Full wwPDB X-ray Structure Validation Report (i)

Jun 17, 2024 – 06:35 AM EDT

PDB ID	:	3KTI
Title	:	Structure of ClpP in complex with ADEP1
Authors	:	Lee, B.-G.; Brotz-Oesterhelt, H.; Song, H.K.
Deposited on	:	2009-11-25
Resolution	:	2.00 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 (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:

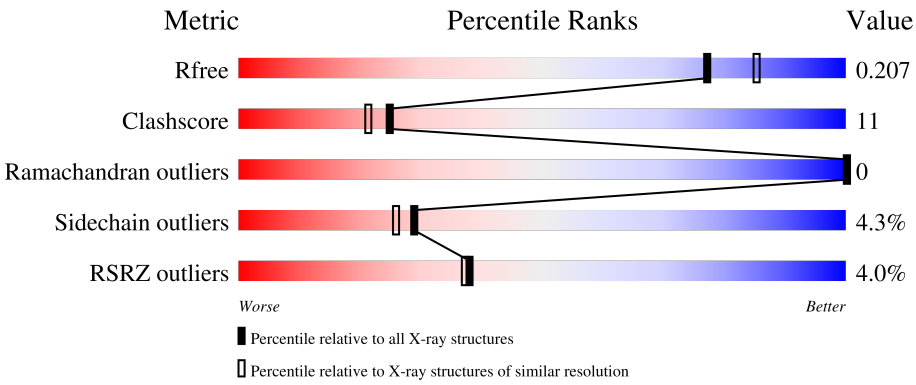
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

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 2.00 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}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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	199	2%, 66%, 20%, •, 13%
1	B	199	4%, 69%, 18%, •, 13%
1	C	199	6%, 67%, 19%, •, 13%
1	D	199	5%, 71%, 14%, •, 13%
1	E	199	3%, 64%, 23%, •, 13%

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Mol	Chain	Length	Quality of chain
1	F	199	3% 66% 20% • 13%
1	G	199	2% 67% 19% • 13%
2	H	7	29% 71%
2	I	7	29% 71%
2	J	7	14% 29% 71%
2	K	7	14% 29% 71%
2	L	7	14% 86%
2	M	7	29% 71%
2	N	7	14% 14% 71% 14%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	NHE	E	505	-	X	-	-
4	NHE	F	506	-	X	-	-
4	NHE	G	507	-	X	-	-

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 10432 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 ATP-dependent Clp protease proteolytic subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	174	1326	839	223	257	7	0	0	0
1	B	174	1326	839	223	257	7	0	0	0
1	C	174	1325	838	223	257	7	0	0	0
1	D	174	1325	838	223	257	7	0	0	0
1	E	174	1330	841	225	257	7	0	0	0
1	F	174	1325	838	223	257	7	0	0	0
1	G	174	1325	838	223	257	7	0	0	0

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	197	HIS	-	EXPRESSION TAG	UNP P80244
A	198	HIS	-	EXPRESSION TAG	UNP P80244
A	199	HIS	-	EXPRESSION TAG	UNP P80244
B	197	HIS	-	EXPRESSION TAG	UNP P80244
B	198	HIS	-	EXPRESSION TAG	UNP P80244
B	199	HIS	-	EXPRESSION TAG	UNP P80244
C	197	HIS	-	EXPRESSION TAG	UNP P80244
C	198	HIS	-	EXPRESSION TAG	UNP P80244
C	199	HIS	-	EXPRESSION TAG	UNP P80244
D	197	HIS	-	EXPRESSION TAG	UNP P80244
D	198	HIS	-	EXPRESSION TAG	UNP P80244
D	199	HIS	-	EXPRESSION TAG	UNP P80244
E	197	HIS	-	EXPRESSION TAG	UNP P80244
E	198	HIS	-	EXPRESSION TAG	UNP P80244
E	199	HIS	-	EXPRESSION TAG	UNP P80244

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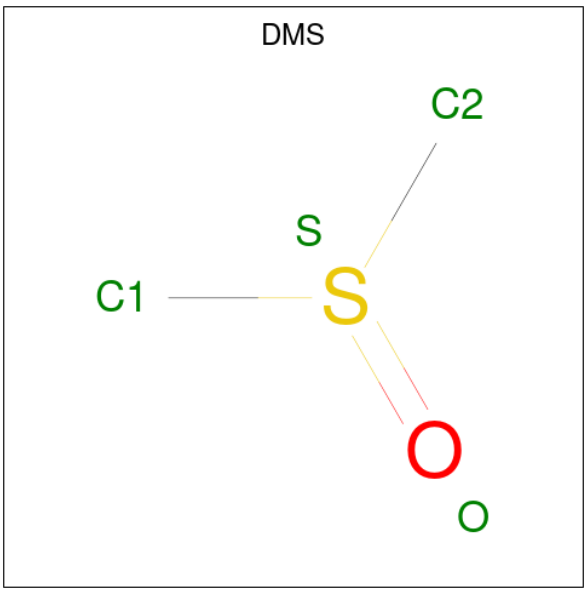
Chain	Residue	Modelled	Actual	Comment	Reference
F	197	HIS	-	EXPRESSION TAG	UNP P80244
F	198	HIS	-	EXPRESSION TAG	UNP P80244
F	199	HIS	-	EXPRESSION TAG	UNP P80244
G	197	HIS	-	EXPRESSION TAG	UNP P80244
G	198	HIS	-	EXPRESSION TAG	UNP P80244
G	199	HIS	-	EXPRESSION TAG	UNP P80244

- Molecule 2 is a protein called Acyldepsipeptide 1.

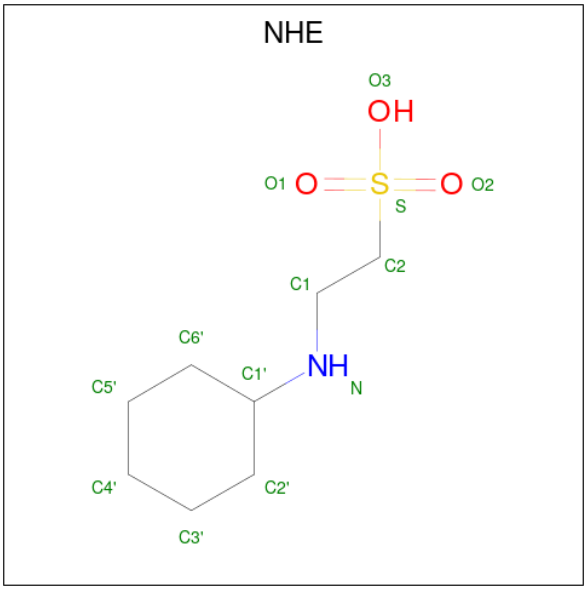
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	H	7	52	38	6	8	0	0	0
2	I	7	52	38	6	8	0	0	0
2	J	7	52	38	6	8	0	0	0
2	K	7	52	38	6	8	0	0	0
2	L	7	52	38	6	8	0	0	0
2	M	7	52	38	6	8	0	0	0
2	N	7	52	38	6	8	0	0	0

- Molecule 3 is DIMETHYL SULFOXIDE (three-letter code: DMS) (formula: C_2H_6OS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O	S		
3	A	1	4	2	1	1	0	0
3	B	1	4	2	1	1	0	0
3	C	1	4	2	1	1	0	0
3	D	1	4	2	1	1	0	0
3	E	1	4	2	1	1	0	0
3	F	1	4	2	1	1	0	0
3	G	1	4	2	1	1	0	0

- Molecule 4 is 2-[N-CYCLOHEXYLAMINO]ETHANE SULFONIC ACID (three-letter code: NHE) (formula: $C_8H_{17}NO_3S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
4	A	1	13	8	1	3	1	0	0
4	B	1	13	8	1	3	1	0	0
4	C	1	13	8	1	3	1	0	0
4	D	1	13	8	1	3	1	0	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
4	E	1	13	8	1	3	1	0	0
			Total	C	N	O	S		
4	F	1	13	8	1	3	1	0	0
			Total	C	N	O	S		
4	G	1	13	8	1	3	1	0	0

- Molecule 5 is water.

