



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 24, 2026 – 02:20 AM UTC

PDB ID : 4EL1 / pdb_00004el1
Title : Crystal structure of oxidized hPDI (abb'xa')
Authors : Wang, C.; Li, W.; Ren, J.; Ke, H.; Gong, W.; Feng, W.; Wang, C.-C.
Deposited on : 2012-04-10
Resolution : 2.88 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

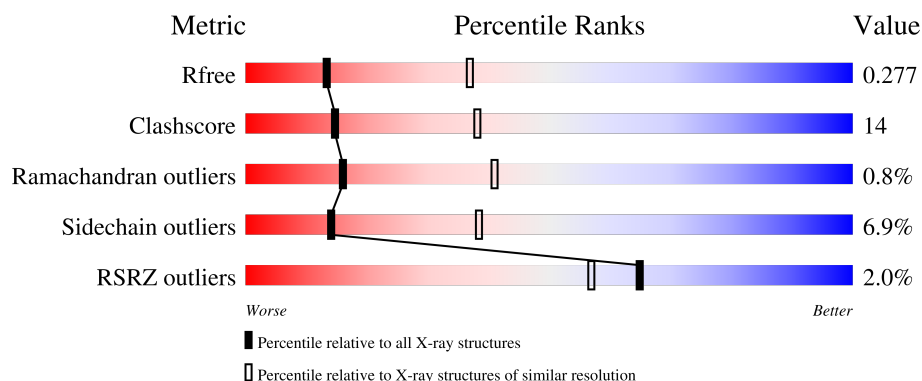
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

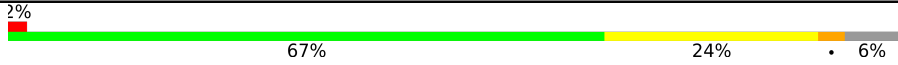
The reported resolution of this entry is 2.88 Å.

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	3557 (2.90-2.86)
Clashscore	190562	3801 (2.90-2.86)
Ramachandran outliers	187476	3699 (2.90-2.86)
Sidechain outliers	187428	3702 (2.90-2.86)
RSRZ outliers	180081	3558 (2.90-2.86)

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	482	
1	B	482	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7222 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 Protein disulfide-isomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	452	Total	C	N	O	S	19	2	0
			3601	2310	589	691	11			
1	B	452	Total	C	N	O	S	20	2	0
			3601	2310	589	691	11			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP P07237
A	-18	HIS	-	expression tag	UNP P07237
A	-17	HIS	-	expression tag	UNP P07237
A	-16	HIS	-	expression tag	UNP P07237
A	-15	HIS	-	expression tag	UNP P07237
A	-14	HIS	-	expression tag	UNP P07237
A	-13	HIS	-	expression tag	UNP P07237
A	-12	SER	-	expression tag	UNP P07237
A	-11	SER	-	expression tag	UNP P07237
A	-10	GLY	-	expression tag	UNP P07237
A	-9	LEU	-	expression tag	UNP P07237
A	-8	GLU	-	expression tag	UNP P07237
A	-7	VAL	-	expression tag	UNP P07237
A	-6	LEU	-	expression tag	UNP P07237
A	-5	PHE	-	expression tag	UNP P07237
A	-4	GLN	-	expression tag	UNP P07237
A	-3	GLY	-	expression tag	UNP P07237
A	-2	PRO	-	expression tag	UNP P07237
A	-1	GLY	-	expression tag	UNP P07237
A	0	SER	-	expression tag	UNP P07237
B	-19	MET	-	expression tag	UNP P07237
B	-18	HIS	-	expression tag	UNP P07237
B	-17	HIS	-	expression tag	UNP P07237
B	-16	HIS	-	expression tag	UNP P07237
B	-15	HIS	-	expression tag	UNP P07237

