Full wwPDB X-ray Structure Validation Report (i)

Dec 2, 2023 – 04:25 pm GMT

PDB ID	:	1UPL
Title	:	Crystal structure of MO25 alpha
Authors	:	Milburn, C.C.; Boudeau, J.; Deak, M.; Alessi, D.R.; Van Aalten, D.M.F.
Deposited on	:	2003-10-07
Resolution	:	2.60 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 (i) 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 (i)) were used in the production of this report:

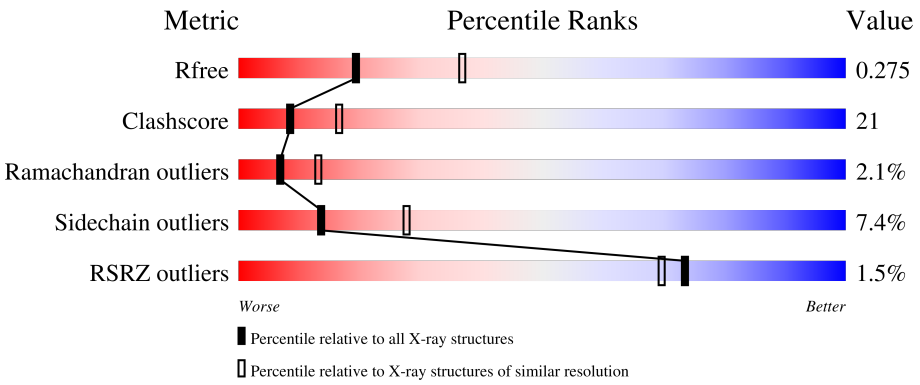
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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 <=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	341	.% 56% 30% 5% 9%
1	B	341	2% 54% 30% 6% • 10%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 5119 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 MO25 PROTEIN.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	312	2567	1650	429	477	3	8	61	0	1
1	B	308	2540	1635	422	472	3	8	57	0	1

- Molecule 2 is water.

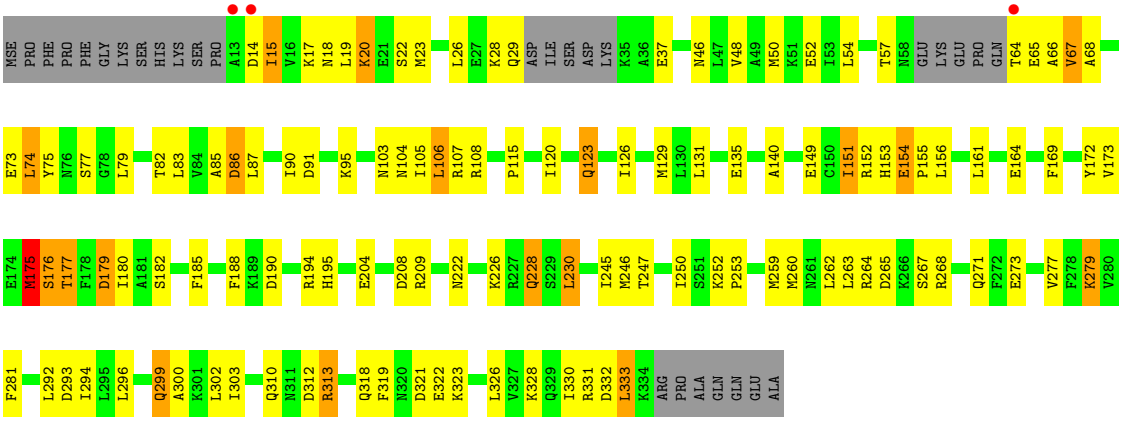
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	8	8	8	0	0
2	B	4	4	4	0	0

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.

