



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 11:28 AM UTC

PDB ID : 9B48 / pdb_00009b48
Title : Asp324Glu variant of phosphate transporter PiPT from Piriformospora indica
Authors : Gupta, M.; Finer-Moore, J.; Stroud, R.M.
Deposited on : 2024-03-20
Resolution : 4.00 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
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

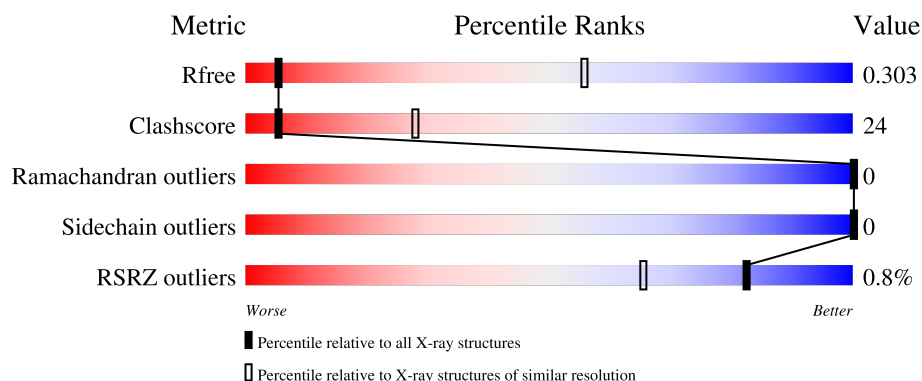
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.00 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	1082 (4.20-3.80)
Clashscore	190562	1129 (4.20-3.80)
Ramachandran outliers	187476	1064 (4.20-3.80)
Sidechain outliers	187428	1055 (4.20-3.80)
RSRZ outliers	180081	1082 (4.20-3.80)

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	530	51% 30% 19%
1	B	530	48% 33% 19%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 13266 atoms, of which 6648 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 Phosphate transporter.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	B	429	Total	C	H	N	O	S	0	0	0
			6633	2186	3324	543	564	16			
1	A	429	Total	C	H	N	O	S	0	0	0
			6633	2186	3324	543	564	16			

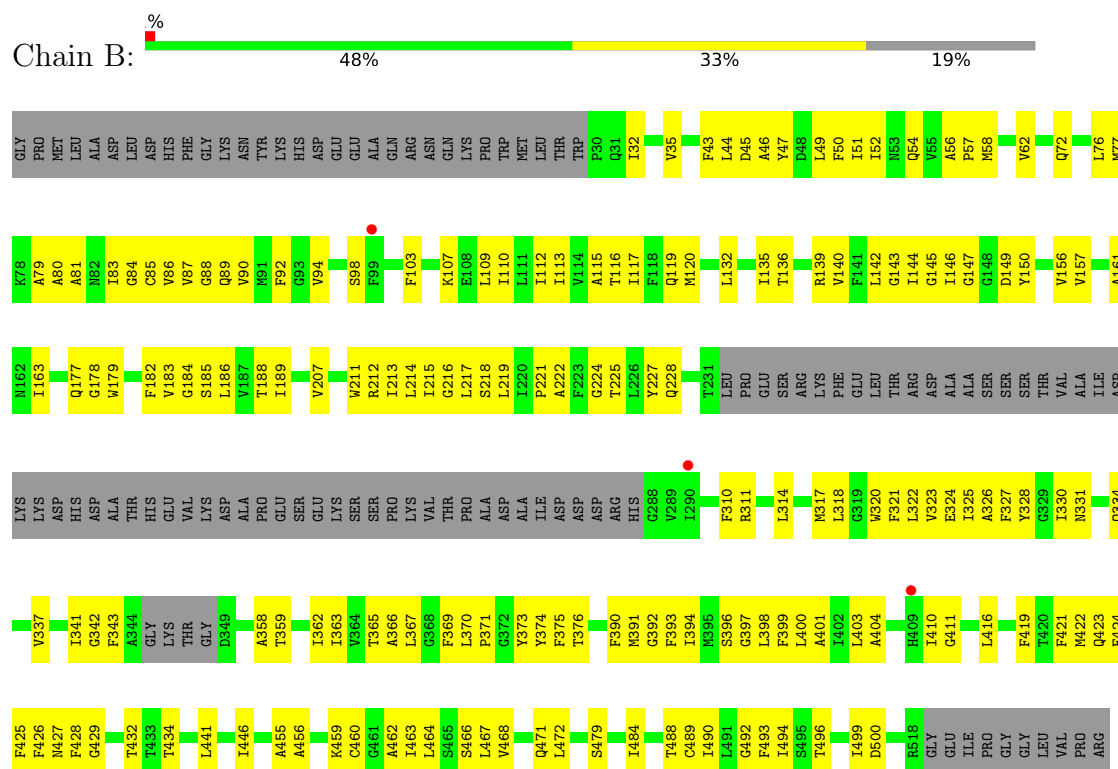
There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-1	GLY	-	expression tag	UNP A8N031
B	0	PRO	-	expression tag	UNP A8N031
B	324	GLU	ASP	engineered mutation	UNP A8N031
B	523	GLY	-	expression tag	UNP A8N031
B	524	GLY	-	expression tag	UNP A8N031
B	525	LEU	-	expression tag	UNP A8N031
B	526	VAL	-	expression tag	UNP A8N031
B	527	PRO	-	expression tag	UNP A8N031
B	528	ARG	-	expression tag	UNP A8N031
A	-1	GLY	-	expression tag	UNP A8N031
A	0	PRO	-	expression tag	UNP A8N031
A	324	GLU	ASP	engineered mutation	UNP A8N031
A	523	GLY	-	expression tag	UNP A8N031
A	524	GLY	-	expression tag	UNP A8N031
A	525	LEU	-	expression tag	UNP A8N031
A	526	VAL	-	expression tag	UNP A8N031
A	527	PRO	-	expression tag	UNP A8N031
A	528	ARG	-	expression tag	UNP A8N031