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	expression tag	UNP P07237
B	-13	HIS	-	expression tag	UNP P07237
B	-12	SER	-	expression tag	UNP P07237
B	-11	SER	-	expression tag	UNP P07237
B	-10	GLY	-	expression tag	UNP P07237
B	-9	LEU	-	expression tag	UNP P07237
B	-8	GLU	-	expression tag	UNP P07237
B	-7	VAL	-	expression tag	UNP P07237
B	-6	LEU	-	expression tag	UNP P07237
B	-5	PHE	-	expression tag	UNP P07237
B	-4	GLN	-	expression tag	UNP P07237
B	-3	GLY	-	expression tag	UNP P07237
B	-2	PRO	-	expression tag	UNP P07237
B	-1	GLY	-	expression tag	UNP P07237
B	0	SER	-	expression tag	UNP P07237

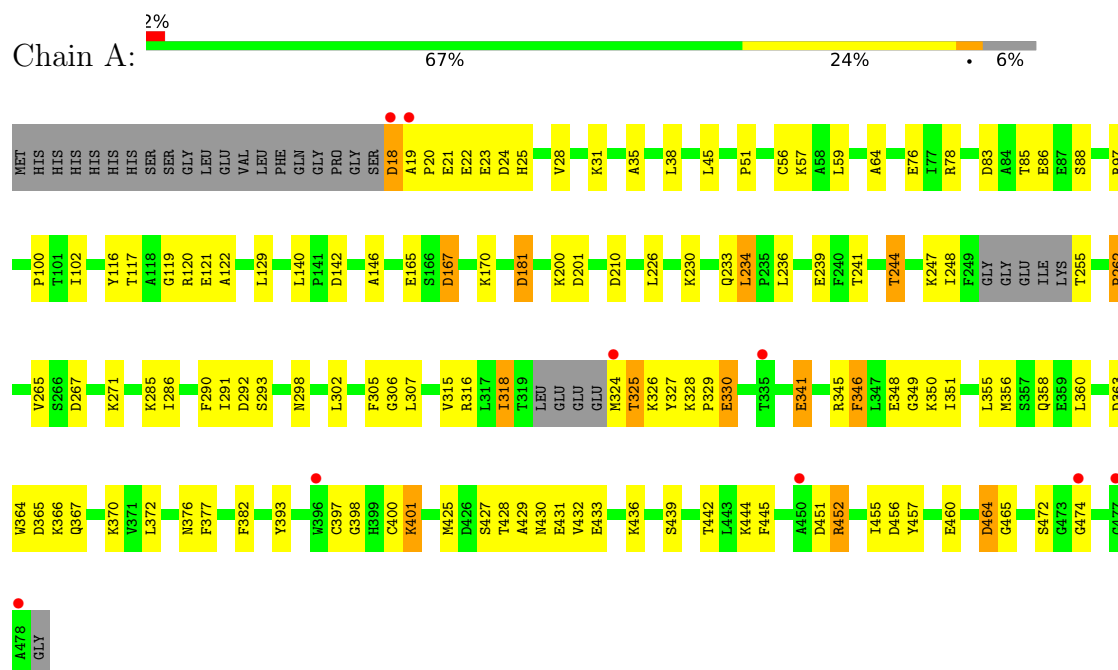
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	10	Total O 10 10	0	0
2	B	10	Total O 10 10	0	0

