



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 25, 2021 – 10:18 AM EDT

PDB ID : 7MI1
Title : X-ray structure of yeast dynein motor domain in the presence of a pyrazolo-pyrimidinone-based compound (compound 20)
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Deposited on : 2021-04-16
Resolution : 4.50 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.23.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.23.2

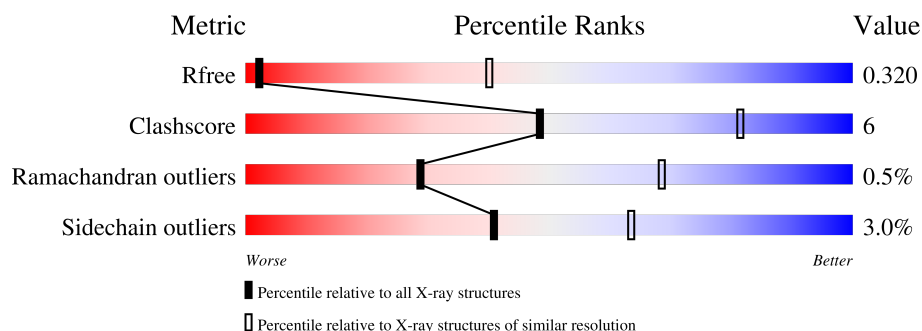
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


The reported resolution of this entry is 4.50 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}	130704	1055 (5.20-3.80)
Clashscore	141614	1123 (5.20-3.80)
Ramachandran outliers	138981	1069 (5.20-3.80)
Sidechain outliers	138945	1050 (5.20-3.80)

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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	2661	 81% 16% ..

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 42528 atoms, of which 21316 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 Chimera protein of Dynein and Endolysin.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	2628	42528	13593	21316	3538	3984	97	0	0	0

There are 31 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1363	GLY	-	expression tag	UNP P36022
A	1849	GLN	GLU	engineered mutation	UNP P36022
A	3120	GLY	-	linker	UNP P36022
A	3121	SER	-	linker	UNP P36022
A	3122	GLY	-	linker	UNP P36022
A	3123	SER	-	linker	UNP P36022
A	3124	GLY	-	linker	UNP P36022
A	3125	SER	-	linker	UNP P36022
A	3178	THR	CYS	conflict	UNP D9IEF7
A	3221	ALA	CYS	conflict	UNP D9IEF7
A	3286	GLY	-	linker	UNP D9IEF7
A	3287	SER	-	linker	UNP D9IEF7
A	3288	GLY	-	linker	UNP D9IEF7
A	3289	SER	-	linker	UNP D9IEF7
A	3290	GLY	-	linker	UNP D9IEF7
A	3291	SER	-	linker	UNP D9IEF7
A	3742	ASP	ASN	conflict	UNP P36022
A	3895	VAL	PHE	conflict	UNP P36022
A	4072	ASP	ASN	conflict	UNP P36022
A	4093	GLY	-	expression tag	UNP P36022
A	4094	SER	-	expression tag	UNP P36022
A	4095	GLY	-	expression tag	UNP P36022
A	4096	SER	-	expression tag	UNP P36022
A	4097	GLY	-	expression tag	UNP P36022
A	4098	SER	-	expression tag	UNP P36022
A	4099	HIS	-	expression tag	UNP P36022
A	4100	HIS	-	expression tag	UNP P36022