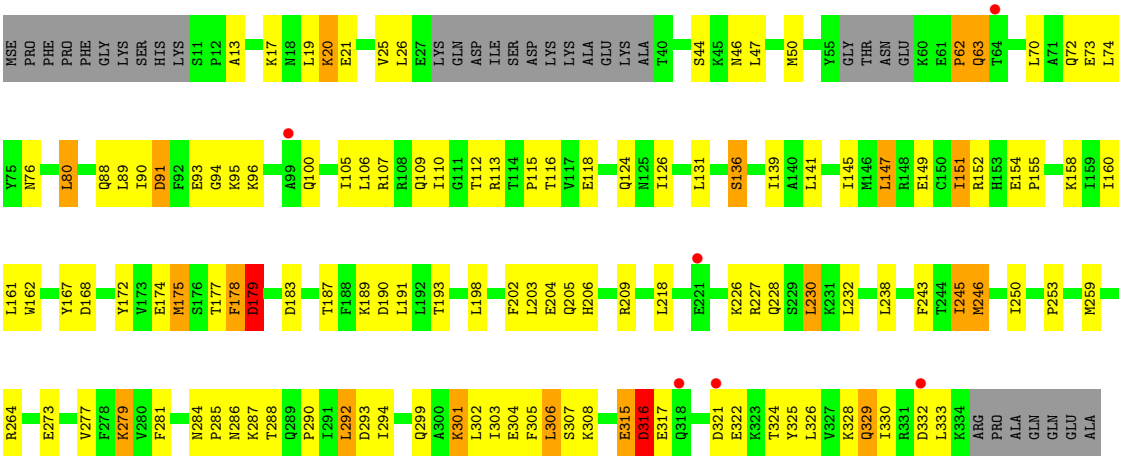
- Molecule 1: MO25 PROTEIN

Chain A:



- Molecule 1: MO25 PROTEIN

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	70.62Å 94.01Å 96.33Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 2.60 24.82 – 2.60	Depositor EDS
% Data completeness (in resolution range)	98.6 (25.00-2.60) 98.5 (24.82-2.60)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.95 (at 2.60Å)	Xtriage
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.212 , 0.280 0.210 , 0.275	Depositor DCC
R_{free} test set	524 reflections (2.62%)	wwPDB-VP
Wilson B-factor (Å^2)	40.7	Xtriage
Anisotropy	0.553	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 44.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.026 for -h,l,k	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	5119	wwPDB-VP
Average B, all atoms (Å^2)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.10% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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	0.62	1/2601 (0.0%)	0.96	10/3485 (0.3%)
1	B	0.64	2/2576 (0.1%)	0.96	14/3455 (0.4%)
All	All	0.63	3/5177 (0.1%)	0.96	24/6940 (0.3%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	264	ARG	NE-CZ	12.42	1.49	1.33
1	B	315	GLU	CB-CG	7.12	1.65	1.52
1	A	108	ARG	CG-CD	6.94	1.69	1.51

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	264	ARG	CD-NE-CZ	-10.31	109.17	123.60
1	B	264	ARG	NE-CZ-NH2	-10.07	115.27	120.30
1	A	293	ASP	CB-CG-OD2	8.24	125.72	118.30
1	B	264	ARG	NE-CZ-NH1	7.43	124.02	120.30
1	A	86	ASP	CB-CG-OD2	7.10	124.69	118.30
1	B	332	ASP	CB-CG-OD2	6.53	124.17	118.30
1	B	333	LEU	O-C-N	-6.45	112.39	122.70
1	A	312	ASP	CB-CG-OD2	6.34	124.01	118.30
1	A	208	ASP	CB-CG-OD2	6.19	123.87	118.30
1	B	293	ASP	CB-CG-OD2	6.17	123.86	118.30
1	B	321	ASP	CB-CG-OD2	6.01	123.71	118.30
1	A	14	ASP	CB-CG-OD2	5.80	123.52	118.30
1	B	316	ASP	CB-CG-OD2	5.76	123.49	118.30
1	A	179	ASP	CB-CG-OD2	5.74	123.47	118.30
1	B	190	ASP	CB-CG-OD2	5.70	123.43	118.30
1	A	333	LEU	O-C-N	-5.64	113.67	122.70
1	A	190	ASP	CB-CG-OD2	5.60	123.34	118.30
1	B	183	ASP	CB-CG-OD2	5.57	123.31	118.30
1	A	91	ASP	CB-CG-OD2	5.46	123.22	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	168	ASP	CB-CG-OD1	5.25	123.03	118.30
1	B	91	ASP	CB-CG-OD2	5.19	122.97	118.30
1	B	161	LEU	CB-CG-CD2	-5.17	102.22	111.00
1	A	106	LEU	CB-CG-CD2	-5.09	102.34	111.00
1	B	179	ASP	CB-CG-OD2	5.04	122.83	118.30

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	2567	0	2606	97	2
1	B	2540	0	2573	112	2
2	A	8	0	0	0	0
2	B	4	0	0	1	0
All	All	5119	0	5179	207	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 21.