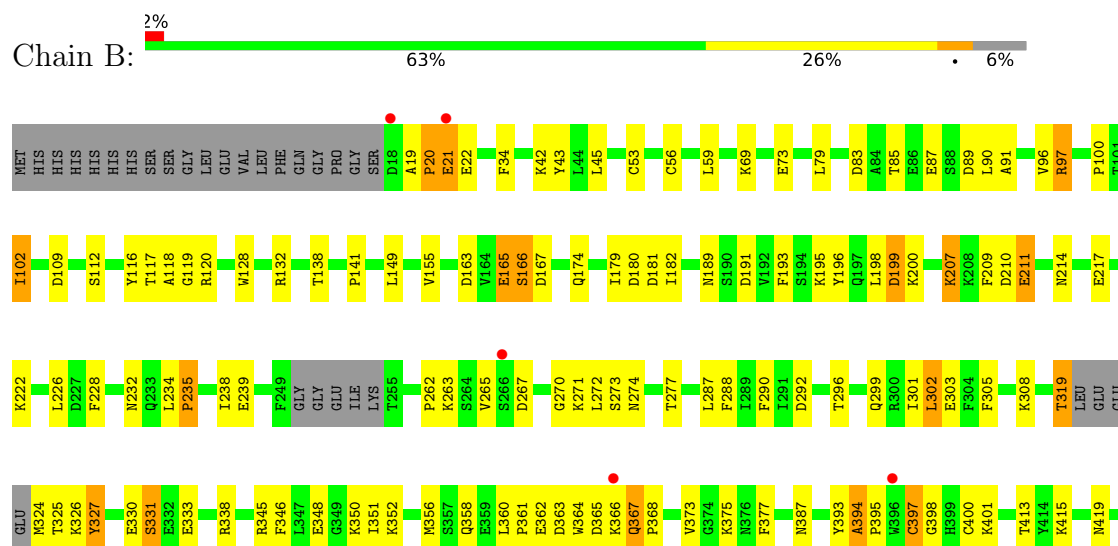
3 Residue-property plots [i](#)

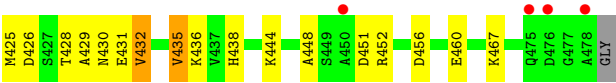
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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



• Molecule 1: Protein disulfide-isomerase





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	160.75Å 34.01Å 156.96Å 90.00° 104.89° 90.00°	Depositor
Resolution (Å)	29.84 – 2.88 29.84 – 2.88	Depositor EDS
% Data completeness (in resolution range)	98.9 (29.84-2.88) 94.5 (29.84-2.88)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.70 (at 2.90Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.261 , 0.281 0.258 , 0.277	Depositor DCC
R_{free} test set	981 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	44.2	Xtriage
Anisotropy	0.216	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 41.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7222	wwPDB-VP
Average B, all atoms (Å ²)	51.0	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 90.49 % 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 2.2531e-08. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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

5 Model quality [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	0.89	2/3690 (0.1%)	0.85	4/4982 (0.1%)
1	B	0.81	1/3690 (0.0%)	0.83	5/4982 (0.1%)
All	All	0.85	3/7380 (0.0%)	0.84	9/9964 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	318	ILE	CA-CB	-8.69	1.43	1.54
1	B	235	PRO	N-CD	6.05	1.56	1.47
1	A	262	PRO	N-CD	5.99	1.56	1.47

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	401	LYS	N-CA-C	-7.62	103.06	114.64
1	B	394	ALA	CA-C-N	6.33	126.02	119.56
1	B	394	ALA	C-N-CA	6.33	126.02	119.56
1	A	318	ILE	CB-CA-C	-6.23	102.62	111.34
1	B	210	ASP	CB-CA-C	-6.23	109.40	116.63
1	A	393	TYR	N-CA-C	6.09	117.45	108.86
1	B	238	ILE	N-CA-C	5.75	116.15	108.11
1	B	393	TYR	N-CA-C	5.22	116.59	109.18
1	A	210	ASP	CB-CA-C	-5.03	110.80	116.63

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	3601	0	3522	96	2
1	B	3601	0	3520	98	2
2	A	10	0	0	0	0
2	B	10	0	0	0	0
All	All	7222	0	7042	194	2

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 14.