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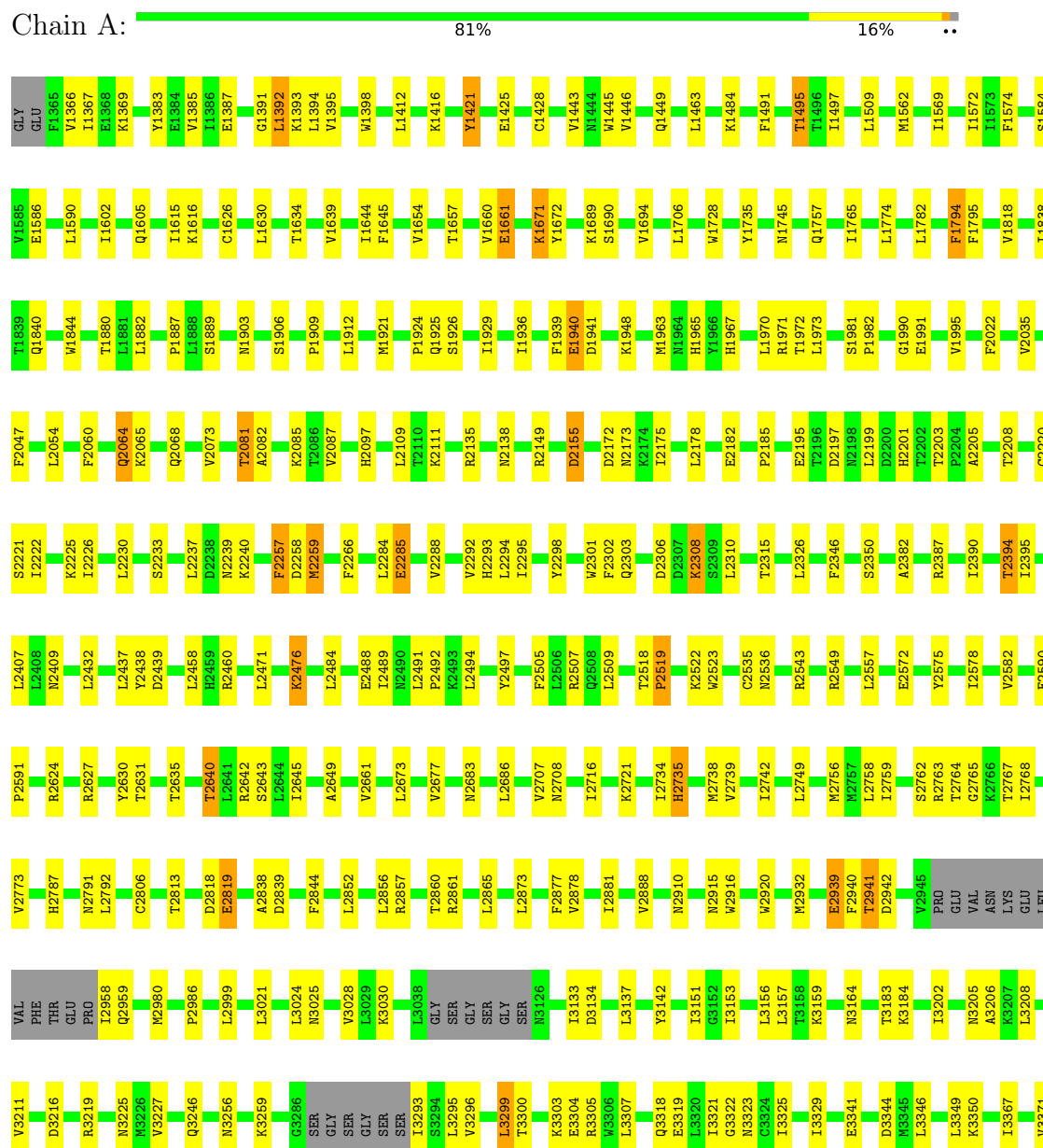
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Chain	Residue	Modelled	Actual	Comment	Reference
A	4101	HIS	-	expression tag	UNP P36022
A	4102	HIS	-	expression tag	UNP P36022
A	4103	HIS	-	expression tag	UNP P36022
A	4104	HIS	-	expression tag	UNP P36022

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. 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. 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: Chimera protein of Dynein and Endolysin



V4081	D3899	V3725	F3611	F3406
T4085	I3903	L3726	V3615	L3407
L4089	L3904	S3727	I3628	L3408
D4093	D3905	E3728	L3632	D3409
S4094	T3906	D3731	F3641	S3412
H4099	F3915	D3743	I3644	T3415
HIS	S3925	L3744	Q3648	L3429
HIS	F3935	R3745	R3655	E3433
HIS	I3939	R3746	V3656	R3439
HIS	V3946	L3747	F3657	V3449
	P3947	V3769	I3658	V3450
	K3953	A3775	K3659	D3459
D3958	D3958	D3776	K3660	F3460
C3959	C3959	V3777	E3663	I3461
D3960	D3960	T3787	T3664	I3466
E3969	E3969	L3803	R3665	K3493
R3970	R3970	E3809	ALA	H3497
V3971	V3971	S3810	ARG	P3501
T3974	T3974	L3811	R3670	L3509
N3975	N3975	K3833	V3671	V3513
I3980	I3980	I3844	D3672	R3517
P3981	P3981	V3851	E3673	F3518
W3982	W3982	L3855	I3674	N3538
R3986	R3986	H3858	L3675	K3541
D3987	D3987	T3862	L3677	L3570
A4012	A4012	K3863	Y3678	Q3574
H4013	H4013	A3864	Y3679	K3577
V4014	V4014	H3868	Q3680	L3578
S4018	S4018	F3871	E3681	E3579
D4019	D4019	K3875	C3684	K3580
N4020	N4020	A3886	T3688	L3583
L4021	L4021	P3887	A3689	K3584
Q4022	Q4022	L3888	L3690	L3587
I4023	I4023	L3889	F3694	K3591
L4033	L4033	R3894	F3703	K3595
L4034	L4034	V3895	Y3706	
Q4035	Q4035	V3896	I3710	
D4072	D4072		Q3714	
R4075	R4075		V3719	
A4078	A4078		L3720	
			T3721	