All (207) 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:23:MSE:HE1	1:A:83:LEU:CD2	1.55	1.36
1:B:301:LYS:HD3	1:B:301:LYS:N	1.49	1.25
1:B:301:LYS:H	1:B:301:LYS:CD	1.35	1.25
1:B:107:ARG:NH2	1:B:152:ARG:HD3	1.52	1.24
1:A:15:ILE:HD12	1:A:15:ILE:H	1.07	1.07
1:A:23:MSE:HE1	1:A:83:LEU:HD21	1.14	1.07
1:B:301:LYS:N	1:B:301:LYS:CD	2.04	1.06
1:B:80:LEU:HD12	1:B:80:LEU:H	1.18	1.02
1:A:20:LYS:HE3	1:A:77:SER:O	1.60	1.01
1:A:23:MSE:CE	1:A:83:LEU:HD21	1.91	1.01
1:B:107:ARG:HH21	1:B:152:ARG:HD3	1.23	0.98

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:19:LEU:HD21	1:B:47:LEU:HD23	1.50	0.94
1:A:23:MSE:HE1	1:A:83:LEU:HD23	1.47	0.94
1:A:23:MSE:CE	1:A:83:LEU:CD2	2.46	0.92
1:A:15:ILE:H	1:A:15:ILE:CD1	1.82	0.90
1:B:90:ILE:HG22	1:B:94:GLY:HA3	1.54	0.89
1:A:19:LEU:HD13	1:A:46:ASN:HB2	1.55	0.88
1:B:107:ARG:NH2	1:B:152:ARG:HH11	1.74	0.85
1:B:315:GLU:O	1:B:316:ASP:HB2	1.77	0.85
1:B:90:ILE:CG2	1:B:94:GLY:HA3	2.06	0.84
1:B:107:ARG:NH2	1:B:152:ARG:CD	2.39	0.84
1:A:54:LEU:HD22	1:A:104:ASN:OD1	1.76	0.84
1:B:19:LEU:HD13	1:B:46:ASN:HB3	1.60	0.84
1:A:175:MSE:HG3	1:A:176:SER:N	1.93	0.82
1:A:15:ILE:HD12	1:A:15:ILE:N	1.92	0.82
1:A:151:ILE:N	1:A:151:ILE:HD13	1.94	0.81
1:B:93:GLU:OE1	1:B:96:LYS:HE2	1.82	0.80
1:B:227:ARG:NH2	1:B:228:GLN:OE1	2.15	0.79
1:A:106:LEU:HD23	1:A:149:GLU:HG3	1.64	0.79
1:B:301:LYS:HD3	1:B:301:LYS:H	0.64	0.79
1:A:19:LEU:HD13	1:A:46:ASN:CB	2.13	0.79
1:B:93:GLU:OE1	1:B:96:LYS:CE	2.32	0.77
1:A:140:ALA:CB	1:A:180:ILE:HD12	2.17	0.75
1:B:301:LYS:N	1:B:301:LYS:HD2	2.00	0.74
1:B:80:LEU:H	1:B:80:LEU:CD1	1.89	0.74
1:A:123:GLN:HB3	1:A:126:ILE:HD12	1.71	0.72
1:A:54:LEU:HD11	1:A:105:ILE:HD11	1.71	0.71
1:B:174:GLU:O	1:B:175:MSE:C	2.29	0.71
1:A:319:PHE:CE2	1:A:323:LYS:HE2	2.27	0.70
1:A:54:LEU:HD11	1:A:105:ILE:CD1	2.22	0.70
1:A:172:TYR:O	1:A:175:MSE:CB	2.40	0.69
1:A:140:ALA:HB3	1:A:180:ILE:HD12	1.75	0.68
1:B:106:LEU:HD13	1:B:116:THR:CG2	2.24	0.68
1:A:172:TYR:O	1:A:175:MSE:HB2	1.94	0.68
1:A:322:GLU:O	1:A:326:LEU:HG	1.93	0.67
1:B:105:ILE:CG2	1:B:116:THR:OG1	2.42	0.67
1:B:107:ARG:NH2	1:B:152:ARG:NH1	2.42	0.67
1:B:19:LEU:HD13	1:B:46:ASN:CB	2.25	0.66
1:B:287:LYS:HD2	1:B:292:LEU:HD12	1.76	0.66
1:B:26:LEU:HB3	1:B:89:LEU:HD13	1.77	0.66
1:B:62:PRO:O	1:B:63:GLN:HB2	1.95	0.66
