1	A	429	ALA	N-CA-C	-5.14	105.57	111.07

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	2549	0	2534	107	0
1	B	2550	0	2534	96	0
All	All	5099	0	5068	190	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 (190) 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:340:MSE:HE1	1:A:376:LEU:HD23	1.48	0.96
1:A:262:ASN:HD21	1:A:266:THR:HG22	1.31	0.93
1:B:468:GLU:O	1:B:472:THR:HB	1.73	0.89
1:A:468:GLU:O	1:A:472:THR:HG22	1.74	0.88
1:A:512:PHE:HB3	1:A:515:GLU:HB3	1.56	0.87
1:B:269:GLN:HA	1:B:272:ILE:HD12	1.58	0.85
1:B:266:THR:HG22	1:B:268:ASP:H	1.42	0.85
1:A:340:MSE:CE	1:A:376:LEU:HD23	2.10	0.80
1:B:515:GLU:HG3	1:B:516:TYR:CD1	2.20	0.77
1:A:301:LEU:HD13	1:A:323:GLU:HB3	1.67	0.76
1:B:249:VAL:O	1:B:253:LYS:HG2	1.86	0.74
1:A:248:GLN:HA	1:A:251:MSE:HE2	1.73	0.70
1:A:466:LEU:O	1:A:470:VAL:HG13	1.91	0.70
1:A:245:GLU:HA	1:A:248:GLN:NE2	2.07	0.70
1:B:332:ILE:O	1:B:337:LYS:HA	1.93	0.69
1:A:489:LEU:HD11	1:A:501:ILE:HG23	1.76	0.67
1:A:537:SER:O	1:A:541:LEU:HB2	1.94	0.67
1:A:259:GLU:OE2	1:A:275:ARG:NH1	2.28	0.67
1:B:485:TRP:HB3	1:B:508:ARG:HG2	1.77	0.67
1:A:493:SER:HA	1:A:501:ILE:HD11	1.76	0.66
1:A:329:GLU:HG2	1:A:346:TYR:OH	1.95	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:288:GLY:HA3	1:A:335:LEU:HD13	1.78	0.65
1:A:242:THR:HG23	1:A:245:GLU:HB2	1.79	0.65
1:A:262:ASN:ND2	1:A:266:THR:HG22	2.07	0.65
1:B:248:GLN:HA	1:B:251:MSE:HE2	1.78	0.64
1:A:471:LEU:HD21	1:A:484:ILE:HB	1.79	0.64
1:A:289:HIS:CD2	1:A:289:HIS:H	2.15	0.63
1:A:447:VAL:O	1:A:451:ILE:HG12	1.99	0.62
1:A:454:LEU:HD22	1:A:466:LEU:HD22	1.83	0.61
1:A:340:MSE:HE2	1:B:535:PRO:HA	1.83	0.60
1:A:340:MSE:HE1	1:A:376:LEU:CD2	2.29	0.60
1:B:309:ALA:HB2	1:B:317:ALA:HB1	1.81	0.60
1:B:513:ARG:O	1:B:517:GLU:HB2	2.01	0.60
1:A:270:THR:O	1:A:274:LYS:HG3	2.01	0.59
1:A:382:MSE:HE1	1:A:415:ALA:HA	1.85	0.59
1:A:252:TRP:O	1:A:256:ILE:HG13	2.02	0.58
1:B:250:ASP:OD1	1:B:254:LYS:HE3	2.03	0.58
1:B:309:ALA:O	1:B:314:MSE:HE1	2.03	0.57
1:B:515:GLU:HG3	1:B:516:TYR:HD1	1.68	0.57
1:A:524:LEU:CD2	1:B:341:LEU:HD11	2.35	0.57
1:B:447:VAL:HG11	1:B:470:VAL:CG1	2.34	0.57
1:B:379:ILE:O	1:B:383:LYS:HG2	2.05	0.56
1:B:264:LEU:O	1:B:265:ARG:C	2.48	0.56
1:B:269:GLN:HB3	1:B:308:LEU:HD21	1.86	0.56
1:B:272:ILE:O	1:B:276:VAL:HG23	2.05	0.56
1:B:319:LEU:O	1:B:323:GLU:HG2	2.06	0.56
1:B:468:GLU:CD	1:B:507:ARG:HH22	2.13	0.56
1:B:265:ARG:O	1:B:265:ARG:HG3	2.05	0.56
1:B:350:GLU:OE2	1:B:358:LYS:NZ	2.39	0.55
1:A:340:MSE:HE1	1:A:376:LEU:HB3	1.88	0.55
1:A:245:GLU:O	1:A:249:VAL:HG23	2.07	0.55
1:B:270:THR:O	1:B:274:LYS:HG3	2.07	0.55
1:B:516:TYR:CG	1:B:523:LEU:HD21	2.41	0.55
1:A:528:TYR:HB2	1:A:535:PRO:CG	2.37	0.55
1:A:531:MSE:O	1:A:532:ASP:OD1	2.25	0.55
1:A:531:MSE:HE2	1:B:410:HIS:CD2	2.41	0.54
1:A:529:LYS:O	1:A:529:LYS:HG3	2.06	0.54
1:A:292:ASP:O	1:A:296:GLU:HG3	2.08	0.54
1:B:309:ALA:HB2	1:B:317:ALA:CB	2.37	0.54
1:B:350:GLU:HA	1:B:350:GLU:OE1	2.08	0.54
1:B:485:TRP:HB3	1:B:508:ARG:CG	2.38	0.54