All (194) 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:244:THR:HG22	1:A:247:LYS:HE3	1.27	1.14
1:B:397:CYS:SG	1:B:400:CYS:CB	2.36	1.14
1:B:397:CYS:SG	1:B:400:CYS:N	2.30	1.04
1:A:244:THR:HG22	1:A:247:LYS:CE	1.87	1.04
1:B:163:ASP:OD1	1:B:165:GLU:HG2	1.61	1.00
1:B:19:ALA:HB1	1:B:20:PRO:HD2	1.48	0.93
1:B:397:CYS:SG	1:B:400:CYS:HB2	2.06	0.92
1:A:117:THR:HG22	1:A:117:THR:O	1.74	0.86
1:B:413:THR:HG21	1:B:467:LYS:CE	2.07	0.85
1:B:179:ILE:HD11	1:B:226:LEU:CD2	2.09	0.82
1:B:163:ASP:OD1	1:B:165:GLU:CG	2.26	0.82
1:A:360:LEU:HD13	1:A:376:ASN:ND2	1.95	0.81
1:B:179:ILE:HD11	1:B:226:LEU:HD22	1.61	0.80
1:A:306:GLY:HA3	1:A:356:MET:CE	2.12	0.80
1:A:19:ALA:N	1:A:20:PRO:CD	2.47	0.77
1:B:42:LYS:HE2	1:B:43:TYR:OH	1.86	0.76
1:A:306:GLY:HA3	1:A:356:MET:HE1	1.68	0.76
1:B:53[A]:CYS:CB	1:B:56[A]:CYS:SG	2.73	0.76
1:A:200:LYS:HG2	1:A:201:ASP:H	1.51	0.75
1:B:102:ILE:HB	1:B:116:TYR:HB3	1.69	0.75
1:A:181:ASP:OD1	1:A:181:ASP:N	2.14	0.74
1:A:326:LYS:O	1:A:327:TYR:HD1	1.69	0.74
1:B:413:THR:HG21	1:B:467:LYS:HE2	1.71	0.73
1:B:141:PRO:O	1:B:189:ASN:ND2	2.21	0.73
1:A:19:ALA:N	1:A:20:PRO:HD3	2.05	0.72
1:A:377:PHE:CD2	1:A:432:VAL:HG21	2.24	0.72
1:A:200:LYS:HG2	1:A:201:ASP:N	2.05	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:179:ILE:CD1	1:B:226:LEU:CD2	2.68	0.70
1:A:325:THR:HG22	1:A:327:TYR:HE1	1.57	0.69
1:A:360:LEU:CD1	1:A:376:ASN:HD22	2.05	0.69
1:B:270:GLY:O	1:B:274:ASN:ND2	2.26	0.69
1:A:358:GLN:OE1	1:A:430:ASN:ND2	2.26	0.69
1:B:363:ASP:O	1:B:365:ASP:N	2.26	0.68
1:B:209:PHE:CD1	1:B:235:PRO:HG3	2.28	0.68
1:A:428:THR:OG1	1:A:429:ALA:N	2.24	0.68
1:B:397:CYS:SG	1:B:400:CYS:CA	2.82	0.68
1:A:397:CYS:CB	1:A:400:CYS:SG	2.80	0.68
1:B:53[A]:CYS:HB3	1:B:56[A]:CYS:SG	2.34	0.68
1:A:56[B]:CYS:SG	1:A:100:PRO:HG3	2.34	0.67
1:A:244:THR:CG2	1:A:247:LYS:HE3	2.17	0.67
1:A:315:VAL:O	1:A:316:ARG:NH1	2.26	0.67
1:A:398:GLY:HA2	1:A:401:LYS:HE2	1.75	0.67
1:B:362:GLU:O	1:B:366:LYS:NZ	2.28	0.66
1:A:363:ASP:HA	1:A:366:LYS:HD3	1.77	0.66
1:A:142:ASP:N	1:A:142:ASP:OD1	2.27	0.66
1:B:91:ALA:HB1	1:B:96:VAL:HG21	1.77	0.66
1:A:360:LEU:CD1	1:A:376:ASN:ND2	2.58	0.66