4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2 ₁	Depositor
Cell constants a, b, c, α , β , γ	135.06Å 157.92Å 179.31Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.66 – 4.50 47.66 – 4.50	Depositor EDS
% Data completeness (in resolution range)	99.5 (47.66-4.50) 99.5 (47.66-4.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.14 (at 4.45Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.276 , 0.320 0.276 , 0.320	Depositor DCC
R_{free} test set	1997 reflections (8.61%)	wwPDB-VP
Wilson B-factor (Å ²)	183.0	Xtriage
Anisotropy	0.178	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 122.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.22$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	42528	wwPDB-VP
Average B, all atoms (Å ²)	228.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.50% 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 [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	0.26	0/21632	0.44	1/29225 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	2941	THR	N-CA-C	5.17	124.97	111.00

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	21212	21316	21315	257	0
All	All	21212	21316	21315	257	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 6.

All (257) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (\AA)	Clash overlap (\AA)
1:A:3665:ARG:NH1	1:A:3673:GLU:OE2	1.97	0.97
1:A:2939:GLU:OE1	1:A:3318:GLN:NE2	2.05	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2624:ARG:NH2	1:A:2910:ASN:O	2.08	0.86
1:A:2064:GLN:OE1	1:A:2065:LYS:N	2.15	0.79
1:A:2677:VAL:HG11	1:A:2686:LEU:HD21	1.64	0.77
1:A:1939:PHE:O	1:A:1941:ASP:N	2.19	0.75
1:A:3660:LYS:O	1:A:3660:LYS:NZ	2.21	0.74
1:A:2458:LEU:HD11	1:A:2484:LEU:HD11	1.69	0.73
1:A:1412:LEU:HD22	1:A:1428:CYS:SG	2.29	0.73
1:A:2155:ASP:OD2	1:A:2507:ARG:NH2	2.22	0.73
1:A:3787:THR:HG22	1:A:3875:MET:HB2	1.69	0.73
1:A:2135:ARG:NH2	1:A:2182:GLU:OE1	2.22	0.73
1:A:2877:PHE:CZ	1:A:2881:ILE:HD11	2.28	0.69
1:A:3946:VAL:HG11	1:A:3953:LYS:HB2	1.74	0.69
1:A:2201:HIS:NE2	1:A:2497:TYR:O	2.25	0.69
1:A:1924:PRO:HB2	1:A:1929:ILE:HD11	1.76	0.67
1:A:3809:GLU:O	1:A:3811:LEU:N	2.28	0.67
1:A:1967:HIS:O	1:A:1972:THR:OG1	2.07	0.67
1:A:2787:HIS:HB3	1:A:3461:ILE:HG23	1.77	0.67
1:A:2460:ARG:NH2	1:A:2839:ASP:OD2	2.28	0.66
1:A:2258:ASP:O	1:A:2259:MET:HB2	1.96	0.65
1:A:2489:ILE:HG22	1:A:2535:CYS:HB3	1.78	0.65
1:A:2838:ALA:HB3	1:A:2878:VAL:HG13	1.79	0.64
1:A:4072:ASP:OD1	1:A:4075:ARG:NH2	2.31	0.64
1:A:2220:CYS:SG	1:A:2221:SER:N	2.71	0.64
1:A:3466:ILE:HD13	1:A:3509:LEU:CD1	2.29	0.63
1:A:2060:PHE:HD2	1:A:2087:VAL:HG11	1.62	0.62
1:A:2222:ILE:HG23	1:A:2284:LEU:HD11	1.80	0.62