1:B:50:MSE:HG3	1:B:70:LEU:HD21	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:23:MSE:CE	1:A:83:LEU:HD23	2.22	0.64
1:A:64:THR:HG23	1:A:64:THR:O	1.97	0.64
1:B:26:LEU:HB3	1:B:89:LEU:CD1	2.27	0.64
1:B:107:ARG:HH21	1:B:152:ARG:CD	2.04	0.64
1:B:162:TRP:HZ2	1:B:205:GLN:NE2	1.97	0.63
1:B:110:ILE:HG22	1:B:110:ILE:O	1.97	0.63
1:A:67:VAL:HG13	1:B:286:ASN:ND2	2.13	0.62
1:A:279:LYS:HE2	1:A:322:GLU:OE2	2.00	0.62
1:B:105:ILE:HG23	1:B:116:THR:OG1	1.99	0.62
1:A:19:LEU:CD1	1:A:46:ASN:HB2	2.30	0.62
1:B:177:THR:O	1:B:179:ASP:N	2.34	0.61
1:A:300:ALA:HA	1:A:303:ILE:HD12	1.84	0.60
1:B:250:ILE:HA	1:B:281:PHE:CE2	2.37	0.60
1:B:107:ARG:HH22	1:B:152:ARG:HH11	1.50	0.60
1:A:230:LEU:HB3	1:A:273:GLU:HB3	1.84	0.59
1:B:284:ASN:O	1:B:287:LYS:NZ	2.36	0.59
1:A:265:ASP:O	1:A:271:GLN:NE2	2.36	0.58
1:B:218:LEU:HD22	1:B:230:LEU:HD13	1.85	0.58
1:A:67:VAL:O	1:A:67:VAL:HG23	2.03	0.58
1:A:318:GLN:O	1:A:322:GLU:HG3	2.04	0.58
1:B:204:GLU:HA	1:B:245:ILE:CD1	2.34	0.58
1:A:279:LYS:HE2	1:A:322:GLU:CD	2.22	0.58
1:A:19:LEU:HD22	1:A:50:MSE:CE	2.33	0.58
1:B:243:PHE:HA	2:B:2004:HOH:O	2.03	0.58
1:B:259:MSE:HG3	1:B:277:VAL:HG12	1.85	0.58
1:B:88:GLN:HG3	1:B:139:ILE:HD13	1.85	0.57
1:A:82:THR:O	1:A:86:ASP:HB2	2.03	0.57
1:B:158:LYS:HG3	1:B:198:LEU:HD11	1.86	0.57
1:A:259:MSE:SE	1:A:277:VAL:HG11	2.54	0.57
1:B:80:LEU:HD12	1:B:80:LEU:N	2.03	0.57
1:B:203:LEU:HB3	1:B:245:ILE:HG12	1.87	0.56
1:B:179:ASP:N	1:B:179:ASP:OD1	2.39	0.56
1:B:246:MSE:HE3	1:B:250:ILE:CD1	2.36	0.56
1:A:154:GLU:N	1:A:155:PRO:CD	2.68	0.56
1:A:172:TYR:O	1:A:175:MSE:N	2.39	0.56
1:B:106:LEU:HD13	1:B:116:THR:HG22	1.87	0.55
1:A:75:TYR:OH	1:A:115:PRO:HB2	2.07	0.55
1:A:228:GLN:HA	1:A:228:GLN:OE1	2.06	0.54
1:B:299:GLN:O	1:B:303:ILE:HG13	2.07	0.54
1:B:19:LEU:CD1	1:B:46:ASN:CB	2.85	0.54
1:A:67:VAL:HG13	1:B:286:ASN:HD21	1.71	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:189:LYS:HG3	1:B:232:LEU:HD13	1.89	0.54
1:B:284:ASN:OD1	1:B:285:PRO:HD2	2.07	0.54
1:B:145:ILE:O	1:B:149:GLU:HG2	2.08	0.54
1:B:288:THR:OG1	1:B:290:PRO:HD2	2.08	0.53
1:B:304:GLU:OE2	1:B:308:LYS:HD3	2.09	0.53
1:B:279:LYS:HE2	1:B:322:GLU:OE2	2.08	0.53
1:B:279:LYS:NZ	1:B:322:GLU:OE2	2.39	0.52
1:B:19:LEU:HD21	1:B:47:LEU:CD2	2.30	0.52
1:B:13:ALA:HA	1:B:73:GLU:HG3	1.90	0.52
1:A:292:LEU:HD11	1:A:333:LEU:HD22	1.92	0.52