1:B:19:ALA:HB1	1:B:20:PRO:CD	2.23	0.66
1:A:365:ASP:HB2	1:A:370:LYS:HE2	1.77	0.66
1:A:326:LYS:C	1:A:327:TYR:HD1	2.04	0.65
1:B:228:PHE:O	1:B:232:ASN:ND2	2.30	0.65
1:A:397:CYS:HB3	1:A:400:CYS:SG	2.36	0.65
1:A:21:GLU:HG3	1:A:28:VAL:HB	1.78	0.65
1:A:349:GLY:O	1:A:351:ILE:N	2.30	0.64
1:B:181:ASP:O	1:B:182:ILE:HD13	1.97	0.64
1:B:267:ASP:HB3	1:B:271:LYS:HZ3	1.62	0.64
1:B:356:MET:HE2	1:B:431:GLU:HG3	1.80	0.64
1:B:367:GLN:HG3	1:B:368:PRO:HD2	1.79	0.64
1:B:363:ASP:HA	1:B:366:LYS:NZ	2.13	0.63
1:A:117:THR:O	1:A:117:THR:CG2	2.47	0.63
1:A:360:LEU:HD13	1:A:376:ASN:HD22	1.58	0.62
1:A:83:ASP:OD2	1:A:86:GLU:HG2	1.99	0.62
1:A:377:PHE:CE2	1:A:432:VAL:HG21	2.35	0.62
1:A:327:TYR:CE2	1:A:346:PHE:HD1	2.17	0.62
1:B:69:LYS:O	1:B:73:GLU:HG3	2.00	0.61
1:B:207:LYS:HE3	1:B:211:GLU:HG3	1.81	0.61
1:B:305:PHE:HA	1:B:326:LYS:HZ3	1.65	0.60
1:A:244:THR:HG22	1:A:247:LYS:HE2	1.76	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:207:LYS:HD2	1:B:209:PHE:O	2.00	0.60
1:B:56[B]:CYS:SG	1:B:100:PRO:HB3	2.42	0.60
1:B:263:LYS:HG3	1:B:292:ASP:OD1	2.02	0.60
1:B:87:GLU:O	1:B:89:ASP:N	2.31	0.59
1:A:21:GLU:CG	1:A:28:VAL:HB	2.34	0.58
1:B:363:ASP:HA	1:B:366:LYS:HZ3	1.68	0.58
1:B:174:GLN:HB3	1:B:222:LYS:HD2	1.85	0.58
1:A:239:GLU:HG2	1:A:290:PHE:CZ	2.39	0.57
1:B:345:ARG:HB2	1:B:351:ILE:HD11	1.85	0.57
1:A:444:LYS:NZ	1:A:456:ASP:OD1	2.33	0.57
1:A:326:LYS:C	1:A:327:TYR:CD1	2.83	0.56
1:A:306:GLY:HA3	1:A:356:MET:HE3	1.85	0.56
1:A:398:GLY:HA2	1:A:401:LYS:HG2	1.87	0.56
1:A:59:LEU:HD13	1:A:120:ARG:HB3	1.88	0.56
1:B:426:ASP:O	1:B:430:ASN:HB2	2.05	0.56
1:A:327:TYR:HE2	1:A:346:PHE:HD1	1.53	0.55
1:A:349:GLY:C	1:A:351:ILE:H	2.14	0.55
1:B:42:LYS:HE2	1:B:43:TYR:CZ	2.40	0.55
1:B:45:LEU:HB3	1:B:79:LEU:HD22	1.88	0.55
1:A:18:ASP:C	1:A:20:PRO:HD2	2.32	0.55
1:A:200:LYS:CG	1:A:201:ASP:H	2.19	0.54
1:B:209:PHE:CE1	1:B:235:PRO:HG3	2.42	0.54
1:A:18:ASP:C	1:A:20:PRO:CD	2.80	0.54
1:B:377:PHE:HB2	1:B:425:MET:HE1	1.89	0.54
1:A:292:ASP:O	1:A:298:ASN:ND2	2.31	0.54
1:A:85:THR:OG1	1:A:86:GLU:OE2	2.22	0.54
1:A:341:GLU:HG3	1:A:345:ARG:HD2	1.88	0.54
1:B:207:LYS:HE2	1:B:214:ASN:ND2	2.23	0.54