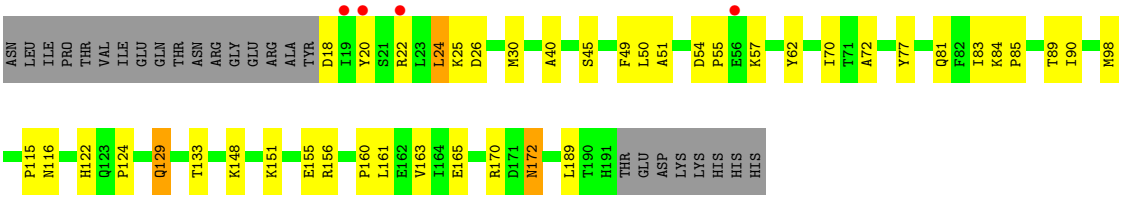
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	111	111	111	0	0
			Total	O		
5	B	117	117	117	0	0
			Total	O		
5	C	69	69	69	0	0
			Total	O		
5	D	105	105	105	0	0
			Total	O		
5	E	78	78	78	0	0
			Total	O		
5	F	109	109	109	0	0
			Total	O		
5	G	75	75	75	0	0
			Total	O		
5	J	1	1	1	0	0
			Total	O		
5	N	2	2	2	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.

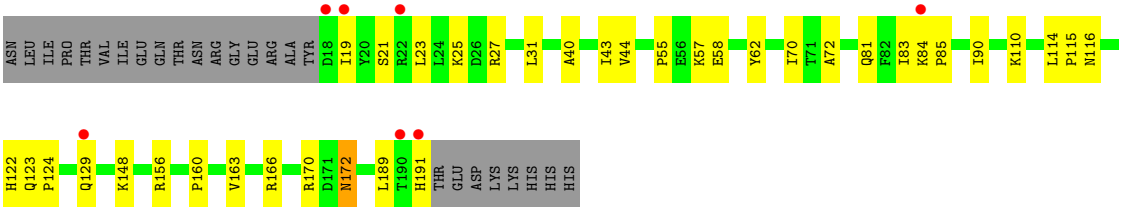
• Molecule 1: ATP-dependent Clp protease proteolytic subunit

Chain A:



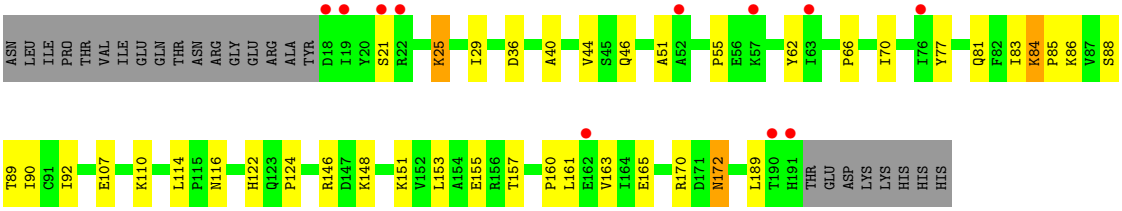
• Molecule 1: ATP-dependent Clp protease proteolytic subunit

Chain B:



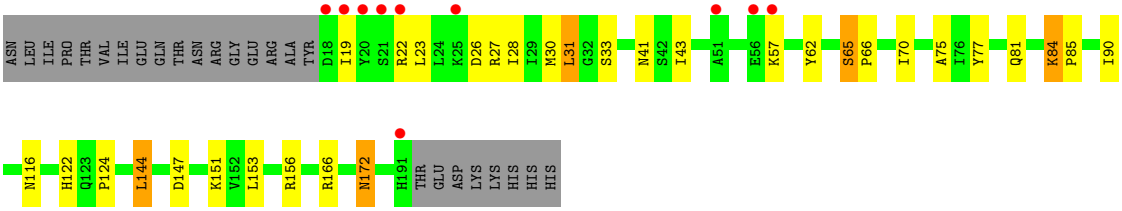
• Molecule 1: ATP-dependent Clp protease proteolytic subunit

Chain C:



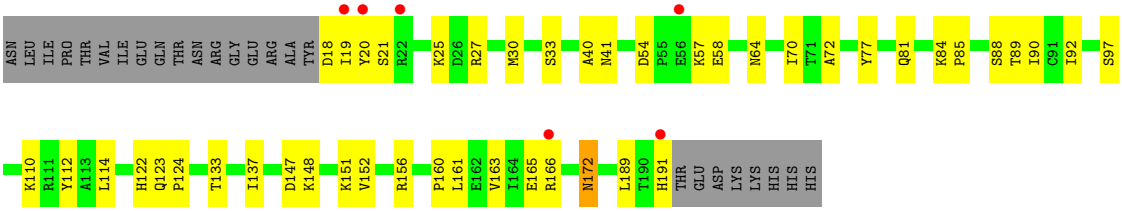
• Molecule 1: ATP-dependent Clp protease proteolytic subunit

Chain D:



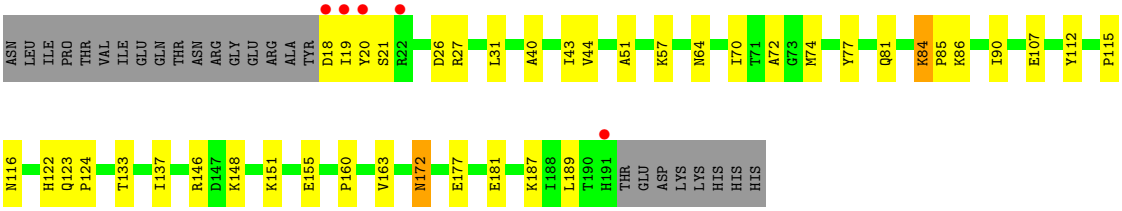
- Molecule 1: ATP-dependent Clp protease proteolytic subunit

Chain E:



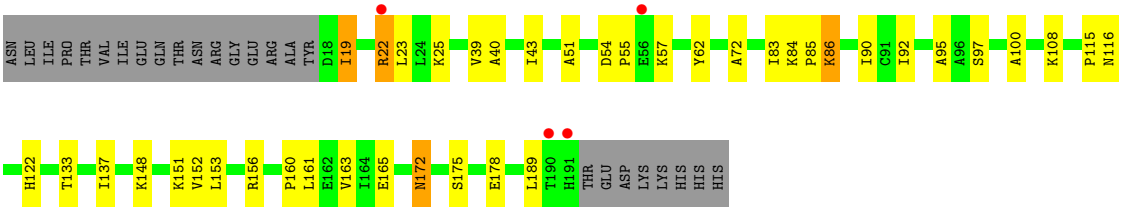
- Molecule 1: ATP-dependent Clp protease proteolytic subunit

Chain F:



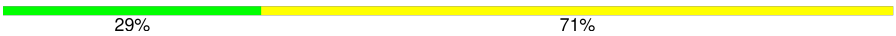
- Molecule 1: ATP-dependent Clp protease proteolytic subunit

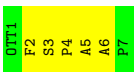
Chain G:



- Molecule 2: Acyldepsipeptide 1

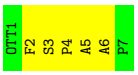
Chain H:





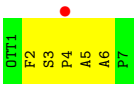
- Molecule 2: Acyldepsipeptide 1

Chain I:



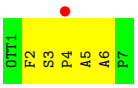
• Molecule 2: Acyldepsipeptide 1

Chain J:



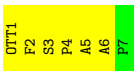
• Molecule 2: Acyldepsipeptide 1

Chain K:



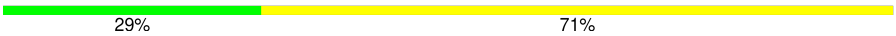
• Molecule 2: Acyldepsipeptide 1

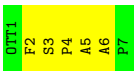
Chain L:



• Molecule 2: Acyldepsipeptide 1

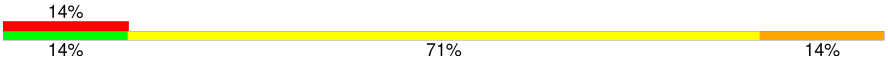
Chain M:

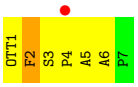




• Molecule 2: Acyldepsipeptide 1

Chain N:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	121.64Å 151.76Å 100.67Å 90.00° 120.03° 90.00°	Depositor
Resolution (Å)	50.00 – 2.00 37.79 – 2.00	Depositor EDS
% Data completeness (in resolution range)	90.9 (50.00-2.00) 98.4 (37.79-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.30 (at 2.00Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.217 , 0.253 0.214 , 0.207	Depositor DCC
R_{free} test set	5210 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å^2)	31.6	Xtriage
Anisotropy	0.424	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 50.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	10432	wwPDB-VP
Average B, all atoms (Å^2)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.84% 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)