1:A:3307:LEU:HD12	1:A:3307:LEU:O	1.99	0.62
1:A:2081:THR:HG22	1:A:2085:LYS:HE3	1.84	0.60
1:A:3296:VAL:HA	1:A:3584:MET:SD	2.42	0.60
1:A:3833:LYS:NZ	1:A:3862:THR:HG21	2.17	0.60
1:A:2258:ASP:O	1:A:2259:MET:CB	2.50	0.59
1:A:2758:LEU:HD23	1:A:2915:ASN:HB3	1.85	0.59
1:A:3628:ILE:HD11	1:A:3679:TYR:CZ	2.38	0.59
1:A:2631:THR:O	1:A:2635:THR:HG22	2.04	0.58
1:A:2806:CYS:HG	1:A:2813:THR:HG1	1.45	0.58
1:A:2138:ASN:ND2	1:A:2185:PRO:O	2.37	0.57
1:A:2230:LEU:HD11	1:A:2257:PHE:HZ	1.69	0.57
1:A:3028:VAL:HG12	1:A:3300:THR:HB	1.86	0.57
1:A:3429:LEU:HD21	1:A:3439:ARG:HB3	1.85	0.57
1:A:1425:GLU:O	1:A:1428:CYS:N	2.37	0.56
1:A:1929:ILE:HD13	1:A:1970:LEU:HD11	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:3346:LEU:HD11	1:A:3350:LYS:HE3	1.87	0.56
1:A:1706:LEU:HD12	1:A:1936:ILE:HD13	1.86	0.56
1:A:1562:MET:CB	1:A:1569:ILE:HD11	2.36	0.56
1:A:1661:GLU:OE2	1:A:1735:TYR:OH	2.22	0.56
1:A:2582:VAL:HG23	1:A:2582:VAL:O	2.06	0.56
1:A:1745:ASN:OD1	1:A:1757:GLN:NE2	2.40	0.56
1:A:1940:GLU:OE1	1:A:1990:GLY:HA2	2.06	0.56
1:A:3641:PHE:HA	1:A:3889:LEU:HD21	1.88	0.55
1:A:2458:LEU:HD21	1:A:2484:LEU:HD21	1.89	0.55
1:A:3935:PHE:HB2	1:A:4014:VAL:HG11	1.89	0.54
1:A:2856:LEU:HD23	1:A:2873:LEU:HB3	1.88	0.54
1:A:2203:THR:HG22	1:A:2205:ALA:H	1.71	0.54
1:A:2818:ASP:OD1	1:A:2819:GLU:N	2.41	0.54
1:A:2197:ASP:N	1:A:2197:ASP:OD1	2.39	0.54
1:A:1991:GLU:O	1:A:1995:VAL:HG23	2.07	0.54
1:A:1925:GLN:O	1:A:1929:ILE:HD12	2.08	0.54
1:A:2522:LYS:HG2	1:A:2523:TRP:H	1.73	0.54
1:A:3211:VAL:HG22	1:A:3246:GLN:HB3	1.90	0.53
1:A:2536:ASN:HB2	1:A:2543:ARG:HE	1.73	0.53
1:A:2762:SER:O	1:A:2764:THR:N	2.42	0.53
1:A:2394:THR:HG22	1:A:2395:ILE:HD12	1.90	0.53
1:A:3971:VAL:HA	1:A:3974:THR:HG22	1.90	0.53
1:A:1366:VAL:HG13	1:A:1369:LYS:HE3	1.91	0.53
1:A:2173:ASN:HB3	1:A:2175:ILE:HG22	1.90	0.52
1:A:1971:ARG:HB2	1:A:2208:THR:HG21	1.90	0.52
1:A:3833:LYS:HZ3	1:A:3862:THR:HG21	1.72	0.52
1:A:2407:LEU:HD12	1:A:2557:LEU:HD11	1.92	0.52
1:A:1395:VAL:HG23	1:A:1398:TRP:NE1	2.25	0.51
1:A:2471:LEU:HD12	1:A:2471:LEU:H	1.76	0.51
1:A:3895:VAL:HG12	1:A:3896:VAL:N	2.26	0.51
1:A:3134:ASP:OD2	1:A:3225:ASN:ND2	2.41	0.51
1:A:3980:ILE:N	1:A:3981:PRO:CD	2.73	0.51
1:A:3775:ALA:HB2	1:A:3803:LEU:HD22	1.93	0.51
1:A:1367:ILE:H	1:A:1367:ILE:HD12	1.76	0.51
1:A:3216:ASP:OD2	1:A:3219:ARG:NH1	2.43	0.51
1:A:3703:PHE:HE2	1:A:3719:VAL:HG21	1.75	0.51
1:A:1840:GLN:OE1	1:A:1887:PRO:HG2	2.11	0.51
1:A:2940:PHE:CB	1:A:3318:GLN:OE1	2.58	0.51
1:A:2073:VAL:HG21	1:A:2199:LEU:HD11	1.93	0.50
1:A:3322:GLY:HA2	1:A:3325:ILE:HD12	1.93	0.50
1:A:3433:GLU:OE2	1:A:3439:ARG:NH2	2.44	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1909:PRO:HD2	1:A:1912:LEU:HD12	1.93	0.50