1:B:95:LYS:HE3	1:B:139:ILE:HG22	1.91	0.51
1:A:226:LYS:HE3	1:A:262:LEU:HD21	1.93	0.51
1:A:65:GLU:HG3	1:A:67:VAL:HG22	1.92	0.51
1:A:250:ILE:HA	1:A:281:PHE:CE2	2.46	0.51
1:A:48:VAL:O	1:A:52:GLU:HG3	2.11	0.50
1:A:103:ASN:O	1:A:107:ARG:HG3	2.12	0.50
1:B:19:LEU:CD1	1:B:46:ASN:HB2	2.41	0.50
1:B:74:LEU:O	1:B:80:LEU:HD11	2.10	0.50
1:B:50:MSE:CG	1:B:70:LEU:HD21	2.41	0.50
1:B:96:LYS:O	1:B:100:GLN:HG3	2.12	0.50
1:A:319:PHE:CZ	1:A:323:LYS:HE2	2.47	0.50
1:B:151:ILE:CD1	1:B:191:LEU:HD21	2.42	0.50
1:A:185:PHE:CZ	1:A:228:GLN:HB3	2.46	0.49
1:B:302:LEU:O	1:B:306:LEU:HG	2.12	0.49
1:A:140:ALA:HB3	1:A:180:ILE:CD1	2.42	0.49
1:A:140:ALA:HB1	1:A:180:ILE:HD12	1.93	0.49
1:B:21:GLU:O	1:B:25:VAL:HG23	2.12	0.49
1:A:185:PHE:CE1	1:A:228:GLN:HB3	2.48	0.48
1:B:167:TYR:CZ	1:B:209:ARG:HG2	2.48	0.48
1:A:85:ALA:HB2	1:A:129:MSE:SE	2.63	0.48
1:B:17:LYS:O	1:B:21:GLU:HG3	2.14	0.48
1:A:86:ASP:O	1:A:87:LEU:C	2.51	0.48
1:A:310:GLN:OE1	1:A:313:ARG:NE	2.42	0.48
1:B:324:THR:O	1:B:325:TYR:C	2.52	0.48
1:B:136:SER:HB2	1:B:139:ILE:HG12	1.96	0.47
1:B:107:ARG:HH22	1:B:152:ARG:HD3	1.67	0.47
1:A:328:LYS:NZ	1:A:332:ASP:OD2	2.40	0.47
1:B:204:GLU:HA	1:B:245:ILE:HD11	1.97	0.47
1:B:230:LEU:HB3	1:B:273:GLU:HB3	1.97	0.47
1:A:28:LYS:O	1:A:29:GLN:HG3	2.15	0.47
1:A:74:LEU:HD11	1:A:105:ILE:HG21	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:20:LYS:O	1:B:20:LYS:HD2	2.15	0.47
1:A:252:LYS:O	1:A:253:PRO:C	2.54	0.46
1:A:23:MSE:SE	1:A:83:LEU:HD23	2.65	0.46
1:A:164:GLU:CD	1:A:209:ARG:HH22	2.18	0.46
1:A:253:PRO:HB3	1:A:294:ILE:HD11	1.97	0.46
1:B:72:GLN:NE2	1:B:76:ASN:OD1	2.48	0.46
1:B:189:LYS:O	1:B:193:THR:OG1	2.25	0.46
1:A:26:LEU:HA	1:A:26:LEU:HD23	1.61	0.46
1:A:65:GLU:CG	1:A:67:VAL:HG22	2.46	0.46
1:B:151:ILE:HD12	1:B:151:ILE:HA	1.75	0.46
1:B:154:GLU:HB3	1:B:155:PRO:HD3	1.98	0.46
1:B:279:LYS:CE	1:B:322:GLU:OE2	2.63	0.46
1:A:149:GLU:OE2	1:A:152:ARG:HD3	2.16	0.46
1:A:328:LYS:O	1:A:332:ASP:N	2.48	0.45
1:B:91:ASP:OD1	1:B:93:GLU:HB2	2.16	0.45
1:A:259:MSE:SE	1:A:277:VAL:CG1	3.15	0.45
1:A:296:LEU:HD21	1:A:333:LEU:HB3	1.99	0.45
1:A:151:ILE:HG22	1:A:151:ILE:O	2.17	0.45
1:A:19:LEU:HD22	1:A:50:MSE:HE1	1.97	0.45
1:A:299:GLN:O	1:A:300:ALA:C	2.55	0.45
1:B:107:ARG:HH21	1:B:152:ARG:NH1	2.12	0.45
1:B:115:PRO:O	1:B:118:GLU:HB2	2.16	0.45
1:B:218:LEU:CD2	1:B:230:LEU:HD13	2.47	0.45