1:B:53[B]:CYS:HB3	1:B:56[B]:CYS:HB2	1.90	0.53
1:A:22:GLU:OE1	1:A:78:ARG:NH2	2.40	0.53
1:B:97:ARG:HH11	1:B:97:ARG:HG2	1.73	0.53
1:B:345:ARG:C	1:B:351:ILE:HD12	2.34	0.53
1:B:305:PHE:HA	1:B:326:LYS:NZ	2.23	0.52
1:A:363:ASP:O	1:A:365:ASP:N	2.42	0.52
1:B:199:ASP:OD1	1:B:199:ASP:N	2.24	0.52
1:A:355:LEU:HD12	1:A:431:GLU:HG3	1.92	0.52
1:B:116:TYR:OH	1:B:119:GLY:O	2.25	0.52
1:B:330:GLU:N	1:B:330:GLU:OE1	2.41	0.52
1:B:387:ASN:OD1	1:B:448:ALA:N	2.41	0.52
1:A:445:PHE:HB3	1:A:455:ILE:HG23	1.91	0.51
1:A:24:ASP:HB3	1:A:64:ALA:HB1	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:398:GLY:HA2	1:B:401:LYS:HE2	1.92	0.51
1:A:200:LYS:CG	1:A:201:ASP:N	2.73	0.51
1:B:333:GLU:HB2	1:B:338:ARG:HD3	1.92	0.51
1:A:102:ILE:HB	1:A:116:TYR:HB3	1.93	0.51
1:B:181:ASP:OD1	1:B:181:ASP:N	2.38	0.51
1:A:233:GLN:O	1:A:234:LEU:HB2	2.11	0.50
1:A:226:LEU:O	1:A:230:LYS:HG2	2.11	0.50
1:B:179:ILE:HG23	1:B:181:ASP:OD1	2.11	0.50
1:B:319:THR:OG1	1:B:325:THR:HG23	2.12	0.50
1:B:432:VAL:O	1:B:436:LYS:HE3	2.12	0.50
1:A:376:ASN:OD1	1:A:377:PHE:N	2.45	0.49
1:B:59:LEU:HD13	1:B:120:ARG:HB3	1.93	0.49
1:A:464:ASP:OD1	1:A:464:ASP:C	2.55	0.49
1:B:387:ASN:N	1:B:419:ASN:O	2.32	0.49
1:A:285:LYS:HB3	1:A:286:ILE:HD12	1.94	0.49
1:A:439:SER:O	1:A:442:THR:OG1	2.31	0.48
1:B:426:ASP:OD2	1:B:428:THR:OG1	2.26	0.48
1:A:262:PRO:HD3	1:A:293:SER:OG	2.13	0.48
1:A:382:PHE:CD1	1:A:452:ARG:HG2	2.48	0.48
1:B:128:TRP:O	1:B:132:ARG:NH1	2.45	0.48
1:A:165:GLU:HA	1:A:170:LYS:HE2	1.95	0.48
1:B:345:ARG:HB2	1:B:351:ILE:CD1	2.44	0.47
1:B:21:GLU:O	1:B:22:GLU:C	2.57	0.47
1:B:163:ASP:O	1:B:166:SER:OG	2.31	0.46
1:A:267:ASP:O	1:A:271:LYS:HG3	2.15	0.46
1:B:262:PRO:HG2	1:B:265:VAL:HB	1.97	0.46
1:A:460:GLU:HG3	1:A:465:GLY:HA3	1.97	0.46
1:A:433:GLU:O	1:A:436:LYS:NZ	2.48	0.46
1:B:109:ASP:OD1	1:B:112:SER:N	2.48	0.46
1:B:432:VAL:HG22	1:B:435:VAL:H	1.80	0.46
1:A:45:LEU:HD22	1:A:129:LEU:HD21	1.98	0.46
1:A:328:LYS:HG3	1:A:329:PRO:HD2	1.97	0.46
1:A:116:TYR:OH	1:A:119:GLY:O	2.23	0.46
1:B:348:GLU:HB2	1:B:350:LYS:HD3	1.96	0.46
1:B:327:TYR:CD2	1:B:327:TYR:N	2.84	0.45
1:A:233:GLN:HG3	1:A:234:LEU:HD13	1.97	0.45