1:A:3329:ILE:HG21	1:A:3349:LEU:HD22	1.92	0.50
1:A:3299:LEU:HD11	1:A:3587:LEU:CB	2.40	0.50
1:A:2315:THR:HG21	1:A:2350:SER:HB3	1.94	0.50
1:A:1387:GLU:HA	1:A:1393:LYS:HA	1.94	0.50
1:A:2488:GLU:HB3	1:A:2491:LEU:HD12	1.94	0.50
1:A:1671:LYS:HD3	1:A:1672:TYR:N	2.26	0.49
1:A:3202:ILE:HD11	1:A:3227:VAL:HG21	1.93	0.49
1:A:3367:ILE:O	1:A:3371:VAL:HG22	2.12	0.49
1:A:2294:LEU:O	1:A:2298:TYR:CD2	2.66	0.49
1:A:2522:LYS:HG2	1:A:2523:TRP:N	2.28	0.49
1:A:1981:SER:HB3	1:A:1982:PRO:HD3	1.94	0.49
1:A:3304:GLU:O	1:A:3307:LEU:HD23	2.13	0.49
1:A:3925:SER:OG	1:A:3969:GLU:OE2	2.29	0.49
1:A:2306:ASP:O	1:A:2310:LEU:HB2	2.12	0.49
1:A:3293:ILE:O	1:A:3296:VAL:HG12	2.12	0.49
1:A:2237:LEU:O	1:A:2240:LYS:O	2.31	0.49
1:A:1630:LEU:HD23	1:A:1634:THR:CG2	2.43	0.49
1:A:3684:CYS:HB2	1:A:3769:VAL:HG11	1.95	0.49
1:A:3412:SER:O	1:A:3415:ILE:HG22	2.12	0.48
1:A:4014:VAL:O	1:A:4021:LEU:HD21	2.13	0.48
1:A:2178:LEU:HD12	1:A:2182:GLU:HB2	1.96	0.48
1:A:1445:TRP:HB2	1:A:1509:LEU:HD13	1.96	0.48
1:A:3429:LEU:HD13	1:A:3450:VAL:HG13	1.94	0.48
1:A:3706:TYR:CZ	1:A:3710:ILE:CD1	2.96	0.48
1:A:3760:LEU:HD21	1:A:4078:ALA:HA	1.95	0.48
1:A:1838:ILE:HD11	1:A:1844:TRP:C	2.34	0.48
1:A:2109:LEU:HD22	1:A:2518:THR:HG22	1.95	0.48
1:A:3570:LEU:HD21	1:A:3583:LEU:HD21	1.95	0.48
1:A:3580:ASN:O	1:A:3584:MET:HG2	2.14	0.48
1:A:3648:GLN:OE1	1:A:3894:ARG:NH2	2.46	0.48
1:A:1562:MET:HB2	1:A:1569:ILE:HD11	1.95	0.48
1:A:3299:LEU:HD11	1:A:3587:LEU:HB3	1.95	0.48
1:A:3725:VAL:HG22	1:A:3731:ASP:HA	1.96	0.48
1:A:2762:SER:C	1:A:2764:THR:H	2.17	0.48
1:A:3678:LEU:HD23	1:A:3678:LEU:C	2.34	0.48
1:A:2233:SER:HB3	1:A:2292:VAL:HG11	1.96	0.47
1:A:2394:THR:HG22	1:A:2395:ILE:H	1.78	0.47
1:A:2791:ASN:OD1	1:A:2792:LEU:N	2.48	0.47
1:A:3580:ASN:OD1	1:A:3580:ASN:N	2.46	0.47
1:A:2035:VAL:HG23	1:A:2054:LEU:HD11	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2940:PHE:HB2	1:A:3318:GLN:OE1	2.15	0.47
1:A:3449:VAL:HG13	1:A:3493:LYS:HB2	1.96	0.47
1:A:1416:LYS:HA	1:A:1421:TYR:CE1	2.50	0.47
1:A:3408:LEU:HD21	1:A:3501:PRO:HG3	1.96	0.47
1:A:3658:ILE:HD12	1:A:3658:ILE:N	2.30	0.47
1:A:2980:MET:HE3	1:A:3341:GLU:HB3	1.96	0.47
1:A:3024:LEU:HD11	1:A:3303:LYS:HE2	1.97	0.47
1:A:4033:LEU:HD13	1:A:4035:GLN:HB2	1.96	0.47
1:A:3133:ILE:HD11	1:A:3295:LEU:HD11	1.96	0.47
1:A:3151:ILE:HB	1:A:3157:LEU:HD11	1.96	0.47
1:A:3321:ILE:HD12	1:A:3321:ILE:H	1.79	0.47
1:A:1495:THR:HG22	1:A:1497:ILE:HG22	1.97	0.46
1:A:1630:LEU:HA	1:A:1634:THR:HG22	1.97	0.46
1:A:1366:VAL:HG12	1:A:1366:VAL:O	2.15	0.46
1:A:3030:LYS:NZ	1:A:3574:GLN:OE1	2.48	0.46
1:A:2640:THR:HG23	1:A:2643:SER:H	1.81	0.46
1:A:1590:LEU:HD13	1:A:1644:ILE:HD13	1.97	0.46
1:A:1782:LEU:HD21	1:A:1794:PHE:CZ	2.50	0.46