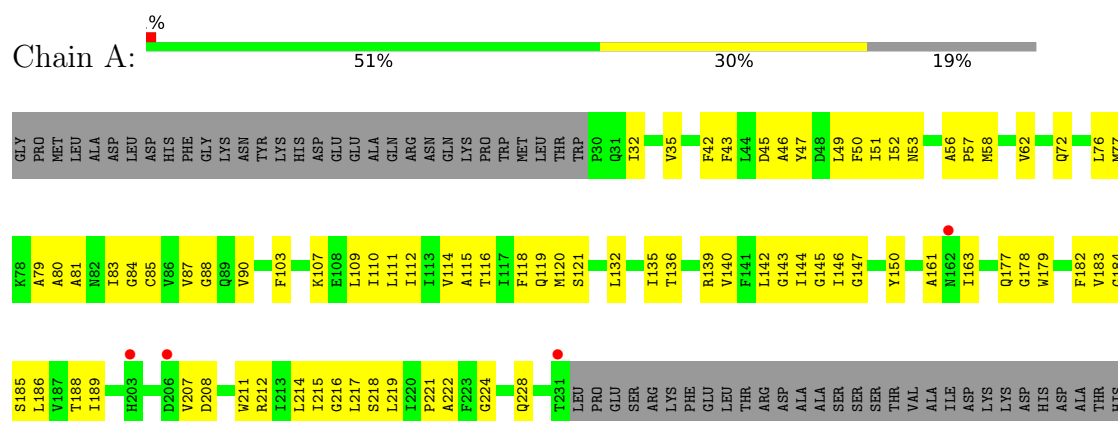
3 Residue-property plots

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: Phosphate transporter



• Molecule 1: Phosphate transporter



K459	C460	G461	A462	I463	L464	S465	S466	L467	V468		Q471	L472	S479		I484	T488	C489		G492	F493	T496	F497	L498	I499	D500	R518	GLY	GLU	ILE	PRO	GLY	GLY	LEU	VAL	PRO	ARG	A358	T359	I362	I363	Y364	T365	A366	L367	G368	F369	L370	P371	G372	Y373	F374	F375	T376	F390	M391	G392	F393	I394	R395	S396	L400	L403	A404	I407	I410	G411	L416	F419	T420	F421	M422	Q423	F424	F425	F426	N427	F428	T434	F435	I436	E440	L441	A455																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																															
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4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	173.39Å 173.39Å 169.12Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	86.69 – 4.00 86.69 – 4.00	Depositor EDS
% Data completeness (in resolution range)	100.0 (86.69-4.00) 86.1 (86.69-4.00)	Depositor EDS
R_{merge}	0.53	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.61 (at 4.01Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.261 , 0.266 0.289 , 0.303	Depositor DCC
R_{free} test set	1595 reflections (9.95%)	wwPDB-VP
Wilson B-factor (Å ²)	115.0	Xtriage
Anisotropy	0.280	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 176.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	0.277 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	13266	wwPDB-VP
Average B, all atoms (Å ²)	187.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.11% 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	0.22	0/3391	0.44	0/4595
1	B	0.22	0/3391	0.43	0/4595
All	All	0.22	0/6782	0.43	0/9190

There are no bond length outliers.

There are no bond angle outliers.

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	3309	3324	3324	154	0
1	B	3309	3324	3324	165	0
All	All	6618	6648	6648	317	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 24.