Bond lengths and bond angles in the following residue types are not validated in this section: MP8, OTT, NHE, DMS, MAA

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	0.35	0/1341	0.58	0/1807
1	B	0.35	0/1341	0.59	0/1807
1	C	0.30	0/1340	0.54	0/1806
1	D	0.32	0/1340	0.58	0/1806
1	E	0.31	0/1346	0.56	0/1814
1	F	0.33	0/1340	0.59	0/1806
1	G	0.31	0/1340	0.55	0/1806
2	H	3.70	8/29 (27.6%)	2.16	1/37 (2.7%)
2	I	3.68	8/29 (27.6%)	2.23	1/37 (2.7%)
2	J	3.67	8/29 (27.6%)	2.20	1/37 (2.7%)
2	K	3.74	8/29 (27.6%)	2.22	1/37 (2.7%)
2	L	3.77	8/29 (27.6%)	2.16	1/37 (2.7%)
2	M	3.70	8/29 (27.6%)	2.19	1/37 (2.7%)
2	N	3.71	8/29 (27.6%)	2.21	1/37 (2.7%)
All	All	0.63	56/9591 (0.6%)	0.64	7/12911 (0.1%)

All (56) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	L	2	PHE	CG-CD2	8.11	1.50	1.38
2	K	2	PHE	CG-CD2	8.03	1.50	1.38
2	N	2	PHE	CG-CD2	7.95	1.50	1.38
2	M	2	PHE	CG-CD2	7.79	1.50	1.38
2	H	2	PHE	CG-CD2	7.72	1.50	1.38
2	J	2	PHE	CG-CD2	7.67	1.50	1.38
2	I	2	PHE	CG-CD2	7.61	1.50	1.38
2	K	2	PHE	CG-CD1	7.35	1.49	1.38
2	I	2	PHE	CG-CD1	7.25	1.49	1.38
2	J	2	PHE	CG-CD1	7.20	1.49	1.38
2	H	2	PHE	CG-CD1	7.18	1.49	1.38
2	L	2	PHE	CG-CD1	7.18	1.49	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	M	2	PHE	CG-CD1	7.13	1.49	1.38
2	N	2	PHE	CG-CD1	7.11	1.49	1.38
2	J	2	PHE	CE1-CZ	6.58	1.49	1.37
2	N	2	PHE	CE1-CZ	6.56	1.49	1.37
2	L	3	SER	C-N	6.53	1.46	1.34
2	L	2	PHE	CE1-CZ	6.49	1.49	1.37
2	I	2	PHE	CE1-CZ	6.49	1.49	1.37
2	H	2	PHE	CE1-CZ	6.48	1.49	1.37
2	K	2	PHE	CE1-CZ	6.47	1.49	1.37
2	M	2	PHE	CE1-CZ	6.42	1.49	1.37
2	H	3	SER	C-N	6.03	1.45	1.34
2	I	3	SER	C-N	6.01	1.45	1.34
2	N	3	SER	C-N	6.01	1.45	1.34
2	M	3	SER	C-N	6.01	1.45	1.34
2	H	2	PHE	CE2-CZ	5.92	1.48	1.37
2	M	4	PRO	CA-C	5.89	1.64	1.52
2	K	3	SER	C-N	5.87	1.45	1.34
2	N	4	PRO	CA-C	5.84	1.64	1.52
2	J	2	PHE	CE2-CZ	5.82	1.48	1.37
2	I	2	PHE	CE2-CZ	5.81	1.48	1.37
2	J	3	SER	C-N	5.81	1.45	1.34
2	K	2	PHE	CE2-CZ	5.81	1.48	1.37
2	L	2	PHE	CE2-CZ	5.79	1.48	1.37
2	M	2	PHE	CE2-CZ	5.77	1.48	1.37
2	M	4	PRO	N-CA	5.77	1.57	1.47
2	N	2	PHE	CE2-CZ	5.75	1.48	1.37
2	J	4	PRO	CA-C	5.72	1.64	1.52
2	K	4	PRO	N-CA	5.61	1.56	1.47
2	L	4	PRO	N-CA	5.58	1.56	1.47
2	N	4	PRO	N-CA	5.54	1.56	1.47
2	L	4	PRO	CA-C	5.50	1.63	1.52
2	H	4	PRO	N-CA	5.49	1.56	1.47
2	K	2	PHE	CD1-CE1	5.48	1.50	1.39
2	H	4	PRO	CA-C	5.48	1.63	1.52
2	K	4	PRO	CA-C	5.43	1.63	1.52
2	I	4	PRO	CA-C	5.39	1.63	1.52
2	H	2	PHE	CD1-CE1	5.34	1.50	1.39
2	I	2	PHE	CD1-CE1	5.33	1.50	1.39
2	I	4	PRO	N-CA	5.31	1.56	1.47
2	N	2	PHE	CD1-CE1	5.25	1.49	1.39
2	J	4	PRO	N-CA	5.19	1.56	1.47
2	M	2	PHE	CD1-CE1	5.18	1.49	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	J	2	PHE	CD1-CE1	5.13	1.49	1.39
2	L	2	PHE	CD1-CE1	5.10	1.49	1.39

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	I	2	PHE	CB-CG-CD2	-5.57	116.90	120.80
2	N	2	PHE	CB-CG-CD2	-5.48	116.96	120.80
2	J	2	PHE	CB-CG-CD2	-5.47	116.97	120.80
2	H	2	PHE	CB-CG-CD2	-5.45	116.99	120.80
2	M	2	PHE	CB-CG-CD2	-5.37	117.04	120.80
2	K	2	PHE	CB-CG-CD2	-5.31	117.08	120.80
2	L	2	PHE	CB-CG-CD2	-5.15	117.20	120.80

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	1326	0	1366	36	0
1	B	1326	0	1366	26	0
1	C	1325	0	1364	36	0
1	D	1325	0	1364	23	0
1	E	1330	0	1369	36	0
1	F	1325	0	1364	38	0
1	G	1325	0	1364	35	0
2	H	52	0	49	1	0
2	I	52	0	49	1	0
2	J	52	0	49	2	0
2	K	52	0	49	1	0
2	L	52	0	49	2	0
2	M	52	0	49	1	0
2	N	52	0	49	4	0
3	A	4	0	6	0	0
3	B	4	0	6	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	4	0	6	0	0
3	D	4	0	6	0	0
3	E	4	0	6	0	0
3	F	4	0	6	0	0
3	G	4	0	6	0	0
4	A	13	0	17	0	0
4	B	13	0	16	1	0
4	C	13	0	17	1	0
4	D	13	0	16	0	0
4	E	13	0	16	0	0
4	F	13	0	16	0	0
4	G	13	0	16	0	0
5	A	111	0	0	3	0
5	B	117	0	0	3	0
5	C	69	0	0	0	0
5	D	105	0	0	3	0
5	E	78	0	0	1	0
5	F	109	0	0	3	0
5	G	75	0	0	1	0
5	J	1	0	0	0	0
5	N	2	0	0	0	0
All	All	10432	0	10056	213	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 11.