1:B:93:GLU:OE1	1:B:96:LYS:HE3	2.13	0.44
1:A:302:LEU:HD23	1:A:302:LEU:HA	1.83	0.44
1:A:169:PHE:O	1:A:173:VAL:HG22	2.18	0.44
1:B:202:PHE:CD1	1:B:206:HIS:ND1	2.86	0.44
1:A:64:THR:O	1:A:64:THR:CG2	2.64	0.44
1:A:204:GLU:HG3	1:A:245:ILE:HD11	1.99	0.43
1:B:238:LEU:HD23	1:B:238:LEU:HA	1.62	0.43
1:A:19:LEU:HD13	1:A:46:ASN:HB3	1.97	0.43
1:B:253:PRO:HB3	1:B:294:ILE:HD11	2.00	0.43
1:A:87:LEU:HD12	1:A:87:LEU:HA	1.86	0.43
1:A:296:LEU:HD23	1:A:296:LEU:HA	1.74	0.43
1:B:162:TRP:CZ2	1:B:205:GLN:NE2	2.81	0.43
1:B:172:TYR:O	1:B:175:MSE:HB2	2.18	0.43
1:B:259:MSE:SE	1:B:277:VAL:HG11	2.69	0.43
1:A:259:MSE:HG3	1:A:277:VAL:HG12	2.01	0.43
1:B:62:PRO:O	1:B:63:GLN:CB	2.65	0.43
1:B:203:LEU:O	1:B:245:ILE:HD11	2.19	0.43
1:B:325:TYR:CE2	1:B:329:GLN:HG3	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:218:LEU:O	1:B:226:LYS:HE3	2.18	0.43
1:A:90:ILE:O	1:A:95:LYS:HE3	2.19	0.43
1:A:151:ILE:O	1:A:195:HIS:HE1	2.02	0.43
1:A:19:LEU:HD23	1:A:79:LEU:HD21	2.01	0.42
1:A:120:ILE:HG22	1:A:156:LEU:HD22	2.01	0.42
1:B:305:PHE:O	1:B:306:LEU:C	2.57	0.42
1:B:246:MSE:HE3	1:B:250:ILE:HD13	2.01	0.42
1:A:176:SER:O	1:A:177:THR:C	2.57	0.42
1:B:326:LEU:O	1:B:330:ILE:HG12	2.20	0.42
1:A:246:MSE:O	1:A:247:THR:C	2.56	0.42
1:A:153:HIS:C	1:A:155:PRO:HD2	2.40	0.42
1:A:135:GLU:HG3	1:A:172:TYR:CE1	2.55	0.42
1:A:172:TYR:O	1:A:175:MSE:CA	2.67	0.41
1:B:154:GLU:N	1:B:155:PRO:CD	2.82	0.41
1:A:151:ILE:HD13	1:A:151:ILE:H	1.82	0.41
1:A:260:MSE:O	1:A:263:LEU:HB2	2.20	0.41
1:B:72:GLN:O	1:B:76:ASN:HB2	2.21	0.41
1:A:18:ASN:O	1:A:22:SER:HB2	2.21	0.41
1:B:20:LYS:HD3	1:B:20:LYS:HA	1.87	0.41
1:A:279:LYS:HB2	1:A:326:LEU:HD11	2.02	0.41
1:B:141:LEU:O	1:B:145:ILE:HD12	2.21	0.41
1:B:284:ASN:HA	1:B:285:PRO:HD3	1.87	0.41
1:B:328:LYS:O	1:B:330:ILE:N	2.53	0.41
1:A:185:PHE:O	1:A:188:PHE:HB3	2.21	0.41
1:B:124:GLN:C	1:B:126:ILE:N	2.73	0.40
1:B:330:ILE:O	1:B:330:ILE:HG22	2.21	0.40
1:A:66:ALA:C	1:A:68:ALA:H	2.24	0.40
1:B:19:LEU:CD1	1:B:46:ASN:HB3	2.39	0.40
1:B:147:LEU:HD13	1:B:187:THR:HG21	2.02	0.40
1:B:306:LEU:O	1:B:307:SER:C	2.59	0.40
1:A:222:ASN:OD1	1:A:222:ASN:C	2.60	0.40
1:B:74:LEU:HD23	1:B:74:LEU:HA	1.91	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:321:ASP:OD2	1:B:152:ARG:NH2[4_556]	1.88	0.32
1:A:194:ARG:NH1	1:B:317:GLU:OE2[4_556]	2.12	0.08