1:B:485:TRP:CB	1:B:508:ARG:HG2	2.37	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:427:SER:O	1:A:431:LYS:HG3	2.07	0.53
1:B:288:GLY:HA3	1:B:335:LEU:CD1	2.38	0.53
1:B:471:LEU:HD21	1:B:484:ILE:HB	1.90	0.53
1:A:515:GLU:HG3	1:A:516:TYR:HD1	1.73	0.53
1:A:275:ARG:HB3	1:A:275:ARG:HH11	1.72	0.53
1:B:306:LYS:O	1:B:310:GLU:HB2	2.09	0.52
1:B:468:GLU:OE1	1:B:507:ARG:NH2	2.43	0.52
1:B:497:ASP:O	1:B:501:ILE:HG13	2.09	0.52
1:A:374:PRO:HB2	1:A:378:TYR:CE1	2.43	0.52
1:B:516:TYR:HB2	1:B:523:LEU:HD11	1.91	0.52
1:B:299:GLN:O	1:B:303:GLN:HG3	2.10	0.52
1:A:521:THR:HG22	1:A:544:LEU:HD13	1.90	0.52
1:B:454:LEU:HD22	1:B:466:LEU:HD22	1.92	0.52
1:A:355:LYS:HB2	1:A:358:LYS:HB3	1.92	0.52
1:A:495:ILE:HG22	1:B:422:CYS:SG	2.50	0.52
1:B:284:LEU:HD22	1:B:294:TRP:CE2	2.45	0.51
1:A:498:LEU:O	1:A:502:LEU:HG	2.10	0.51
1:B:516:TYR:CD1	1:B:523:LEU:HD21	2.45	0.51
1:A:253:LYS:NZ	1:A:296:GLU:OE2	2.43	0.51
1:A:266:THR:OG1	1:A:267:GLU:N	2.43	0.51
1:A:448:LEU:HG	1:A:484:ILE:CD1	2.41	0.50
1:B:451:ILE:HD13	1:B:467:PHE:CE2	2.46	0.50
1:A:524:LEU:HD23	1:B:341:LEU:HD11	1.93	0.50
1:B:288:GLY:HA3	1:B:335:LEU:HD13	1.94	0.50
1:A:515:GLU:HG3	1:A:516:TYR:CD1	2.47	0.50
1:B:360:HIS:HD2	1:B:381:TYR:OH	1.94	0.50
1:B:361:SER:O	1:B:365:ARG:HG3	2.12	0.49
1:A:459:GLU:HG3	1:A:462:ASN:HB2	1.95	0.49
1:A:275:ARG:NH1	1:A:275:ARG:HB3	2.28	0.49
1:A:260:LYS:HE2	1:A:300:TYR:CD1	2.48	0.49
1:A:528:TYR:HB2	1:A:535:PRO:HG2	1.94	0.48
1:B:447:VAL:HG11	1:B:470:VAL:HG11	1.95	0.48
1:A:334:THR:O	1:A:337:LYS:HE3	2.13	0.48
1:A:478:PRO:HD2	1:A:479:GLU:OE1	2.13	0.48
1:B:355:LYS:O	1:B:358:LYS:HB3	2.12	0.48
1:B:374:PRO:HG3	1:B:406:ARG:CZ	2.43	0.48
1:B:512:PHE:O	1:B:513:ARG:C	2.56	0.48
1:B:425:ASP:CG	1:B:428:VAL:HG23	2.38	0.48
1:B:391:ILE:O	1:B:395:ARG:HG3	2.14	0.48
1:B:448:LEU:HD22	1:B:487:ARG:CZ	2.44	0.48
1:B:419:GLU:HA	1:B:419:GLU:OE1	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:530:PHE:O	1:A:531:MSE:C	2.56	0.47
1:B:337:LYS:O	1:B:343:TYR:OH	2.28	0.47
1:B:419:GLU:O	1:B:423:SER:HB2	2.14	0.47
1:A:259:GLU:CD	1:A:275:ARG:NH1	2.72	0.47
1:A:408:ARG:HH21	1:A:410:HIS:CE1	2.32	0.47
1:A:425:ASP:OD2	1:A:425:ASP:C	2.58	0.47
1:B:382:MSE:HE2	1:B:418:MSE:CE	2.45	0.47
1:A:269:GLN:HB3	1:A:308:LEU:HD21	1.97	0.47
1:A:309:ALA:HA	1:A:314:MSE:HE2	1.97	0.47
1:A:454:LEU:HD12	1:A:457:LEU:HD12	1.96	0.47
1:B:447:VAL:HG11	1:B:470:VAL:HG13	1.96	0.47
1:A:372:ILE:C	1:A:372:ILE:HD12	2.40	0.46
1:A:497:ASP:HA	1:B:387:ARG:HD2	1.98	0.46
1:B:266:THR:HG22	1:B:268:ASP:N	2.20	0.46
1:B:304:SER:O	1:B:308:LEU:HG	2.16	0.46
1:A:536:CYS:HB2	1:A:541:LEU:HD13	1.96	0.46
1:B:325:ALA:HB2	1:B:353:ARG:HH12	1.80	0.46
1:A:398:PHE:HE2	1:A:411:VAL:HG13	1.81	0.46
1:A:292:ASP:OD1	1:A:292:ASP:N	2.49	0.46
1:B:514:GLU:O	1:B:515:GLU:C	2.59	0.46
1:A:242:THR:HG23	1:A:245:GLU:CB	2.44	0.45
1:B:382:MSE:HE1	1:B:415:ALA:HA	1.97	0.45
1:A:528:TYR:HB2	1:A:535:PRO:HG3	1.98	0.45
1:B:517:GLU:HA	1:B:517:GLU:OE1	2.16	0.45
1:A:425:ASP:OD2	1:A:427:SER:HB3	2.16	0.45