All (213) 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:F:84:LYS:HZ2	1:F:84:LYS:H	1.13	0.96
1:C:84:LYS:HE3	1:C:84:LYS:H	1.38	0.89
1:A:116:ASN:HD22	1:G:148:LYS:HD2	1.40	0.85
1:E:41:ASN:HD21	1:F:64:ASN:HD22	1.25	0.85
1:D:41:ASN:HD21	1:E:64:ASN:HD22	1.26	0.84
1:E:133:THR:HG21	5:F:280:HOH:O	1.77	0.83
1:C:86:LYS:HD2	1:C:107:GLU:OE1	1.80	0.82
1:C:84:LYS:H	1:C:84:LYS:CE	1.94	0.80
1:F:84:LYS:H	1:F:84:LYS:NZ	1.80	0.80
1:A:148:LYS:HD2	1:B:116:ASN:HD22	1.48	0.78
1:F:123:GLN:HE22	1:F:146:ARG:HH21	1.29	0.78
1:E:148:LYS:HD2	1:F:116:ASN:HD22	1.48	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:84:LYS:HE3	1:C:84:LYS:N	1.99	0.78
1:F:148:LYS:HD2	1:G:116:ASN:HD22	1.49	0.76
1:C:70:ILE:HD11	1:C:124:PRO:HB3	1.67	0.74
1:E:70:ILE:HD11	1:E:124:PRO:HB3	1.69	0.74
1:G:22:ARG:HE	1:G:25:LYS:HD2	1.54	0.72
1:G:151:LYS:NZ	1:G:151:LYS:HB2	2.05	0.71
1:E:27:ARG:NE	1:E:57:LYS:HB3	2.06	0.71
1:F:123:GLN:NE2	1:F:146:ARG:HH21	1.89	0.71
1:E:88:SER:HB3	1:E:110:LYS:HB3	1.74	0.70
1:F:84:LYS:HZ2	1:F:84:LYS:N	1.88	0.69
1:A:160:PRO:HG2	1:A:163:VAL:HG23	1.76	0.67
1:E:89:THR:C	1:E:90:ILE:HD12	2.15	0.67
1:A:116:ASN:ND2	1:G:148:LYS:HD2	2.10	0.67
1:F:160:PRO:HG2	1:F:163:VAL:HG23	1.77	0.66
1:C:90:ILE:HD11	2:K:6:ALA:HB1	1.77	0.66
1:A:160:PRO:HG2	1:A:163:VAL:CG2	2.27	0.65
1:E:160:PRO:HG2	1:E:163:VAL:HG23	1.79	0.64
1:F:133:THR:O	1:F:137:ILE:HG12	1.96	0.64
1:E:161:LEU:O	1:E:165:GLU:HG3	1.97	0.63
1:D:62:TYR:CE1	1:D:90:ILE:HD12	2.33	0.63
1:D:33:SER:O	1:D:65:SER:HB2	1.99	0.62
1:C:151:LYS:O	1:C:155:GLU:HG3	1.98	0.62
1:C:161:LEU:O	1:C:165:GLU:HG3	1.99	0.62
1:C:160:PRO:HG2	1:C:163:VAL:HG23	1.82	0.62
1:G:160:PRO:HG2	1:G:163:VAL:HG23	1.82	0.61
1:F:84:LYS:H	1:F:84:LYS:CE	2.14	0.61
1:C:77:TYR:O	1:C:81:GLN:HG2	2.01	0.60
1:F:160:PRO:HG2	1:F:163:VAL:CG2	2.32	0.60
1:C:90:ILE:N	1:C:90:ILE:HD12	2.16	0.60
1:G:115:PRO:HD3	1:G:189:LEU:O	2.01	0.60
1:A:70:ILE:HG12	1:A:98:MET:CE	2.31	0.60
1:E:90:ILE:HD11	2:M:6:ALA:HB1	1.82	0.60
1:C:29:ILE:HD12	1:C:46:GLN:HB3	1.84	0.60
1:E:41:ASN:ND2	1:F:64:ASN:HD22	1.98	0.60
1:F:90:ILE:HD13	1:F:112:TYR:HB2	1.84	0.60
1:A:70:ILE:HG12	1:A:98:MET:HE3	1.84	0.59
1:F:146:ARG:HD2	5:F:455:HOH:O	2.03	0.59
1:G:161:LEU:O	1:G:165:GLU:HG3	2.02	0.59
1:E:90:ILE:HD12	1:E:90:ILE:N	2.18	0.59
1:F:40:ALA:HB2	1:F:72:ALA:HB1	1.84	0.58
1:D:27:ARG:NE	1:D:57:LYS:HB3	2.17	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:70:ILE:HD11	1:D:124:PRO:HB3	1.84	0.58
1:F:31:LEU:HD13	1:F:43:ILE:HD11	1.84	0.58
1:B:148:LYS:HD2	1:C:116:ASN:HD22	1.68	0.58
1:D:41:ASN:ND2	1:E:64:ASN:HD22	1.98	0.57
1:G:62:TYR:CE1	1:G:90:ILE:HD12	2.38	0.57
1:G:152:VAL:O	1:G:156:ARG:HG2	2.03	0.57
1:C:40:ALA:O	1:C:44:VAL:HG23	2.04	0.57
1:C:51:ALA:HB1	1:C:84:LYS:NZ	2.18	0.57
1:A:22:ARG:NH2	1:A:25:LYS:HD2	2.20	0.57
1:A:133:THR:HG21	5:B:238:HOH:O	2.04	0.57
1:D:147:ASP:O	1:D:151:LYS:HG3	2.05	0.56
1:E:160:PRO:HG2	1:E:163:VAL:CG2	2.36	0.56
1:F:90:ILE:HD11	2:J:6:ALA:CB	2.35	0.56
1:F:115:PRO:HD3	1:F:189:LEU:O	2.04	0.56
1:A:161:LEU:O	1:A:165:GLU:HG3	2.06	0.56
1:C:89:THR:C	1:C:90:ILE:HD12	2.26	0.56
1:E:133:THR:O	1:E:137:ILE:HG12	2.06	0.56
1:F:31:LEU:HD13	1:F:43:ILE:CD1	2.37	0.55
1:D:22:ARG:HG3	2:L:1:OTT:H8A	1.88	0.55
1:F:187:LYS:HG3	5:F:328:HOH:O	2.06	0.55
1:D:31:LEU:HG	1:D:43:ILE:CD1	2.37	0.54
1:B:62:TYR:CE1	1:B:90:ILE:HD12	2.43	0.54
1:A:62:TYR:CE1	1:A:90:ILE:HD13	2.43	0.54
1:F:19:ILE:HG23	1:F:20:TYR:CD2	2.43	0.54
1:C:172:ASN:C	1:C:172:ASN:HD22	2.10	0.54
1:F:27:ARG:NE	1:F:57:LYS:HB3	2.22	0.53
1:G:108:LYS:HE2	1:G:156:ARG:O	2.08	0.53
1:D:77:TYR:O	1:D:81:GLN:HG2	2.09	0.53
1:E:172:ASN:C	1:E:172:ASN:HD22	2.11	0.53
1:B:55:PRO:HA	1:B:85:PRO:HG3	1.90	0.53
1:F:77:TYR:O	1:F:81:GLN:HG2	2.07	0.53
1:A:77:TYR:O	1:A:81:GLN:HG2	2.09	0.52
1:C:148:LYS:HD2	1:D:116:ASN:HD22	1.75	0.52
1:G:55:PRO:HA	1:G:85:PRO:HG3	1.92	0.52
1:B:172:ASN:C	1:B:172:ASN:HD22	2.11	0.51
1:C:160:PRO:HG2	1:C:163:VAL:CG2	2.39	0.51
1:G:22:ARG:HH21	1:G:25:LYS:HB2	1.74	0.51
1:A:90:ILE:N	1:A:90:ILE:HD12	2.25	0.51
1:C:84:LYS:HB2	1:C:85:PRO:HD3	1.93	0.51
1:E:54:ASP:HB3	1:E:57:LYS:HB2	1.92	0.51
1:B:115:PRO:HD3	1:B:189:LEU:O	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:151:LYS:HB2	1:G:151:LYS:HZ2	1.76	0.51
1:B:31:LEU:HD13	1:B:43:ILE:HD11	1.93	0.51
1:E:27:ARG:CZ	1:E:57:LYS:HB3	2.41	0.51
1:B:31:LEU:HD13	1:B:43:ILE:CD1	2.41	0.50
1:B:27:ARG:NE	1:B:57:LYS:HB3	2.26	0.50
1:D:144:LEU:HD23	5:D:219:HOH:O	2.11	0.50
1:B:191:HIS:CG	5:B:584:HOH:O	2.63	0.50
1:B:81:GLN:NE2	1:B:156:ARG:HH22	2.10	0.50
1:E:84:LYS:HB2	1:E:85:PRO:HD3	1.94	0.49
1:A:62:TYR:CD1	1:A:90:ILE:HD13	2.47	0.49
1:G:160:PRO:HG2	1:G:163:VAL:CG2	2.41	0.49
1:A:40:ALA:HB2	1:A:72:ALA:HB1	1.93	0.49
1:A:70:ILE:HD11	1:A:124:PRO:HB3	1.93	0.49
1:E:152:VAL:O	1:E:156:ARG:HG2	2.12	0.49
1:G:90:ILE:HD11	2:N:6:ALA:HB3	1.94	0.49
1:B:19:ILE:O	1:B:23:LEU:HD13	2.12	0.49
1:C:124:PRO:HD2	1:C:146:ARG:HG3	1.95	0.49
1:A:55:PRO:HA	1:A:85:PRO:HG3	1.95	0.49
1:B:90:ILE:HD11	2:H:6:ALA:HB3	1.95	0.49
1:E:148:LYS:HD2	1:F:116:ASN:ND2	2.23	0.49
1:D:172:ASN:C	1:D:172:ASN:HD22	2.17	0.48
1:B:70:ILE:HD11	1:B:124:PRO:HB3	1.93	0.48
1:G:22:ARG:HG3	2:N:1:OTT:C8	2.43	0.48
1:C:55:PRO:HA	1:C:85:PRO:HG3	1.96	0.48
1:G:92:ILE:HG13	1:G:92:ILE:O	2.13	0.48
1:A:90:ILE:HD11	2:I:6:ALA:HB1	1.94	0.48
1:C:62:TYR:CD1	1:C:90:ILE:HD13	2.49	0.48