All (317) 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:85:CYS:SG	1:A:459:LYS:NZ	2.52	0.81
1:B:392:GLY:O	1:B:396:SER:OG	1.98	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:143:GLY:O	1:B:147:GLY:N	2.15	0.79
1:A:376:THR:OG1	1:A:427:ASN:O	2.03	0.77
1:B:119:GLN:HE22	1:B:142:LEU:HD22	1.51	0.75
1:A:365:THR:HA	1:A:369:PHE:HB3	1.68	0.75
1:B:365:THR:HA	1:B:369:PHE:HB3	1.69	0.75
1:A:143:GLY:O	1:A:147:GLY:N	2.20	0.74
1:B:139:ARG:HA	1:B:142:LEU:HB3	1.71	0.73
1:A:219:LEU:HA	1:A:222:ALA:HB3	1.69	0.73
1:A:139:ARG:HA	1:A:142:LEU:HB3	1.69	0.71
1:B:462:ALA:O	1:B:466:SER:OG	2.05	0.71
1:B:219:LEU:HA	1:B:222:ALA:HB3	1.71	0.71
1:A:45:ASP:O	1:A:49:LEU:N	2.20	0.71
1:B:72:GLN:O	1:B:76:LEU:HD13	1.91	0.70
1:A:140:VAL:O	1:A:144:ILE:N	2.21	0.70
1:A:334:GLN:HE22	1:A:365:THR:HG23	1.57	0.70
1:B:320:TRP:HD1	1:B:324:GLU:HG2	1.56	0.70
1:A:330:ILE:O	1:A:334:GLN:N	2.26	0.69
1:A:462:ALA:O	1:A:466:SER:OG	2.11	0.69
1:A:119:GLN:HE22	1:A:142:LEU:HD22	1.57	0.68
1:B:334:GLN:HE22	1:B:365:THR:HG23	1.58	0.68
1:B:464:LEU:O	1:B:468:VAL:HG12	1.92	0.68
1:B:376:THR:OG1	1:B:427:ASN:O	2.08	0.68
1:B:132:LEU:O	1:B:136:THR:HG23	1.95	0.67
1:A:132:LEU:O	1:A:136:THR:HG23	1.95	0.67
1:B:112:ILE:HG23	1:B:145:GLY:HA3	1.78	0.66
1:A:112:ILE:O	1:A:116:THR:N	2.29	0.66
1:B:45:ASP:O	1:B:49:LEU:N	2.23	0.66
1:B:334:GLN:OE1	1:B:423:GLN:NE2	2.29	0.66
1:B:376:THR:OG1	1:B:428:PHE:HA	1.96	0.66
1:B:422:MET:O	1:B:426:PHE:N	2.29	0.66
1:A:489:CYS:O	1:A:493:PHE:N	2.28	0.65
1:A:178:GLY:O	1:A:182:PHE:N	2.26	0.65
1:B:492:GLY:O	1:B:496:THR:OG1	2.10	0.64
1:B:489:CYS:O	1:B:493:PHE:N	2.26	0.64
1:B:330:ILE:O	1:B:334:GLN:N	2.30	0.64
1:B:140:VAL:O	1:B:144:ILE:N	2.24	0.64
1:A:109:LEU:HD12	1:A:112:ILE:HD11	1.79	0.64
1:B:178:GLY:O	1:B:182:PHE:N	2.28	0.63
1:A:464:LEU:O	1:A:468:VAL:HG12	1.98	0.63
1:A:182:PHE:CE1	1:A:362:ILE:HD11	2.33	0.63
1:A:62:VAL:HG22	1:A:207:VAL:HG13	1.81	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:392:GLY:O	1:A:396:SER:OG	2.12	0.62
1:A:331:ASN:OD1	1:A:369:PHE:CE1	2.53	0.62
1:A:391:MET:HA	1:A:394:ILE:HG22	1.82	0.61
1:B:182:PHE:CE1	1:B:362:ILE:HD11	2.36	0.61
1:A:376:THR:OG1	1:A:428:PHE:HA	2.01	0.61
1:B:320:TRP:CD1	1:B:324:GLU:HG2	2.36	0.61
1:A:49:LEU:O	1:A:52:ILE:HG22	2.01	0.61
1:B:81:ALA:O	1:B:85:CYS:HB2	2.01	0.60
1:A:50:PHE:HE2	1:A:177:GLN:HG3	1.66	0.60
1:A:460:CYS:HA	1:A:463:ILE:HG12	1.82	0.60
1:A:81:ALA:O	1:A:85:CYS:HB2	2.02	0.60
1:A:314:LEU:HD22	1:A:498:LEU:HD13	1.83	0.59