1:B:256:ILE:HG22	1:B:260:LYS:HE3	1.99	0.45
1:A:524:LEU:HG	1:B:341:LEU:HD11	1.99	0.45
1:A:391:ILE:HG21	1:A:395:ARG:NH2	2.32	0.45
1:A:546:TYR:OH	1:A:548:ASP:HB2	2.17	0.45
1:A:408:ARG:NH2	1:B:533:LEU:HD21	2.32	0.45
1:A:410:HIS:CD2	1:B:531:MSE:HE2	2.51	0.45
1:A:417:LEU:HD22	1:A:421:TYR:HE1	1.81	0.44
1:A:540:GLU:OE2	1:B:339:ASN:ND2	2.50	0.44
1:A:382:MSE:HB3	1:A:418:MSE:HE1	1.97	0.44
1:A:491:PHE:CZ	1:A:495:ILE:HD11	2.53	0.44
1:B:374:PRO:HB2	1:B:378:TYR:CE1	2.53	0.44
1:A:516:TYR:CD2	1:A:523:LEU:HB2	2.53	0.44
1:A:517:GLU:HA	1:A:517:GLU:OE1	2.16	0.44
1:A:526:ASP:HA	1:A:529:LYS:HG2	1.99	0.44
1:B:302:GLU:HG3	1:B:349:TYR:CE1	2.52	0.44
1:A:271:LEU:O	1:A:275:ARG:HG3	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:546:TYR:CZ	1:A:548:ASP:HB2	2.52	0.43
1:B:365:ARG:O	1:B:368:ALA:HB3	2.18	0.43
1:B:512:PHE:HB3	1:B:515:GLU:HB3	2.01	0.43
1:B:531:MSE:O	1:B:531:MSE:HG3	2.19	0.43
1:A:459:GLU:CD	1:A:462:ASN:HD22	2.26	0.43
1:A:537:SER:OG	1:A:540:GLU:HB2	2.18	0.43
1:B:407:THR:CG2	1:B:411:VAL:HG21	2.48	0.43
1:B:531:MSE:O	1:B:532:ASP:OD1	2.37	0.43
1:B:528:TYR:HB2	1:B:535:PRO:CG	2.48	0.43
1:B:256:ILE:HD13	1:B:280:TYR:CD2	2.54	0.43
1:B:262:ASN:C	1:B:262:ASN:HD22	2.27	0.43
1:A:340:MSE:HE3	1:A:344:PHE:CZ	2.53	0.43
1:A:460:ASP:OD2	1:A:460:ASP:N	2.51	0.43
1:A:507:ARG:O	1:A:510:THR:HB	2.19	0.43
1:A:255:TYR:CD2	1:A:255:TYR:C	2.97	0.43
1:A:331:ALA:HB1	1:A:342:LEU:HD13	2.01	0.43
1:B:524:LEU:O	1:B:527:ARG:HB3	2.18	0.42
1:B:539:SER:O	1:B:542:LYS:HB3	2.19	0.42
1:B:485:TRP:CG	1:B:508:ARG:HG2	2.55	0.42
1:B:425:ASP:OD1	1:B:428:VAL:HG23	2.19	0.42
1:A:303:GLN:H	1:A:303:GLN:HG2	1.68	0.42
1:B:273:THR:HG21	1:B:320:PHE:CE2	2.54	0.42
1:A:483:GLU:O	1:A:487:ARG:HG2	2.20	0.41
1:A:479:GLU:H	1:A:479:GLU:CD	2.22	0.41
1:B:447:VAL:CG1	1:B:470:VAL:HG11	2.49	0.41
1:A:277:MSE:O	1:A:281:GLU:HG2	2.20	0.41
1:B:248:GLN:CA	1:B:251:MSE:HE2	2.49	0.41
1:B:375:THR:O	1:B:379:ILE:HG13	2.20	0.41
1:B:528:TYR:HB2	1:B:535:PRO:HG3	2.03	0.41
1:A:264:LEU:O	1:A:265:ARG:HB2	2.20	0.41
1:A:457:LEU:HD23	1:A:457:LEU:HA	1.92	0.41
1:B:312:GLY:O	1:B:314:MSE:N	2.54	0.41
1:A:295:TYR:CZ	1:B:524:LEU:HD22	2.56	0.41
1:A:490:ALA:O	1:A:493:SER:HB3	2.21	0.41
1:A:419:GLU:HG3	1:A:428:VAL:HG12	2.03	0.41
1:A:464:ARG:NE	1:A:492:GLU:OE2	2.49	0.41
1:B:277:MSE:SE	1:B:301:LEU:HG	2.70	0.41
1:B:464:ARG:NE	1:B:492:GLU:OE2	2.51	0.41
1:A:382:MSE:HE1	1:A:415:ALA:CA	2.50	0.41
1:A:309:ALA:HA	1:A:314:MSE:CE	2.51	0.40
1:A:419:GLU:O	1:A:423:SER:HB2	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:340:MSE:HE3	1:A:344:PHE:CE1	2.56	0.40
1:A:348:ASP:OD2	1:B:527:ARG:NH1	2.54	0.40
1:A:381:TYR:CE1	1:A:397:ILE:HD12	2.57	0.40
1:A:323:GLU:O	1:A:327:ILE:HG12	2.22	0.40
1:A:422:CYS:SG	1:B:495:ILE:HG22	2.62	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	306/308 (99%)	291 (95%)	13 (4%)	2 (1%)	18	47
1	B	306/308 (99%)	285 (93%)	16 (5%)	5 (2%)	7	27
All	All	612/616 (99%)	576 (94%)	29 (5%)	7 (1%)	11	36