1:A:2518:THR:OG1	1:A:2519:PRO:HD3	2.15	0.46
1:A:3587:LEU:O	1:A:3591:LYS:HG2	2.14	0.46
1:A:2195:GLU:OE2	1:A:2549:ARG:NH2	2.49	0.46
1:A:2734:ILE:H	1:A:2734:ILE:HD12	1.81	0.46
1:A:3663:GLU:C	1:A:3665:ARG:H	2.19	0.46
1:A:2432:LEU:HD22	1:A:2438:TYR:HB2	1.98	0.46
1:A:3690:LEU:HD23	1:A:3694:PHE:HB3	1.98	0.46
1:A:3183:THR:OG1	1:A:3184:LYS:N	2.50	0.45
1:A:1562:MET:HB3	1:A:1569:ILE:HD11	1.97	0.45
1:A:1586:GLU:HG3	1:A:1765:ILE:H	1.82	0.45
1:A:1939:PHE:O	1:A:1940:GLU:C	2.54	0.45
1:A:3024:LEU:O	1:A:3028:VAL:HG13	2.16	0.45
1:A:1385:VAL:HG21	1:A:1491:PHE:CD1	2.51	0.45
1:A:1392:LEU:HD23	1:A:1484:LYS:HG2	1.99	0.45
1:A:3021:LEU:HA	1:A:3024:LEU:HD12	1.99	0.45
1:A:3025:ASN:O	1:A:3028:VAL:HG22	2.17	0.45
1:A:3656:VAL:HG22	1:A:3681:GLU:HG3	1.99	0.45
1:A:3864:ALA:O	1:A:3868:HIS:N	2.49	0.45
1:A:3721:THR:O	1:A:3725:VAL:HG23	2.16	0.45
1:A:1569:ILE:HA	1:A:1584:SER:HA	1.98	0.44
1:A:2575:TYR:HA	1:A:2578:ILE:HG12	1.99	0.44
1:A:3406:PHE:HB2	1:A:3513:VAL:HG11	1.99	0.44
1:A:3688:THR:HG21	1:A:3777:VAL:HG21	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2060:PHE:CD2	1:A:2087:VAL:HG11	2.48	0.44
1:A:3137:LEU:HD13	1:A:3153:ILE:HG13	1.99	0.44
1:A:2293:HIS:NE2	1:A:2409:ASN:HB3	2.32	0.44
1:A:1995:VAL:HG22	1:A:2022:PHE:CE2	2.53	0.44
1:A:2742:ILE:HG23	1:A:2773:VAL:HG22	1.99	0.44
1:A:1882:LEU:HG	1:A:1882:LEU:O	2.17	0.44
1:A:3202:ILE:HG23	1:A:3208:LEU:HB2	1.99	0.44
1:A:2877:PHE:CE2	1:A:2881:ILE:HD11	2.53	0.44
1:A:3307:LEU:HD12	1:A:3307:LEU:C	2.37	0.43
1:A:3611:PHE:O	1:A:3615:VAL:HG13	2.18	0.43
1:A:3706:TYR:CE2	1:A:3710:ILE:HD12	2.53	0.43
1:A:1654:VAL:O	1:A:1657:THR:HG22	2.18	0.43
1:A:2308:LYS:HE3	1:A:2308:LYS:HA	2.00	0.43
1:A:3655:ARG:NH1	1:A:3681:GLU:OE1	2.51	0.43
1:A:1774:LEU:HD21	1:A:1921:MET:CE	2.48	0.43
1:A:3409:ASP:HB3	1:A:3518:PHE:HB2	2.00	0.43
1:A:3844:ILE:HD11	1:A:3855:LEU:HD22	1.99	0.43
1:A:3895:VAL:CG1	1:A:3896:VAL:N	2.82	0.43
1:A:1443:VAL:O	1:A:1446:VAL:HG12	2.18	0.43
1:A:2735:HIS:ND1	1:A:2738:MET:HB2	2.34	0.43
1:A:2707:VAL:HG12	1:A:2708:ASN:H	1.83	0.43
1:A:3946:VAL:N	1:A:3947:PRO:CD	2.82	0.43
1:A:1572:ILE:HG23	1:A:1574:PHE:CE2	2.54	0.43
1:A:2627:ARG:O	1:A:2631:THR:HG23	2.18	0.43
1:A:2857:ARG:HG2	1:A:2861:ARG:HD3	2.00	0.43
1:A:3156:LEU:HD21	1:A:3159:LYS:HG3	2.00	0.43
1:A:1690:SER:HB3	1:A:1694:VAL:HG13	2.01	0.43
1:A:1929:ILE:HD12	1:A:1929:ILE:H	1.84	0.42
1:A:3319:GLU:O	1:A:3323:ASN:ND2	2.37	0.42
1:A:3986:ARG:HD3	1:A:4012:ALA:HA	2.00	0.42
1:A:2756:MET:O	1:A:2888:VAL:HA	2.19	0.42
1:A:3409:ASP:OD1	1:A:3409:ASP:N	2.52	0.42
1:A:1689:LYS:O	1:A:1689:LYS:HG3	2.18	0.42
1:A:2222:ILE:HG22	1:A:2226:ILE:HD11	2.01	0.42
1:A:2476:LYS:H	1:A:2476:LYS:HD3	1.84	0.42
1:A:3408:LEU:HD23	1:A:3517:HIS:CE1	2.55	0.42