1:A:84:GLY:O	1:A:88:GLY:N	2.30	0.59
1:B:331:ASN:OD1	1:B:369:PHE:CE1	2.55	0.59
1:B:391:MET:HA	1:B:394:ILE:HG22	1.84	0.58
1:A:323:VAL:O	1:A:327:PHE:N	2.24	0.58
1:B:330:ILE:HD13	1:B:426:PHE:CE2	2.39	0.58
1:A:224:GLY:O	1:A:228:GLN:HG2	2.04	0.58
1:B:50:PHE:HE2	1:B:177:GLN:HG3	1.67	0.58
1:B:460:CYS:HA	1:B:463:ILE:HG12	1.86	0.58
1:A:322:LEU:HA	1:A:325:ILE:HD12	1.86	0.58
1:B:112:ILE:O	1:B:116:THR:OG1	2.13	0.56
1:A:103:PHE:O	1:A:107:LYS:HD3	2.06	0.56
1:A:179:TRP:O	1:A:183:VAL:N	2.32	0.56
1:A:182:PHE:O	1:A:186:LEU:N	2.34	0.56
1:B:343:PHE:HE2	1:B:416:LEU:HD13	1.70	0.56
1:B:404:ALA:O	1:B:479:SER:OG	2.20	0.56
1:B:87:VAL:HA	1:B:90:VAL:HG12	1.87	0.56
1:A:212:ARG:O	1:A:216:GLY:N	2.32	0.56
1:A:326:ALA:O	1:A:330:ILE:HD12	2.06	0.56
1:A:467:LEU:O	1:A:471:GLN:NE2	2.39	0.56
1:B:52:ILE:HD11	1:B:77:MET:HE3	1.88	0.56
1:B:363:ILE:HG23	1:B:367:LEU:HD12	1.87	0.56
1:B:424:PHE:O	1:B:428:PHE:N	2.33	0.56
1:B:330:ILE:HD13	1:B:426:PHE:HE2	1.70	0.55
1:B:49:LEU:O	1:B:52:ILE:HG22	2.07	0.55
1:B:459:LYS:HA	1:B:462:ALA:HB3	1.88	0.55
1:A:50:PHE:HE2	1:A:177:GLN:CG	2.19	0.55
1:A:321:PHE:O	1:A:325:ILE:HG13	2.06	0.55
1:A:363:ILE:HG23	1:A:367:LEU:HD12	1.87	0.55
1:A:72:GLN:O	1:A:76:LEU:HD13	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:490:ILE:O	1:B:494:ILE:N	2.28	0.55
1:A:139:ARG:O	1:A:143:GLY:N	2.34	0.55
1:A:459:LYS:O	1:A:463:ILE:HG12	2.07	0.54
1:A:310:PHE:CE2	1:A:314:LEU:HD12	2.43	0.54
1:A:43:PHE:O	1:A:47:TYR:N	2.28	0.54
1:A:488:THR:O	1:A:492:GLY:N	2.38	0.54
1:A:321:PHE:CE1	1:A:464:LEU:HD12	2.42	0.53
1:A:47:TYR:O	1:A:51:ILE:HB	2.09	0.53
1:B:161:ALA:HB1	1:B:163:ILE:HG22	1.88	0.53
1:B:179:TRP:O	1:B:183:VAL:N	2.34	0.53
1:B:337:VAL:HG11	1:B:419:PHE:CE1	2.43	0.53
1:A:370:LEU:C	1:A:370:LEU:HD23	2.33	0.53
1:A:422:MET:O	1:A:426:PHE:N	2.41	0.53
1:B:459:LYS:O	1:B:463:ILE:HG12	2.09	0.53
1:A:46:ALA:CB	1:A:177:GLN:HA	2.39	0.53
1:B:182:PHE:O	1:B:186:LEU:N	2.35	0.52
1:A:112:ILE:O	1:A:116:THR:OG1	2.20	0.52
1:A:215:ILE:O	1:A:218:SER:OG	2.22	0.52
1:A:337:VAL:HG11	1:A:419:PHE:CE1	2.45	0.52
1:B:112:ILE:O	1:B:116:THR:N	2.35	0.52
1:A:184:GLY:O	1:A:188:THR:OG1	2.15	0.52
1:B:43:PHE:HZ	1:B:183:VAL:HG11	1.75	0.52
1:B:221:PRO:O	1:B:225:THR:N	2.39	0.52
1:B:441:LEU:HD11	1:B:499:ILE:HD12	1.92	0.52
1:A:463:ILE:O	1:A:467:LEU:HB2	2.10	0.52
1:B:76:LEU:O	1:B:80:ALA:N	2.43	0.51
1:A:50:PHE:CE1	1:A:331:ASN:HB3	2.45	0.51
1:A:43:PHE:HZ	1:A:183:VAL:HG11	1.75	0.51