1:B:179:ILE:CD1	1:B:226:LEU:HD22	2.36	0.45
1:A:372:LEU:HB2	1:A:425:MET:HB2	1.98	0.45
1:B:356:MET:HG2	1:B:429:ALA:O	2.17	0.45
1:A:244:THR:HB	1:A:248:ILE:HD11	1.98	0.45
1:B:34:PHE:CD1	1:B:90:LEU:HD21	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:287:LEU:HD12	1:B:288:PHE:N	2.32	0.45
1:A:327:TYR:CE2	1:A:346:PHE:CD1	3.03	0.44
1:B:195:LYS:HD3	1:B:196:TYR:CE2	2.53	0.44
1:B:299:GLN:HA	1:B:302:LEU:HB2	2.00	0.44
1:B:373:VAL:HG23	1:B:375:LYS:H	1.82	0.44
1:B:352:LYS:HB3	1:B:352:LYS:HE2	1.66	0.44
1:A:51:PRO:HD3	1:A:83:ASP:OD1	2.18	0.44
1:A:255:THR:CG2	1:A:286:ILE:HG13	2.47	0.44
1:B:356:MET:HE1	1:B:428:THR:O	2.18	0.44
1:A:325:THR:HG22	1:A:327:TYR:CE1	2.45	0.44
1:B:325:THR:HG23	1:B:327:TYR:HE2	1.83	0.43
1:B:189:ASN:O	1:B:193:PHE:HD2	2.01	0.43
1:B:360:LEU:HD13	1:B:373:VAL:HG22	2.00	0.43
1:B:83:ASP:OD1	1:B:85:THR:OG1	2.34	0.43
1:B:179:ILE:HG22	1:B:180:ASP:N	2.34	0.43
1:A:360:LEU:HD11	1:A:376:ASN:HD22	1.80	0.43
1:B:331:SER:HB3	1:B:338:ARG:HH21	1.83	0.43
1:A:140:LEU:HD13	1:A:146:ALA:HA	2.01	0.43
1:A:24:ASP:O	1:A:25:HIS:HB2	2.19	0.43
1:A:457:TYR:OH	1:A:460:GLU:O	2.37	0.42
1:B:273:SER:O	1:B:277:THR:HG23	2.20	0.42
1:B:358:GLN:OE1	1:B:430:ASN:ND2	2.51	0.42
1:B:444:LYS:NZ	1:B:456:ASP:OD1	2.52	0.42
1:B:239:GLU:HG2	1:B:290:PHE:CZ	2.55	0.42
1:B:368:PRO:HB3	1:B:415:LYS:HB3	2.01	0.42
1:B:394:ALA:HA	1:B:395:PRO:HD3	1.75	0.42
1:A:167:ASP:OD1	1:A:167:ASP:N	2.52	0.42
1:A:451:ASP:OD1	1:A:451:ASP:N	2.42	0.42
1:A:31:LYS:HE3	1:A:31:LYS:HB2	1.83	0.42
1:A:236:LEU:HD12	1:A:236:LEU:HA	1.93	0.42
1:A:121:GLU:HG2	1:A:122:ALA:N	2.34	0.42
1:A:472:SER:C	1:A:474:GLY:H	2.28	0.42
1:B:360:LEU:HA	1:B:361:PRO:HD3	1.85	0.42
1:A:318:ILE:HA	1:A:326:LYS:HA	2.02	0.41
1:B:198:LEU:HD22	1:B:200:LYS:O	2.20	0.41
1:B:308:LYS:HE2	1:B:308:LYS:HB3	1.92	0.41
1:A:35:ALA:HA	1:A:38:LEU:HD12	2.03	0.41
1:B:327:TYR:N	1:B:327:TYR:HD2	2.18	0.41
1:B:20:PRO:HB2	1:B:21:GLU:H	1.63	0.41
1:A:305:PHE:O	1:A:356:MET:HE1	2.21	0.41
1:A:330:GLU:OE2	1:A:345:ARG:NH1	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:363:ASP:O	1:B:366:LYS:N	2.46	0.40
1:A:302:LEU:HD23	1:A:307:LEU:O	2.21	0.40