1:F:90:ILE:CD1	1:F:112:TYR:HB2	2.42	0.48
1:G:172:ASN:C	1:G:172:ASN:HD22	2.16	0.48
1:B:83:ILE:HB	1:B:85:PRO:HD2	1.95	0.47
1:B:166:ARG:HD2	5:B:222:HOH:O	2.12	0.47
1:F:151:LYS:O	1:F:155:GLU:HG3	2.14	0.47
1:E:19:ILE:HG23	1:E:20:TYR:CD1	2.48	0.47
1:E:58:GLU:OE2	1:E:110:LYS:HE2	2.15	0.47
1:G:83:ILE:HB	1:G:85:PRO:HD2	1.97	0.47
1:D:81:GLN:NE2	1:D:156:ARG:HH22	2.13	0.47
1:G:151:LYS:HB2	1:G:151:LYS:HZ3	1.78	0.47
1:C:92:ILE:O	1:C:92:ILE:HG13	2.15	0.47
4:C:503:NHE:H3'1	1:D:66:PRO:HB3	1.97	0.46
1:C:62:TYR:CE1	1:C:90:ILE:HD13	2.50	0.46
1:B:160:PRO:HG2	1:B:163:VAL:HG23	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:45:SER:HB3	1:B:19:ILE:HG12	1.98	0.46
1:A:89:THR:C	1:A:90:ILE:HD12	2.36	0.46
1:C:88:SER:HB3	1:C:110:LYS:CG	2.45	0.46
1:C:88:SER:HB3	1:C:110:LYS:HG2	1.98	0.45
1:A:22:ARG:CZ	1:A:25:LYS:HD2	2.46	0.45
1:B:81:GLN:HE22	1:B:156:ARG:HH22	1.64	0.45
1:C:84:LYS:N	1:C:85:PRO:CD	2.79	0.45
1:B:40:ALA:O	1:B:44:VAL:HG23	2.16	0.45
1:A:24:LEU:CD2	1:A:50:LEU:HD21	2.47	0.45
1:E:41:ASN:HD21	1:F:64:ASN:ND2	2.04	0.45
1:A:115:PRO:HD3	1:A:189:LEU:O	2.16	0.45
1:B:58:GLU:OE2	1:B:110:LYS:HE3	2.16	0.45
1:E:147:ASP:O	1:E:151:LYS:HG3	2.16	0.45
1:G:51:ALA:HA	1:G:85:PRO:HG2	1.98	0.45
1:E:40:ALA:HB2	1:E:72:ALA:HB1	1.99	0.45
1:F:90:ILE:HD11	2:J:6:ALA:HB1	2.00	0.44
1:G:19:ILE:HD13	1:G:19:ILE:O	2.17	0.44
1:A:129:GLN:HE21	1:A:129:GLN:HB3	1.64	0.44
1:G:54:ASP:OD2	1:G:57:LYS:HB2	2.17	0.44
1:A:172:ASN:HD22	1:A:172:ASN:C	2.21	0.44
1:C:51:ALA:HA	1:C:85:PRO:HG2	1.99	0.44
1:A:83:ILE:HB	1:A:85:PRO:HD2	2.00	0.44
1:F:70:ILE:HD11	1:F:124:PRO:HB3	2.00	0.44
1:E:70:ILE:CD1	1:E:124:PRO:HB3	2.43	0.43
1:F:70:ILE:O	1:F:74:MET:HG2	2.18	0.43
1:G:39:VAL:O	1:G:43:ILE:HG12	2.18	0.43
1:B:40:ALA:HB2	1:B:72:ALA:HB1	1.99	0.43
1:F:18:ASP:N	1:F:21:SER:HG	2.16	0.43
1:C:114:LEU:HD23	1:C:189:LEU:HB2	2.00	0.43
1:F:177:GLU:O	1:F:181:GLU:HG3	2.18	0.43
1:A:49:PHE:HB2	1:B:19:ILE:HD11	2.01	0.43
1:A:156:ARG:HD2	5:A:251:HOH:O	2.18	0.43
1:F:172:ASN:C	1:F:172:ASN:HD22	2.22	0.43
1:G:133:THR:O	1:G:137:ILE:HG12	2.18	0.43
1:D:90:ILE:HD11	2:L:6:ALA:HB3	2.01	0.43
1:A:155:GLU:HG3	5:A:452:HOH:O	2.19	0.42
1:F:86:LYS:HD3	1:F:107:GLU:HG2	2.01	0.42
1:F:40:ALA:O	1:F:44:VAL:HG23	2.19	0.42
1:D:23:LEU:HD12	1:D:28:ILE:HD12	2.00	0.42
1:A:84:LYS:N	1:A:85:PRO:CD	2.82	0.42
1:D:19:ILE:O	1:D:23:LEU:HD23	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:153:LEU:HD23	1:D:153:LEU:HA	1.88	0.42
1:G:189:LEU:HD13	2:N:2:PHE:HB3	2.02	0.42
1:B:84:LYS:N	1:B:85:PRO:CD	2.83	0.42
5:D:220:HOH:O	1:E:33:SER:HA	2.19	0.42
1:E:166:ARG:HG3	5:E:571:HOH:O	2.18	0.42
1:G:19:ILE:O	1:G:23:LEU:HD13	2.20	0.42
1:G:84:LYS:N	1:G:85:PRO:CD	2.83	0.42
1:C:21:SER:O	1:C:25:LYS:HG2	2.19	0.42
1:B:21:SER:O	1:B:25:LYS:HG3	2.19	0.41
1:C:83:ILE:HB	1:C:85:PRO:HD2	2.01	0.41
1:C:29:ILE:CD1	1:C:46:GLN:HB3	2.48	0.41
1:C:153:LEU:O	1:C:157:THR:HG23	2.20	0.41
1:F:51:ALA:HB1	1:F:84:LYS:NZ	2.35	0.41
1:E:112:TYR:CD2	1:E:189:LEU:HD21	2.55	0.41
1:G:86:LYS:HD3	5:G:215:HOH:O	2.20	0.41
1:G:153:LEU:HD12	1:G:153:LEU:HA	1.95	0.41
1:G:175:SER:OG	1:G:178:GLU:HG3	2.21	0.41
4:B:502:NHE:H3'1	1:C:66:PRO:HB3	2.01	0.41
1:F:84:LYS:N	1:F:85:PRO:CD	2.82	0.41
1:A:51:ALA:HA	1:A:85:PRO:HG2	2.03	0.41
1:A:54:ASP:OD1	1:A:57:LYS:HB2	2.21	0.41
1:A:148:LYS:HD2	1:B:116:ASN:ND2	2.27	0.41
1:G:40:ALA:HB2	1:G:72:ALA:HB1	2.02	0.41
1:A:70:ILE:HG12	1:A:98:MET:HE1	2.01	0.41
1:D:75:ALA:HB1	1:E:92:ILE:HG22	2.02	0.41
1:E:18:ASP:HB3	1:E:21:SER:OG	2.21	0.41
1:E:92:ILE:HG23	1:E:114:LEU:HD12	2.03	0.41
1:A:18:ASP:OD2	1:A:20:TYR:HB2	2.21	0.40
1:A:151:LYS:HD2	5:A:437:HOH:O	2.20	0.40
1:D:31:LEU:HG	1:D:43:ILE:HD11	2.03	0.40
1:D:166:ARG:HD2	5:D:230:HOH:O	2.21	0.40
1:E:19:ILE:HG23	1:E:20:TYR:N	2.36	0.40
1:D:84:LYS:N	1:D:85:PRO:CD	2.84	0.40
1:E:77:TYR:O	1:E:81:GLN:HG2	2.22	0.40
1:G:90:ILE:HD11	2:N:6:ALA:CB	2.50	0.40
1:G:95:ALA:O	1:G:100:ALA:HB2	2.20	0.40
1:C:36:ASP:C	1:C:36:ASP:OD1	2.60	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	172/199 (86%)	165 (96%)	7 (4%)	0	100	100
1	B	172/199 (86%)	166 (96%)	6 (4%)	0	100	100
1	C	172/199 (86%)	165 (96%)	7 (4%)	0	100	100
1	D	172/199 (86%)	169 (98%)	3 (2%)	0	100	100
1	E	172/199 (86%)	165 (96%)	7 (4%)	0	100	100
1	F	172/199 (86%)	166 (96%)	6 (4%)	0	100	100
1	G	172/199 (86%)	167 (97%)	5 (3%)	0	100	100
2	H	3/7 (43%)	2 (67%)	1 (33%)	0	100	100
2	I	3/7 (43%)	2 (67%)	1 (33%)	0	100	100
2	J	3/7 (43%)	2 (67%)	1 (33%)	0	100	100
2	K	3/7 (43%)	2 (67%)	1 (33%)	0	100	100
2	L	3/7 (43%)	2 (67%)	1 (33%)	0	100	100
2	M	3/7 (43%)	2 (67%)	1 (33%)	0	100	100
2	N	3/7 (43%)	2 (67%)	1 (33%)	0	100	100
All	All	1225/1442 (85%)	1177 (96%)	48 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	141/165 (86%)	134 (95%)	7 (5%)	24	20
1	B	141/165 (86%)	135 (96%)	6 (4%)	29	26
1	C	141/165 (86%)	136 (96%)	5 (4%)	36	35
1	D	141/165 (86%)	133 (94%)	8 (6%)	20	16
1	E	142/165 (86%)	135 (95%)	7 (5%)	25	21
1	F	141/165 (86%)	137 (97%)	4 (3%)	43	44
1	G	141/165 (86%)	135 (96%)	6 (4%)	29	26
2	H	3/3 (100%)	3 (100%)	0	100	100
2	I	3/3 (100%)	3 (100%)	0	100	100
2	J	3/3 (100%)	3 (100%)	0	100	100
2	K	3/3 (100%)	3 (100%)	0	100	100
2	L	3/3 (100%)	3 (100%)	0	100	100
2	M	3/3 (100%)	3 (100%)	0	100	100
2	N	3/3 (100%)	3 (100%)	0	100	100
All	All	1009/1176 (86%)	966 (96%)	43 (4%)	29	26