1:A:62:VAL:CG2	1:A:207:VAL:HG13	2.41	0.51
1:A:211:TRP:CE3	1:A:215:ILE:HD12	2.46	0.51
1:B:393:PHE:CD1	1:B:489:CYS:HA	2.46	0.51
1:B:47:TYR:O	1:B:51:ILE:HB	2.11	0.51
1:B:463:ILE:O	1:B:467:LEU:HB2	2.11	0.51
1:A:424:PHE:O	1:A:428:PHE:N	2.31	0.51
1:B:484:ILE:O	1:B:488:THR:HG23	2.10	0.51
1:B:321:PHE:O	1:B:325:ILE:HG13	2.12	0.50
1:B:365:THR:O	1:B:370:LEU:HB2	2.11	0.50
1:B:468:VAL:HG22	1:B:472:LEU:HG	1.93	0.50
1:B:467:LEU:O	1:B:471:GLN:NE2	2.44	0.50
1:B:463:ILE:O	1:B:467:LEU:N	2.33	0.50
1:A:404:ALA:O	1:A:479:SER:OG	2.23	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:84:GLY:O	1:B:88:GLY:N	2.35	0.50
1:A:217:LEU:O	1:A:221:PRO:HD2	2.11	0.50
1:A:146:ILE:O	1:A:150:TYR:HB2	2.12	0.50
1:A:310:PHE:CZ	1:A:314:LEU:HD12	2.46	0.50
1:B:45:ASP:OD2	1:B:149:ASP:OD2	2.30	0.50
1:A:208:ASP:HA	1:A:211:TRP:CD1	2.47	0.50
1:A:459:LYS:O	1:A:463:ILE:N	2.30	0.50
1:A:189:ILE:HG23	1:A:358:ALA:HB1	1.93	0.49
1:A:110:ILE:O	1:A:114:VAL:HG23	2.12	0.49
1:B:146:ILE:O	1:B:150:TYR:HB2	2.13	0.49
1:B:396:SER:O	1:B:400:LEU:N	2.42	0.49
1:B:488:THR:O	1:B:492:GLY:N	2.38	0.49
1:A:83:ILE:HD11	1:A:467:LEU:HD21	1.95	0.49
1:A:87:VAL:HA	1:A:90:VAL:HG12	1.94	0.49
1:B:370:LEU:HD21	1:B:374:TYR:HE2	1.76	0.49
1:A:459:LYS:HA	1:A:462:ALA:HB3	1.95	0.49
1:B:50:PHE:HE2	1:B:177:GLN:CG	2.25	0.49
1:A:403:LEU:O	1:A:407:ILE:HB	2.13	0.48
1:B:227:TYR:HB2	1:A:118:PHE:CE2	2.48	0.48
1:B:139:ARG:O	1:B:143:GLY:N	2.36	0.48
1:B:178:GLY:HA2	1:B:369:PHE:CE2	2.48	0.48
1:B:394:ILE:O	1:B:398:LEU:HG	2.12	0.48
1:A:88:GLY:HA2	1:A:144:ILE:HG23	1.94	0.48
1:B:212:ARG:O	1:B:216:GLY:N	2.37	0.48
1:B:320:TRP:CD1	1:B:324:GLU:CG	2.97	0.48
1:A:109:LEU:CD1	1:A:112:ILE:HD11	2.43	0.48
1:B:85:CYS:SG	1:B:459:LYS:NZ	2.63	0.48
1:B:322:LEU:HA	1:B:325:ILE:HD12	1.96	0.47
1:A:76:LEU:O	1:A:80:ALA:N	2.46	0.47
1:A:369:PHE:O	1:A:427:ASN:ND2	2.38	0.47
1:A:463:ILE:O	1:A:467:LEU:N	2.41	0.47
1:B:211:TRP:CE3	1:B:215:ILE:HD12	2.50	0.47
1:B:392:GLY:HA3	1:B:429:GLY:HA3	1.97	0.47
1:B:224:GLY:O	1:B:228:GLN:HG2	2.15	0.47
1:B:317:MET:O	1:B:321:PHE:CB	2.62	0.47
1:A:330:ILE:HD13	1:A:426:PHE:CE2	2.49	0.47
1:B:86:VAL:HA	1:B:89:GLN:HG2	1.97	0.47
1:B:103:PHE:O	1:B:107:LYS:HD3	2.14	0.47
1:B:362:ILE:O	1:B:366:ALA:N	2.36	0.47
1:B:341:ILE:HG22	1:B:342:GLY:H	1.79	0.47
1:B:311:ARG:NH1	1:B:500:ASP:OD1	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:GLN:HG3	1:A:135:ILE:HG23	1.96	0.47
1:B:359:THR:HA	1:B:362:ILE:HG22	1.96	0.47
1:A:463:ILE:HA	1:A:466:SER:HB2	1.97	0.47