1:A:3671:VAL:O	1:A:3674:ILE:HG22	2.19	0.42
1:A:3851:VAL:HG13	1:A:3855:LEU:HB3	2.00	0.42
1:A:2109:LEU:CD2	1:A:2518:THR:HG22	2.49	0.42
1:A:2225:LYS:HD3	1:A:2284:LEU:HD12	2.01	0.42
1:A:2505:PHE:CE2	1:A:2509:LEU:HD11	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2958:ILE:HG13	1:A:2959:GLN:H	1.85	0.42
1:A:3946:VAL:HG11	1:A:3953:LYS:CB	2.46	0.42
1:A:1394:LEU:HD22	1:A:1449:GLN:OE1	2.19	0.42
1:A:4021:LEU:HD23	1:A:4023:ILE:HG13	2.02	0.42
1:A:2759:ILE:O	1:A:2759:ILE:HG22	2.18	0.42
1:A:3886:ALA:N	1:A:3887:PRO:HD2	2.34	0.42
1:A:2266:PHE:HD1	1:A:2326:LEU:HD21	1.84	0.42
1:A:2387:ARG:O	1:A:2390:ILE:HG22	2.20	0.42
1:A:1660:VAL:HG11	1:A:1728:TRP:CH2	2.55	0.42
1:A:2382:ALA:HB1	1:A:2630:TYR:HE1	1.85	0.41
1:A:2649:ALA:HB1	1:A:2673:LEU:HD21	2.02	0.41
1:A:3632:LEU:HD13	1:A:3644:ILE:HD13	2.01	0.41
1:A:4085:THR:O	1:A:4089:LEU:HG	2.19	0.41
1:A:3743:ASP:OD1	1:A:3743:ASP:N	2.53	0.41
1:A:2749:LEU:HD12	1:A:2773:VAL:HG12	2.03	0.41
1:A:3409:ASP:OD1	1:A:3497:HIS:NE2	2.53	0.41
1:A:1392:LEU:HD23	1:A:1484:LYS:HA	2.02	0.41
1:A:1645:PHE:CB	1:A:1765:ILE:HG22	2.50	0.41
1:A:2225:LYS:NZ	1:A:2285:GLU:OE2	2.35	0.41
1:A:3903:ILE:HD12	1:A:3903:ILE:H	1.85	0.41
1:A:2572:GLU:HG3	1:A:2590:GLU:HA	2.03	0.41
1:A:2661:VAL:HG12	1:A:2916:TRP:CE2	2.56	0.41
1:A:2787:HIS:HB2	1:A:3459:ASP:HB2	2.01	0.41
1:A:1626:CYS:SG	1:A:1639:VAL:HG11	2.61	0.41
1:A:1795:PHE:O	1:A:1795:PHE:CD1	2.74	0.41
1:A:2437:LEU:H	1:A:2437:LEU:HD12	1.85	0.41
1:A:2642:ARG:O	1:A:2645:ILE:HG13	2.21	0.41
1:A:4081:VAL:O	1:A:4085:THR:HG23	2.20	0.41
1:A:4089:LEU:HB3	1:A:4094:SER:OG	2.20	0.41
1:A:3256:ASN:HA	1:A:3259:LYS:HZ2	1.86	0.41
1:A:3538:ASN:HB3	1:A:3541:MET:HG2	2.03	0.41
1:A:2047:PHE:CE2	1:A:2082:ALA:HB1	2.56	0.41
1:A:2491:LEU:N	1:A:2492:PRO:CD	2.84	0.41
1:A:2590:GLU:N	1:A:2591:PRO:HD2	2.35	0.41
1:A:2856:LEU:HD21	1:A:2877:PHE:HB2	2.03	0.41
1:A:2860:THR:HG22	1:A:2865:LEU:O	2.21	0.41
1:A:2765:GLY:CA	1:A:2768:ILE:HG22	2.51	0.41
1:A:3676:TRP:CD1	1:A:3706:TYR:CZ	3.09	0.41
1:A:3833:LYS:HZ2	1:A:3858:HIS:CD2	2.38	0.41
1:A:1926:SER:HB2	1:A:1973:LEU:HD21	2.02	0.40
1:A:2097:HIS:HB2	1:A:2149:ARG:HE	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:4018:SER:O	1:A:4020:ASN:N	2.54	0.40
1:A:1963:MET:HG2	1:A:1965:HIS:CE1	2.57	0.40
1:A:2844:PHE:CG	1:A:2852:LEU:HD22	2.55	0.40
1:A:3727:SER:OG	1:A:3728:GLU:N	2.54	0.40
1:A:1615:ILE:HG13	1:A:1616:LYS:N	2.37	0.40
1:A:1948:LYS:NZ	1:A:1991:GLU:OE1	2.54	0.40
1:A:2716:ILE:HD12	1:A:2739:VAL:HG22	2.02	0.40
1:A:3689:ALA:C	1:A:3690:LEU:HD12	2.42	0.40
1:A:3205:ASN:O	1:A:3206:ALA:HB3	2.21	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	2618/2661 (98%)	2464 (94%)	141 (5%)	13 (0%)	29 68