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	LEU
1	A	26	ASP
1	A	30	MET
1	A	122	HIS
1	A	129	GLN
1	A	170	ARG
1	A	172	ASN
1	B	114	LEU
1	B	122	HIS
1	B	123	GLN
1	B	129	GLN
1	B	170	ARG
1	B	172	ASN
1	C	25	LYS
1	C	84	LYS
1	C	122	HIS
1	C	170	ARG
1	C	172	ASN
1	D	26	ASP

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Mol	Chain	Res	Type
1	D	30	MET
1	D	31	LEU
1	D	65	SER
1	D	84	LYS
1	D	122	HIS
1	D	144	LEU
1	D	172	ASN
1	E	25	LYS
1	E	30	MET
1	E	97	SER
1	E	122	HIS
1	E	123	GLN
1	E	172	ASN
1	E	191	HIS
1	F	26	ASP
1	F	84	LYS
1	F	122	HIS
1	F	172	ASN
1	G	19	ILE
1	G	22	ARG
1	G	86	LYS
1	G	97	SER
1	G	122	HIS
1	G	172	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	ASN
1	A	116	ASN
1	A	129	GLN
1	A	172	ASN
1	B	38	ASN
1	B	81	GLN
1	B	116	ASN
1	B	123	GLN
1	B	129	GLN
1	B	172	ASN
1	C	116	ASN
1	C	123	GLN
1	C	172	ASN
1	D	38	ASN

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Mol	Chain	Res	Type
1	D	41	ASN
1	D	81	GLN
1	D	116	ASN
1	D	123	GLN
1	D	172	ASN
1	E	41	ASN
1	E	116	ASN
1	E	123	GLN
1	E	172	ASN
1	E	191	HIS
1	F	38	ASN
1	F	81	GLN
1	F	116	ASN
1	F	123	GLN
1	F	172	ASN
1	G	116	ASN
1	G	172	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)

14 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	MAA	J	5	2	4,5,6	3.35	3 (75%)	2,5,7	0.96	0
2	MAA	H	5	2	4,5,6	3.46	3 (75%)	2,5,7	0.92	0
2	MP8	L	7	2	6,8,9	1.10	0	3,10,12	1.22	0
2	MP8	M	7	2	6,8,9	1.05	0	3,10,12	1.27	0
2	MP8	I	7	2	6,8,9	1.02	0	3,10,12	1.24	0
2	MP8	N	7	2	6,8,9	1.02	0	3,10,12	1.24	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	MAA	K	5	2	4,5,6	3.45	4 (100%)	2,5,7	0.96	0
2	MP8	H	7	2	6,8,9	1.01	0	3,10,12	1.32	0
2	MP8	K	7	2	6,8,9	1.00	0	3,10,12	1.26	0
2	MAA	L	5	2	4,5,6	3.45	3 (75%)	2,5,7	0.92	0
2	MAA	M	5	2	4,5,6	3.34	3 (75%)	2,5,7	0.96	0
2	MAA	I	5	2	4,5,6	3.26	3 (75%)	2,5,7	1.01	0
2	MAA	N	5	2	4,5,6	3.40	4 (100%)	2,5,7	0.89	0
2	MP8	J	7	2	6,8,9	0.95	0	3,10,12	1.20	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MAA	J	5	2	-	0/2/4/6	-
2	MAA	H	5	2	-	0/2/4/6	-
2	MP8	L	7	2	-	0/0/11/13	0/1/1/1
2	MP8	M	7	2	-	0/0/11/13	0/1/1/1
2	MP8	I	7	2	-	0/0/11/13	0/1/1/1
2	MP8	N	7	2	-	0/0/11/13	0/1/1/1
2	MAA	K	5	2	-	0/2/4/6	-
2	MP8	H	7	2	-	0/0/11/13	0/1/1/1
2	MP8	K	7	2	-	0/0/11/13	0/1/1/1
2	MAA	L	5	2	-	0/2/4/6	-
2	MAA	M	5	2	-	0/2/4/6	-
2	MAA	I	5	2	-	0/2/4/6	-
2	MAA	N	5	2	-	0/2/4/6	-
2	MP8	J	7	2	-	0/0/11/13	0/1/1/1

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	5	MAA	CA-N	5.43	1.56	1.47
2	K	5	MAA	CA-N	5.39	1.56	1.47
2	L	5	MAA	CA-N	5.31	1.56	1.47
2	J	5	MAA	CA-N	5.27	1.56	1.47
2	N	5	MAA	CA-N	5.22	1.56	1.47
2	M	5	MAA	CA-N	5.17	1.56	1.47
2	I	5	MAA	CA-N	5.10	1.56	1.47
2	L	5	MAA	CM-N	3.04	1.54	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	M	5	MAA	CM-N	3.01	1.54	1.46
2	N	5	MAA	CM-N	2.96	1.54	1.46
2	K	5	MAA	CM-N	2.91	1.54	1.46
2	H	5	MAA	CM-N	2.88	1.53	1.46
2	J	5	MAA	CM-N	2.87	1.53	1.46
2	I	5	MAA	CM-N	2.66	1.53	1.46
2	H	5	MAA	CB-CA	2.50	1.58	1.51
2	K	5	MAA	CB-CA	2.49	1.58	1.51
2	L	5	MAA	CB-CA	2.48	1.58	1.51
2	N	5	MAA	CB-CA	2.43	1.58	1.51
2	I	5	MAA	CB-CA	2.35	1.58	1.51
2	M	5	MAA	CB-CA	2.34	1.58	1.51
2	J	5	MAA	CB-CA	2.28	1.57	1.51
2	N	5	MAA	O-C	2.05	1.27	1.20
2	K	5	MAA	O-C	2.01	1.27	1.20