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	515	GLU
1	A	531	MSE
1	B	265	ARG
1	B	313	ASP
1	B	514	GLU
1	A	515	GLU
1	B	531	MSE

5.3.2 Protein sidechains [i](#)

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	270/261 (103%)	258 (96%)	12 (4%)	25	59
1	B	270/261 (103%)	261 (97%)	9 (3%)	33	69
All	All	540/522 (103%)	519 (96%)	21 (4%)	28	64

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	289	HIS
1	A	302	GLU
1	A	303	GLN
1	A	323	GLU
1	A	355	LYS
1	A	386	ARG
1	A	411	VAL
1	A	445	GLU
1	A	460	ASP
1	A	470	VAL
1	A	507	ARG
1	A	532	ASP
1	B	391	ILE
1	B	403	GLU
1	B	407	THR
1	B	458	ASN
1	B	472	THR
1	B	479	GLU
1	B	508	ARG
1	B	529	LYS
1	B	532	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	248	GLN
1	A	257	GLN
1	A	262	ASN
1	A	289	HIS
1	A	316	ASN
1	A	380	GLN

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Mol	Chain	Res	Type
1	A	458	ASN
1	A	462	ASN
1	B	247	GLN
1	B	257	GLN
1	B	303	GLN
1	B	316	ASN
1	B	326	ASN
1	B	360	HIS
1	B	380	GLN
1	B	458	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](#)

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	299/308 (97%)	-0.05	7 (2%) 61 51	28, 48, 94, 106	0
1	B	299/308 (97%)	0.04	7 (2%) 61 51	29, 54, 99, 109	0
All	All	598/616 (97%)	-0.00	14 (2%) 61 51	28, 51, 97, 109	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	549	VAL	5.1
1	A	516	TYR	4.0
1	B	516	TYR	3.8
1	B	242	THR	2.9
1	B	549	VAL	2.7
1	B	461	ASN	2.6
1	A	243	PRO	2.5
1	A	548	ASP	2.3
1	B	512	PHE	2.2
1	A	511	ALA	2.2
1	B	266	THR	2.2
1	B	475	SER	2.1
1	A	513	ARG	2.1
1	A	510	THR	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 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.