All (2) 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:18:ASP:O	1:B:438:HIS:NE2[1_545]	1.49	0.71
1:A:18:ASP:O	1:B:438:HIS:CE1[1_545]	1.90	0.30

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	448/482 (93%)	420 (94%)	24 (5%)	4 (1%)	14	38
1	B	448/482 (93%)	428 (96%)	17 (4%)	3 (1%)	18	44
All	All	896/964 (93%)	848 (95%)	41 (5%)	7 (1%)	16	41

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	348	GLU
1	A	350	LYS
1	B	364	TRP
1	A	364	TRP
1	B	20	PRO
1	A	234	LEU
1	B	118	ALA

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	386/408 (95%)	366 (95%)	20 (5%)	21	50
1	B	386/408 (95%)	353 (92%)	33 (8%)	10	28
All	All	772/816 (95%)	719 (93%)	53 (7%)	14	38

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	ASP
1	A	23	GLU
1	A	76	GLU
1	A	88	SER
1	A	97	ARG
1	A	167	ASP
1	A	181	ASP
1	A	241	THR
1	A	244	THR
1	A	265	VAL
1	A	291	ILE
1	A	324	MET
1	A	325	THR
1	A	330	GLU
1	A	341	GLU
1	A	346	PHE
1	A	367	GLN
1	A	427	SER
1	A	452	ARG
1	A	464	ASP
1	B	21	GLU
1	B	97	ARG
1	B	102	ILE
1	B	117	THR
1	B	138	THR
1	B	149	LEU
1	B	155	VAL

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Mol	Chain	Res	Type
1	B	165	GLU
1	B	166	SER
1	B	167	ASP
1	B	191	ASP
1	B	199	ASP
1	B	207	LYS
1	B	211	GLU
1	B	217	GLU
1	B	234	LEU
1	B	272	LEU
1	B	296	THR
1	B	301	ILE
1	B	302	LEU
1	B	303	GLU
1	B	319	THR
1	B	324	MET
1	B	327	TYR
1	B	331	SER
1	B	346	PHE
1	B	367	GLN
1	B	397	CYS
1	B	432	VAL
1	B	435	VAL
1	B	451	ASP
1	B	452	ARG
1	B	460	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	55	HIS
1	A	214	ASN
1	A	232	ASN
1	A	274	ASN
1	A	387	ASN
1	A	419	ASN
1	B	214	ASN
1	B	231	HIS
1	B	354	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 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	452/482 (93%)	0.34	9 (1%) 65 57	24, 51, 79, 118	12 (2%)
1	B	452/482 (93%)	0.28	9 (1%) 65 57	24, 49, 73, 105	12 (2%)
All	All	904/964 (93%)	0.31	18 (1%) 65 57	24, 50, 76, 118	24 (2%)

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	21	GLU	3.2
1	B	478	ALA	3.2
1	A	478	ALA	3.0
1	B	366	LYS	2.9
1	B	475	GLN	2.9
1	B	266	SER	2.8
1	A	396	TRP	2.7
1	B	396	TRP	2.7
1	A	18	ASP	2.6
1	A	19	ALA	2.5
1	A	477	GLY	2.5
1	B	476	ASP	2.2
1	A	335	THR	2.2
1	A	450	ALA	2.2
1	A	474	GLY	2.2
1	B	18	ASP	2.2
1	A	324	MET	2.2
1	B	450	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 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.