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

14 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	DMS	E	305	-	3,3,3	0.42	0	3,3,3	0.91	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	DMS	B	302	-	3,3,3	0.45	0	3,3,3	0.89	0
3	DMS	F	306	-	3,3,3	0.45	0	3,3,3	0.91	0
3	DMS	A	301	-	3,3,3	0.43	0	3,3,3	0.92	0
4	NHE	G	507	-	13,13,13	2.26	3 (23%)	16,17,17	5.82	12 (75%)
3	DMS	C	303	-	3,3,3	0.39	0	3,3,3	0.91	0
4	NHE	E	505	-	13,13,13	2.13	2 (15%)	16,17,17	5.85	12 (75%)
4	NHE	A	501	-	13,13,13	2.29	2 (15%)	16,17,17	5.75	10 (62%)
4	NHE	D	504	-	13,13,13	2.23	3 (23%)	16,17,17	5.86	12 (75%)
4	NHE	F	506	-	13,13,13	2.23	2 (15%)	16,17,17	5.83	12 (75%)
4	NHE	B	502	-	13,13,13	2.30	2 (15%)	16,17,17	5.85	11 (68%)
3	DMS	D	304	-	3,3,3	0.45	0	3,3,3	0.90	0
3	DMS	G	307	-	3,3,3	0.42	0	3,3,3	0.92	0
4	NHE	C	503	-	13,13,13	2.37	2 (15%)	16,17,17	5.82	11 (68%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NHE	G	507	-	-	5/7/15/15	0/1/1/1
4	NHE	B	502	-	-	5/7/15/15	0/1/1/1
4	NHE	E	505	-	-	5/7/15/15	0/1/1/1
4	NHE	A	501	-	-	3/7/15/15	0/1/1/1
4	NHE	D	504	-	-	3/7/15/15	0/1/1/1
4	NHE	F	506	-	-	5/7/15/15	0/1/1/1
4	NHE	C	503	-	-	3/7/15/15	0/1/1/1

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	503	NHE	C2-S	-7.44	1.67	1.77
4	B	502	NHE	C2-S	-7.11	1.67	1.77
4	A	501	NHE	C2-S	-7.09	1.67	1.77
4	G	507	NHE	C2-S	-6.96	1.67	1.77
4	F	506	NHE	C2-S	-6.93	1.67	1.77
4	D	504	NHE	C2-S	-6.90	1.67	1.77
4	E	505	NHE	C2-S	-6.48	1.68	1.77
4	B	502	NHE	O1-S	2.06	1.50	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	501	NHE	O1-S	2.06	1.50	1.45
4	F	506	NHE	O1-S	2.05	1.50	1.45
4	D	504	NHE	O2-S	2.05	1.50	1.45
4	C	503	NHE	O2-S	2.04	1.50	1.45
4	G	507	NHE	O1-S	2.04	1.50	1.45
4	E	505	NHE	O1-S	2.02	1.50	1.45
4	G	507	NHE	O2-S	2.02	1.50	1.45
4	D	504	NHE	O1-S	2.01	1.50	1.45

All (80) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	B	502	NHE	O3-S-C2	-13.22	80.13	106.00
4	D	504	NHE	O3-S-C2	-13.15	80.28	106.00
4	F	506	NHE	O3-S-C2	-12.97	80.64	106.00
4	E	505	NHE	O3-S-C2	-12.94	80.69	106.00
4	G	507	NHE	O3-S-C2	-12.86	80.85	106.00
4	A	501	NHE	O3-S-C2	-12.86	80.85	106.00
4	C	503	NHE	O3-S-C2	-12.85	80.87	106.00
4	F	506	NHE	O1-S-C2	11.78	124.52	106.73
4	E	505	NHE	O1-S-C2	11.75	124.49	106.73
4	D	504	NHE	O1-S-C2	11.73	124.46	106.73
4	G	507	NHE	O1-S-C2	11.57	124.21	106.73
4	B	502	NHE	O1-S-C2	11.49	124.10	106.73
4	A	501	NHE	O1-S-C2	11.47	124.06	106.73
4	C	503	NHE	O1-S-C2	11.41	123.97	106.73
4	D	504	NHE	O3-S-O1	-7.49	92.65	111.40
4	E	505	NHE	O3-S-O1	-7.44	92.79	111.40
4	F	506	NHE	O3-S-O1	-7.41	92.87	111.40
4	G	507	NHE	O3-S-O1	-7.29	93.15	111.40
4	C	503	NHE	O3-S-O2	-7.29	93.17	111.40
4	B	502	NHE	O3-S-O1	-7.28	93.18	111.40
4	F	506	NHE	O3-S-O2	-7.26	93.24	111.40
4	A	501	NHE	O3-S-O2	-7.25	93.26	111.40
4	G	507	NHE	O3-S-O2	-7.22	93.34	111.40
4	B	502	NHE	O3-S-O2	-7.22	93.34	111.40
4	A	501	NHE	O3-S-O1	-7.20	93.38	111.40
4	C	503	NHE	O3-S-O1	-7.18	93.43	111.40
4	D	504	NHE	O3-S-O2	-7.07	93.71	111.40
4	E	505	NHE	O3-S-O2	-6.97	93.95	111.40
4	C	503	NHE	C1-N-C1'	6.89	127.44	114.18
4	E	505	NHE	C1-N-C1'	6.54	126.78	114.18

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	G	507	NHE	C1-N-C1'	6.48	126.65	114.18
4	C	503	NHE	O2-S-C2	6.47	116.50	106.73
4	D	504	NHE	C1-N-C1'	6.39	126.48	114.18
4	A	501	NHE	C1-N-C1'	6.31	126.34	114.18
4	G	507	NHE	O2-S-C2	6.28	116.22	106.73
4	B	502	NHE	C1-N-C1'	6.27	126.25	114.18
4	F	506	NHE	C1-N-C1'	6.24	126.19	114.18
4	E	505	NHE	O2-S-C2	6.21	116.11	106.73
4	A	501	NHE	O2-S-C2	6.12	115.97	106.73
4	B	502	NHE	O2-S-C2	6.08	115.91	106.73
4	F	506	NHE	O2-S-C2	6.00	115.79	106.73
4	D	504	NHE	O2-S-C2	5.95	115.72	106.73
4	B	502	NHE	C3'-C2'-C1'	4.06	118.37	111.09
4	D	504	NHE	C3'-C2'-C1'	4.01	118.28	111.09
4	E	505	NHE	C3'-C2'-C1'	3.89	118.07	111.09
4	F	506	NHE	C3'-C2'-C1'	3.83	117.96	111.09
4	G	507	NHE	C3'-C2'-C1'	3.82	117.94	111.09
4	A	501	NHE	C3'-C2'-C1'	3.67	117.67	111.09
4	C	503	NHE	C3'-C2'-C1'	3.58	117.51	111.09
4	C	503	NHE	C5'-C6'-C1'	3.44	117.26	111.09
4	E	505	NHE	C5'-C6'-C1'	3.43	117.25	111.09
4	G	507	NHE	C5'-C6'-C1'	3.35	117.10	111.09
4	D	504	NHE	C5'-C6'-C1'	3.26	116.94	111.09
4	B	502	NHE	C5'-C6'-C1'	3.26	116.93	111.09
4	B	502	NHE	C6'-C1'-C2'	3.25	116.41	110.80
4	F	506	NHE	C5'-C6'-C1'	3.24	116.90	111.09
4	A	501	NHE	C5'-C6'-C1'	3.12	116.69	111.09
4	G	507	NHE	C6'-C1'-C2'	3.09	116.14	110.80
4	E	505	NHE	C6'-C1'-C2'	3.09	116.14	110.80
4	A	501	NHE	C6'-C1'-C2'	3.07	116.11	110.80
4	F	506	NHE	C6'-C1'-C2'	3.02	116.01	110.80
4	D	504	NHE	C6'-C1'-C2'	3.01	116.00	110.80
4	C	503	NHE	C6'-C1'-C2'	2.93	115.86	110.80
4	B	502	NHE	C4'-C3'-C2'	2.27	116.08	111.42
4	E	505	NHE	C5'-C4'-C3'	2.21	117.75	111.19
4	E	505	NHE	C4'-C3'-C2'	2.18	115.89	111.42
4	G	507	NHE	C5'-C4'-C3'	2.17	117.63	111.19
4	D	504	NHE	C4'-C3'-C2'	2.16	115.87	111.42
4	C	503	NHE	C5'-C4'-C3'	2.15	117.58	111.19
4	C	503	NHE	C4'-C3'-C2'	2.12	115.78	111.42
4	B	502	NHE	C5'-C4'-C3'	2.11	117.47	111.19
4	F	506	NHE	C5'-C4'-C3'	2.11	117.44	111.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	D	504	NHE	C4'-C5'-C6'	2.10	115.74	111.42
4	G	507	NHE	C4'-C3'-C2'	2.10	115.73	111.42
4	E	505	NHE	C4'-C5'-C6'	2.09	115.72	111.42
4	F	506	NHE	C4'-C5'-C6'	2.08	115.69	111.42
4	F	506	NHE	C4'-C3'-C2'	2.07	115.67	111.42
4	D	504	NHE	C5'-C4'-C3'	2.05	117.29	111.19
4	A	501	NHE	C4'-C3'-C2'	2.05	115.62	111.42
4	G	507	NHE	C4'-C5'-C6'	2.03	115.59	111.42

There are no chirality outliers.