1:B:326:ALA:O	1:B:330:ILE:HD12	2.14	0.46
1:B:217:LEU:O	1:B:221:PRO:HD2	2.16	0.46
1:B:324:GLU:OE1	1:B:328:TYR:CD2	2.68	0.46
1:B:369:PHE:O	1:B:427:ASN:ND2	2.38	0.46
1:B:327:PHE:CE1	1:B:331:ASN:ND2	2.83	0.46
1:A:421:PHE:O	1:A:425:PHE:CD2	2.69	0.46
1:A:462:ALA:O	1:A:466:SER:N	2.37	0.46
1:B:318:LEU:O	1:B:322:LEU:N	2.37	0.46
1:A:311:ARG:NH2	1:A:500:ASP:OD1	2.40	0.46
1:A:324:GLU:HA	1:A:327:PHE:HB3	1.97	0.46
1:A:499:ILE:HG22	1:A:500:ASP:O	2.15	0.46
1:B:310:PHE:CZ	1:B:314:LEU:HD12	2.51	0.46
1:B:320:TRP:O	1:B:324:GLU:HG2	2.16	0.46
1:B:359:THR:O	1:B:362:ILE:HG22	2.15	0.46
1:A:370:LEU:HD21	1:A:374:TYR:HE2	1.79	0.46
1:B:362:ILE:O	1:B:366:ALA:CB	2.64	0.46
1:A:43:PHE:O	1:A:47:TYR:CB	2.64	0.46
1:A:320:TRP:HE1	1:A:324:GLU:CD	2.24	0.46
1:B:43:PHE:O	1:B:47:TYR:N	2.34	0.45
1:A:32:ILE:O	1:A:35:VAL:HG22	2.17	0.45
1:A:365:THR:O	1:A:370:LEU:CB	2.65	0.45
1:B:88:GLY:HA3	1:B:144:ILE:HA	1.97	0.45
1:B:390:PHE:HD1	1:B:493:PHE:CD1	2.34	0.45
1:B:462:ALA:O	1:B:466:SER:CB	2.64	0.45
1:A:47:TYR:O	1:A:51:ILE:N	2.41	0.45
1:A:84:GLY:O	1:A:144:ILE:HA	2.16	0.45
1:B:109:LEU:HA	1:B:112:ILE:HG12	1.99	0.45
1:A:359:THR:HA	1:A:362:ILE:HG22	1.98	0.45
1:B:115:ALA:O	1:B:119:GLN:HB2	2.16	0.45
1:B:323:VAL:HG11	1:B:434:THR:OG1	2.17	0.45
1:A:314:LEU:HD23	1:A:314:LEU:C	2.42	0.45
1:A:468:VAL:HG22	1:A:472:LEU:HG	1.99	0.45
1:B:216:GLY:O	1:B:219:LEU:HB2	2.16	0.45
1:A:366:ALA:O	1:A:371:PRO:HD3	2.17	0.45
1:B:83:ILE:HD11	1:B:467:LEU:HD21	1.98	0.45
1:A:115:ALA:O	1:A:119:GLN:HB2	2.17	0.45
1:B:213:ILE:O	1:B:217:LEU:HG	2.17	0.45
1:B:393:PHE:O	1:B:397:GLY:N	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:455:ALA:O	1:B:459:LYS:HG3	2.17	0.45
1:A:343:PHE:HE2	1:A:416:LEU:HD13	1.82	0.45
1:A:373:TYR:CE2	1:A:427:ASN:ND2	2.85	0.44
1:A:393:PHE:CD1	1:A:489:CYS:HA	2.52	0.44
1:B:79:ALA:O	1:B:83:ILE:N	2.37	0.44
1:B:116:THR:O	1:B:120:MET:N	2.40	0.44
1:B:375:PHE:CD1	1:B:424:PHE:CE1	3.05	0.44
1:A:327:PHE:CE1	1:A:331:ASN:ND2	2.85	0.44
1:B:112:ILE:HB	1:B:142:LEU:HD12	1.99	0.44
1:B:112:ILE:HG22	1:B:142:LEU:HD12	1.99	0.44
1:A:492:GLY:O	1:A:496:THR:OG1	2.23	0.44
1:B:463:ILE:HA	1:B:466:SER:HB2	1.98	0.44
1:A:462:ALA:O	1:A:466:SER:CB	2.65	0.44
1:B:310:PHE:CE2	1:B:314:LEU:HD12	2.53	0.44
1:B:365:THR:O	1:B:370:LEU:CB	2.66	0.44
1:B:317:MET:O	1:B:321:PHE:HB3	2.17	0.43
1:A:359:THR:O	1:A:362:ILE:HG22	2.18	0.43
1:A:410:ILE:HG13	1:A:411:GLY:H	1.82	0.43
1:B:47:TYR:O	1:B:51:ILE:N	2.38	0.43