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	306/341 (90%)	283 (92%)	17 (6%)	6 (2%)	7 14
1	B	302/341 (89%)	275 (91%)	20 (7%)	7 (2%)	6 11
All	All	608/682 (89%)	558 (92%)	37 (6%)	13 (2%)	7 13

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	175	MSE
1	B	178	PHE
1	B	316	ASP
1	A	67	VAL
1	A	177	THR
1	A	299	GLN
1	B	63	GLN
1	A	57	THR
1	B	306	LEU
1	A	175	MSE
1	A	176	SER
1	B	329	GLN
1	B	62	PRO

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	285/303 (94%)	262 (92%)	23 (8%)	11 23

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	284/303 (94%)	265 (93%)	19 (7%)	16	33
All	All	569/606 (94%)	527 (93%)	42 (7%)	13	28

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	ILE
1	A	17	LYS
1	A	20	LYS
1	A	37	GLU
1	A	73	GLU
1	A	74	LEU
1	A	123	GLN
1	A	131	LEU
1	A	151	ILE
1	A	154	GLU
1	A	161	LEU
1	A	175	MSE
1	A	179	ASP
1	A	182	SER
1	A	228	GLN
1	A	230	LEU
1	A	264	ARG
1	A	267	SER
1	A	268	ARG
1	A	279	LYS
1	A	313	ARG
1	A	330	ILE
1	A	331	ARG
1	B	20	LYS
1	B	44	SER
1	B	80	LEU
1	B	109	GLN
1	B	112	THR
1	B	113	ARG
1	B	131	LEU
1	B	136	SER
1	B	147	LEU
1	B	151	ILE
1	B	160	ILE
1	B	178	PHE
1	B	179	ASP

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Mol	Chain	Res	Type
1	B	230	LEU
1	B	245	ILE
1	B	246	MSE
1	B	279	LYS
1	B	292	LEU
1	B	301	LYS

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

Mol	Chain	Res	Type
1	A	76	ASN
1	A	153	HIS
1	A	195	HIS
1	A	205	GLN
1	A	261	ASN
1	A	320	ASN
1	B	46	ASN
1	B	125	ASN
1	B	205	GLN
1	B	261	ASN
1	B	286	ASN
1	B	298	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 monosaccharides 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	304/341 (89%)	-0.40	3 (0%)	82	80	4, 16, 36, 51	16 (5%)
1	B	300/341 (87%)	-0.27	6 (2%)	65	60	4, 18, 41, 47	15 (5%)
All	All	604/682 (88%)	-0.33	9 (1%)	73	70	4, 17, 40, 51	31 (5%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	14	ASP	4.0
1	B	64	THR	3.9
1	B	321	ASP	2.5
1	B	332	ASP	2.3
1	A	64	THR	2.2
1	A	13	ALA	2.1
1	B	221	GLU	2.1
1	B	99	ALA	2.0
1	B	318	GLN	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 monosaccharides 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.