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1940	GLU
1	A	3810	SER
1	A	4099	HIS
1	A	1391	GLY
1	A	2259	MET
1	A	2763	ARG
1	A	2301	TRP
1	A	2519	PRO
1	A	3947	PRO
1	A	2986	PRO
1	A	3664	THR
1	A	4093	GLY

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Mol	Chain	Res	Type
1	A	2288	VAL

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	2381/2406 (99%)	2309 (97%)	72 (3%)	41 63

All (72) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1383	TYR
1	A	1392	LEU
1	A	1421	TYR
1	A	1463	LEU
1	A	1495	THR
1	A	1602	ILE
1	A	1605	GLN
1	A	1661	GLU
1	A	1671	LYS
1	A	1794	PHE
1	A	1818	VAL
1	A	1880	THR
1	A	1889	SER
1	A	1903	ASN
1	A	1906	SER
1	A	2064	GLN
1	A	2068	GLN
1	A	2081	THR
1	A	2111	LYS
1	A	2155	ASP
1	A	2172	ASP
1	A	2239	ASN
1	A	2257	PHE
1	A	2285	GLU
1	A	2295	ILE

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Mol	Chain	Res	Type
1	A	2302	PHE
1	A	2303	GLN
1	A	2308	LYS
1	A	2346	PHE
1	A	2394	THR
1	A	2439	ASP
1	A	2476	LYS
1	A	2494	LEU
1	A	2640	THR
1	A	2683	ASN
1	A	2721	LYS
1	A	2735	HIS
1	A	2767	THR
1	A	2819	GLU
1	A	2920	TRP
1	A	2932	MET
1	A	2939	GLU
1	A	2941	THR
1	A	2942	ASP
1	A	2999	LEU
1	A	3142	TYR
1	A	3164	ASN
1	A	3299	LEU
1	A	3305	ARG
1	A	3344	ASP
1	A	3538	ASN
1	A	3577	MET
1	A	3578	LEU
1	A	3580	ASN
1	A	3595	MET
1	A	3660	LYS
1	A	3714	GLN
1	A	3745	ARG
1	A	3747	LEU
1	A	3750	TYR
1	A	3871	PHE
1	A	3899	ASP
1	A	3905	ASP
1	A	3906	THR
1	A	3915	PHE
1	A	3939	ILE
1	A	3958	ASP

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Mol	Chain	Res	Type
1	A	3960	ASP
1	A	3975	ASN
1	A	3982	TRP
1	A	3987	ASP
1	A	4033	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1899	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.