1:B:50:PHE:CE1	1:B:331:ASN:HB3	2.54	0.43
1:B:321:PHE:CE1	1:B:464:LEU:HD12	2.53	0.43
1:B:490:ILE:HA	1:B:493:PHE:HB3	2.00	0.43
1:A:365:THR:O	1:A:370:LEU:HB2	2.18	0.43
1:A:396:SER:O	1:A:400:LEU:N	2.43	0.43
1:B:46:ALA:CB	1:B:177:GLN:HA	2.48	0.43
1:A:455:ALA:O	1:A:459:LYS:HG3	2.18	0.43
1:B:44:LEU:HD12	1:B:218:SER:HB2	2.01	0.43
1:B:45:ASP:CG	1:B:146:ILE:HG23	2.44	0.43
1:A:341:ILE:HG22	1:A:342:GLY:H	1.83	0.43
1:B:366:ALA:O	1:B:371:PRO:HD3	2.19	0.43
1:B:456:ALA:O	1:B:460:CYS:N	2.40	0.43
1:A:424:PHE:O	1:A:428:PHE:CB	2.67	0.43
1:A:424:PHE:O	1:A:428:PHE:HB3	2.18	0.43
1:B:424:PHE:O	1:B:428:PHE:CB	2.66	0.43
1:B:446:ILE:C	1:B:446:ILE:HD12	2.44	0.43
1:A:58:MET:SD	1:A:214:LEU:HD11	2.59	0.43
1:A:375:PHE:CD1	1:A:424:PHE:CE1	3.06	0.43
1:A:390:PHE:O	1:A:394:ILE:HG22	2.19	0.43
1:B:112:ILE:O	1:B:116:THR:CB	2.67	0.43
1:B:227:TYR:OH	1:A:121:SER:O	2.22	0.43
1:A:316:SER:N	1:A:499:ILE:HD11	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:161:ALA:HB1	1:A:163:ILE:HG22	2.01	0.43
1:A:182:PHE:O	1:A:185:SER:N	2.52	0.43
1:B:113:ILE:O	1:B:117:ILE:N	2.39	0.42
1:B:94:VAL:O	1:B:98:SER:N	2.49	0.42
1:B:184:GLY:O	1:B:188:THR:OG1	2.25	0.42
1:A:109:LEU:HA	1:A:112:ILE:HG12	2.01	0.42
1:A:112:ILE:HG23	1:A:145:GLY:HA3	2.01	0.42
1:A:116:THR:O	1:A:120:MET:N	2.40	0.42
1:B:390:PHE:CE1	1:B:496:THR:HB	2.54	0.42
1:B:393:PHE:HD1	1:B:489:CYS:HA	1.84	0.42
1:B:459:LYS:O	1:B:463:ILE:N	2.30	0.42
1:B:56:ALA:N	1:B:57:PRO:HD2	2.35	0.42
1:A:330:ILE:HD13	1:A:426:PHE:HE2	1.85	0.42
1:B:43:PHE:O	1:B:47:TYR:CB	2.68	0.42
1:A:88:GLY:HA3	1:A:144:ILE:HA	2.00	0.42
1:B:410:ILE:HG13	1:B:411:GLY:H	1.84	0.42
1:B:424:PHE:O	1:B:428:PHE:HB3	2.19	0.42
1:A:140:VAL:O	1:A:144:ILE:HG13	2.20	0.42
1:A:211:TRP:HE3	1:A:215:ILE:HD12	1.84	0.42
1:A:323:VAL:HG13	1:A:324:GLU:N	2.35	0.42
1:B:58:MET:SD	1:B:214:LEU:HD11	2.60	0.42
1:B:422:MET:O	1:B:425:PHE:N	2.53	0.42
1:A:50:PHE:O	1:A:53:ASN:HB3	2.20	0.42
1:A:323:VAL:HG11	1:A:434:THR:OG1	2.20	0.42
1:B:178:GLY:HA2	1:B:369:PHE:HE2	1.85	0.41
1:B:460:CYS:O	1:B:464:LEU:HG	2.20	0.41
1:B:54:GLN:NE2	1:B:185:SER:HA	2.35	0.41
1:A:112:ILE:HB	1:A:142:LEU:HD12	2.02	0.41
1:A:321:PHE:CZ	1:A:464:LEU:HD12	2.54	0.41
1:B:399:PHE:O	1:B:403:LEU:HB2	2.20	0.41
1:A:362:ILE:O	1:A:366:ALA:CB	2.69	0.41
1:A:369:PHE:CZ	1:A:373:TYR:OH	2.73	0.41
1:A:179:TRP:HA	1:A:182:PHE:HB3	2.02	0.41
1:A:484:ILE:O	1:A:488:THR:HG23	2.20	0.41
1:B:50:PHE:CE2	1:B:177:GLN:CG	3.03	0.41