All (29) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	501	NHE	N-C1-C2-S
4	A	501	NHE	C1-C2-S-O3
4	B	502	NHE	C2'-C1'-N-C1
4	B	502	NHE	N-C1-C2-S
4	B	502	NHE	C1-C2-S-O3
4	C	503	NHE	N-C1-C2-S
4	C	503	NHE	C1-C2-S-O1
4	D	504	NHE	N-C1-C2-S
4	D	504	NHE	C1-C2-S-O3
4	E	505	NHE	C1-C2-S-O3
4	G	507	NHE	C2'-C1'-N-C1
4	G	507	NHE	N-C1-C2-S
4	G	507	NHE	C1-C2-S-O3
4	C	503	NHE	C1-C2-S-O3
4	F	506	NHE	C1-C2-S-O3
4	F	506	NHE	C2'-C1'-N-C1
4	F	506	NHE	C6'-C1'-N-C1
4	A	501	NHE	C1-C2-S-O1
4	B	502	NHE	C1-C2-S-O1
4	D	504	NHE	C1-C2-S-O1
4	E	505	NHE	C1-C2-S-O1
4	F	506	NHE	C1-C2-S-O1
4	G	507	NHE	C1-C2-S-O1
4	F	506	NHE	N-C1-C2-S
4	B	502	NHE	C6'-C1'-N-C1
4	E	505	NHE	C6'-C1'-N-C1
4	G	507	NHE	C6'-C1'-N-C1
4	E	505	NHE	C2'-C1'-N-C1
4	E	505	NHE	N-C1-C2-S

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	502	NHE	1	0
4	C	503	NHE	1	0

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	174/199 (87%)	-0.03	4 (2%)	60	59	23, 32, 53, 80	0
1	B	174/199 (87%)	-0.10	7 (4%)	38	37	24, 32, 53, 78	0
1	C	174/199 (87%)	0.29	11 (6%)	20	19	27, 37, 64, 82	0
1	D	174/199 (87%)	0.04	10 (5%)	23	23	25, 34, 60, 81	0
1	E	174/199 (87%)	0.11	6 (3%)	45	44	27, 35, 57, 76	0
1	F	174/199 (87%)	-0.07	5 (2%)	51	50	24, 33, 54, 72	0
1	G	174/199 (87%)	0.01	4 (2%)	60	59	25, 36, 58, 75	0
2	H	4/7 (57%)	0.15	0	100	100	40, 46, 47, 48	0
2	I	4/7 (57%)	0.21	0	100	100	41, 47, 48, 50	0
2	J	4/7 (57%)	0.54	1 (25%)	0	0	35, 45, 48, 49	0
2	K	4/7 (57%)	0.62	1 (25%)	0	0	46, 55, 58, 59	0
2	L	4/7 (57%)	0.44	0	100	100	47, 55, 57, 58	0
2	M	4/7 (57%)	0.15	0	100	100	45, 48, 48, 52	0
2	N	4/7 (57%)	1.30	1 (25%)	0	0	47, 56, 58, 60	0
All	All	1246/1442 (86%)	0.05	50 (4%)	38	37	23, 34, 59, 82	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	191	HIS	8.5
1	B	191	HIS	4.8
1	D	19	ILE	4.8
1	B	19	ILE	4.6
1	A	19	ILE	4.4
2	N	4	PRO	4.1
1	D	22	ARG	3.8
1	E	19	ILE	3.7

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Mol	Chain	Res	Type	RSRZ
1	C	19	ILE	3.6
1	D	25	LYS	3.6
1	E	191	HIS	3.4
1	D	18	ASP	3.3
1	A	20	TYR	3.2
1	C	22	ARG	3.2
1	F	22	ARG	3.2
1	B	84	LYS	3.1
1	D	20	TYR	3.1
1	E	20	TYR	3.1
1	F	18	ASP	3.0
1	F	19	ILE	2.9
1	G	22	ARG	2.9
1	E	22	ARG	2.9
1	F	20	TYR	2.9
1	C	18	ASP	2.8
1	D	21	SER	2.8
1	B	18	ASP	2.8
1	A	22	ARG	2.7
1	G	191	HIS	2.7
1	C	21	SER	2.6
1	C	191	HIS	2.6
1	F	191	HIS	2.5
1	B	22	ARG	2.5
1	B	129	GLN	2.4
2	J	4	PRO	2.4
1	G	56	GLU	2.4
1	E	56	GLU	2.3
1	C	57	LYS	2.3
1	C	63	ILE	2.3
1	A	56	GLU	2.3
1	C	162	GLU	2.3
1	G	190	THR	2.3
1	C	190	THR	2.2
1	B	190	THR	2.2
2	K	4	PRO	2.2
1	E	166	ARG	2.2
1	D	56	GLU	2.1
1	C	52	ALA	2.1
1	D	51	ALA	2.1
1	C	76	ILE	2.0
1	D	57	LYS	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	MAA	K	5	6/7	0.85	0.19	58,59,59,60	0
2	MP8	N	7	8/9	0.85	0.17	53,55,55,55	0
2	MAA	N	5	6/7	0.86	0.16	58,59,59,60	0
2	MP8	K	7	8/9	0.89	0.12	57,57,58,58	0
2	MP8	L	7	8/9	0.90	0.25	56,56,56,57	0
2	MAA	L	5	6/7	0.91	0.16	56,57,57,57	0
2	MP8	H	7	8/9	0.94	0.12	47,47,49,49	0
2	MAA	M	5	6/7	0.94	0.14	49,50,51,51	0
2	MAA	I	5	6/7	0.94	0.12	48,49,50,52	0
2	MP8	J	7	8/9	0.94	0.12	45,46,46,47	0
2	MP8	M	7	8/9	0.95	0.09	42,45,45,46	0
2	MAA	H	5	6/7	0.95	0.13	44,45,46,46	0
2	MAA	J	5	6/7	0.95	0.16	47,47,48,49	0
2	MP8	I	7	8/9	0.96	0.11	46,48,48,49	0

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	NHE	D	504	13/13	0.88	0.19	63,66,67,68	0
4	NHE	B	502	13/13	0.89	0.18	59,64,65,65	0
3	DMS	G	307	4/4	0.90	0.18	64,65,65,66	0
3	DMS	B	302	4/4	0.92	0.20	63,63,64,64	0
3	DMS	F	306	4/4	0.94	0.19	67,68,68,70	0
3	DMS	D	304	4/4	0.94	0.17	63,64,65,65	0
4	NHE	F	506	13/13	0.94	0.15	59,61,62,63	0
4	NHE	C	503	13/13	0.95	0.13	35,40,45,46	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	NHE	A	501	13/13	0.95	0.13	41,49,54,54	0
3	DMS	E	305	4/4	0.95	0.19	64,65,65,67	0
4	NHE	E	505	13/13	0.96	0.12	41,46,48,48	0
4	NHE	G	507	13/13	0.96	0.10	38,41,42,43	0
3	DMS	C	303	4/4	0.97	0.15	65,65,66,67	0
3	DMS	A	301	4/4	0.98	0.17	58,59,59,60	0

6.5 Other polymers (i)

There are no such residues in this entry.