1:B:89:GLN:OE1	1:B:459:LYS:HE3	2.21	0.41
1:B:156:VAL:HG23	1:B:157:VAL:HG13	2.03	0.41
1:B:373:TYR:CE2	1:B:427:ASN:ND2	2.89	0.41
1:A:56:ALA:N	1:A:57:PRO:HD2	2.36	0.41
1:A:112:ILE:O	1:A:116:THR:CB	2.69	0.41
1:B:32:ILE:O	1:B:35:VAL:HG22	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:83:ILE:O	1:B:86:VAL:HG12	2.20	0.41
1:A:42:PHE:O	1:A:46:ALA:N	2.31	0.41
1:B:43:PHE:O	1:B:47:TYR:HB2	2.21	0.41
1:B:375:PHE:CD1	1:B:424:PHE:HE1	2.38	0.41
1:A:43:PHE:O	1:A:47:TYR:HB2	2.21	0.41
1:A:77:MET:CB	1:A:136:THR:HG22	2.51	0.41
1:A:107:LYS:HA	1:A:110:ILE:HG12	2.03	0.41
1:A:422:MET:O	1:A:425:PHE:N	2.54	0.41
1:A:436:ILE:HG22	1:A:440:GLU:OE1	2.20	0.41
1:B:88:GLY:CA	1:B:144:ILE:HA	2.50	0.41
1:B:421:PHE:O	1:B:425:PHE:CD2	2.74	0.41
1:A:77:MET:HB2	1:A:136:THR:HG22	2.03	0.41
1:A:441:LEU:HD11	1:A:499:ILE:HD12	2.02	0.41
1:B:88:GLY:O	1:B:92:PHE:HB2	2.21	0.40
1:B:376:THR:HG21	1:B:432:THR:HG21	2.03	0.40
1:A:79:ALA:O	1:A:83:ILE:N	2.40	0.40
1:B:107:LYS:HA	1:B:110:ILE:HG12	2.02	0.40
1:B:135:ILE:O	1:B:139:ARG:HG3	2.20	0.40
1:A:47:TYR:OH	1:A:214:LEU:O	2.37	0.40
1:A:370:LEU:HD21	1:A:374:TYR:CE2	2.55	0.40
1:B:62:VAL:CG2	1:B:207:VAL:HG13	2.52	0.40
1:B:189:ILE:HG23	1:B:358:ALA:HB1	2.03	0.40
1:B:140:VAL:O	1:B:144:ILE:HG13	2.22	0.40
1:B:370:LEU:CD2	1:B:374:TYR:HE2	2.35	0.40
1:B:397:GLY:O	1:B:401:ALA:N	2.42	0.40
1:A:58:MET:CE	1:A:214:LEU:HD11	2.52	0.40
1:A:111:LEU:O	1:A:115:ALA:CB	2.70	0.40
1:A:112:ILE:HB	1:A:142:LEU:CD1	2.52	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	423/530 (80%)	411 (97%)	12 (3%)	0	100	100
1	B	423/530 (80%)	413 (98%)	10 (2%)	0	100	100
All	All	846/1060 (80%)	824 (97%)	22 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	338/433 (78%)	338 (100%)	0	100	100
1	B	338/433 (78%)	338 (100%)	0	100	100
All	All	676/866 (78%)	676 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	119	GLN
1	B	331	ASN
1	B	334	GLN
1	B	423	GLN
1	A	119	GLN
1	A	129	ASN
1	A	203	HIS
1	A	331	ASN
1	A	334	GLN
1	A	423	GLN
1	A	470	ASN
1	A	471	GLN

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	429/530 (80%)	-0.54	4 (0%) 81 64	127, 174, 284, 372	0
1	B	429/530 (80%)	-0.60	3 (0%) 84 68	124, 177, 284, 374	0
All	All	858/1060 (80%)	-0.57	7 (0%) 82 65	124, 175, 284, 374	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	203	HIS	4.3
1	B	290	ILE	3.2
1	A	231	THR	2.5
1	A	162	ASN	2.4
1	B	409	HIS	2.3
1	A	206	ASP	2.1
1	B	99	PHE	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.