



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 07:01 AM GMT

PDB ID : 1D0X
Title : DICTYOSTELIUM MYOSIN S1DC (MOTOR DOMAIN FRAGMENT)
COMPLEXED WITH M-NITROPHENYL AMINOETHYLDIPHOSPHA
TEBERYLLIUM TRIFLUORIDE.
Authors : Gulick, A.M.; Bauer, C.B.; Thoden, J.B.; Pate, E.; Yount, R.G.; Rayment, I.
Deposited on : 1999-09-15
Resolution : 2.00 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

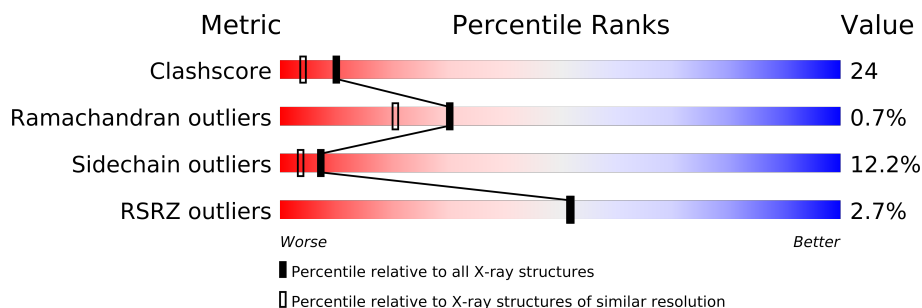
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

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))
Clashscore	79885	6188 (2.00-2.00)
Ramachandran outliers	78287	6102 (2.00-2.00)
Sidechain outliers	78261	6100 (2.00-2.00)
RSRZ outliers	66119	4890 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	761	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6643 atoms, of which 0 are hydrogen and 0 are deuterium.

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 MYOSIN S1DC MOTOR DOMAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	740	Total	C	N	O	S	0	35	0
			5887	3744	1017	1110	16			

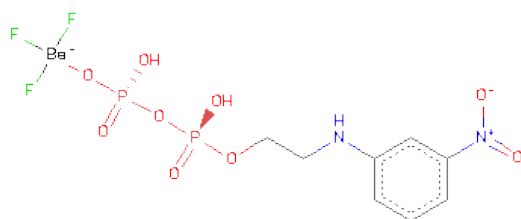
There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	312	CYS	TYR	SEE REMARK 999	UNP P08799
A	760	PRO	GLN	ENGINEERED	UNP P08799
A	761	ASN	ARG	ENGINEERED	UNP P08799

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		

- Molecule 3 is M-NITROPHENYL AMINOETHYLDIPHOSPHATEBERYLLIUM TRIFLUORIDE (three-letter code: MNQ) (formula: C₈H₁₁BeF₃N₂O₉P₂).



Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
			Total	Be	C	F	N	O	P		
3	A	1	25	1	8	3	2	9	2	0	0

- Molecule 4 is water.

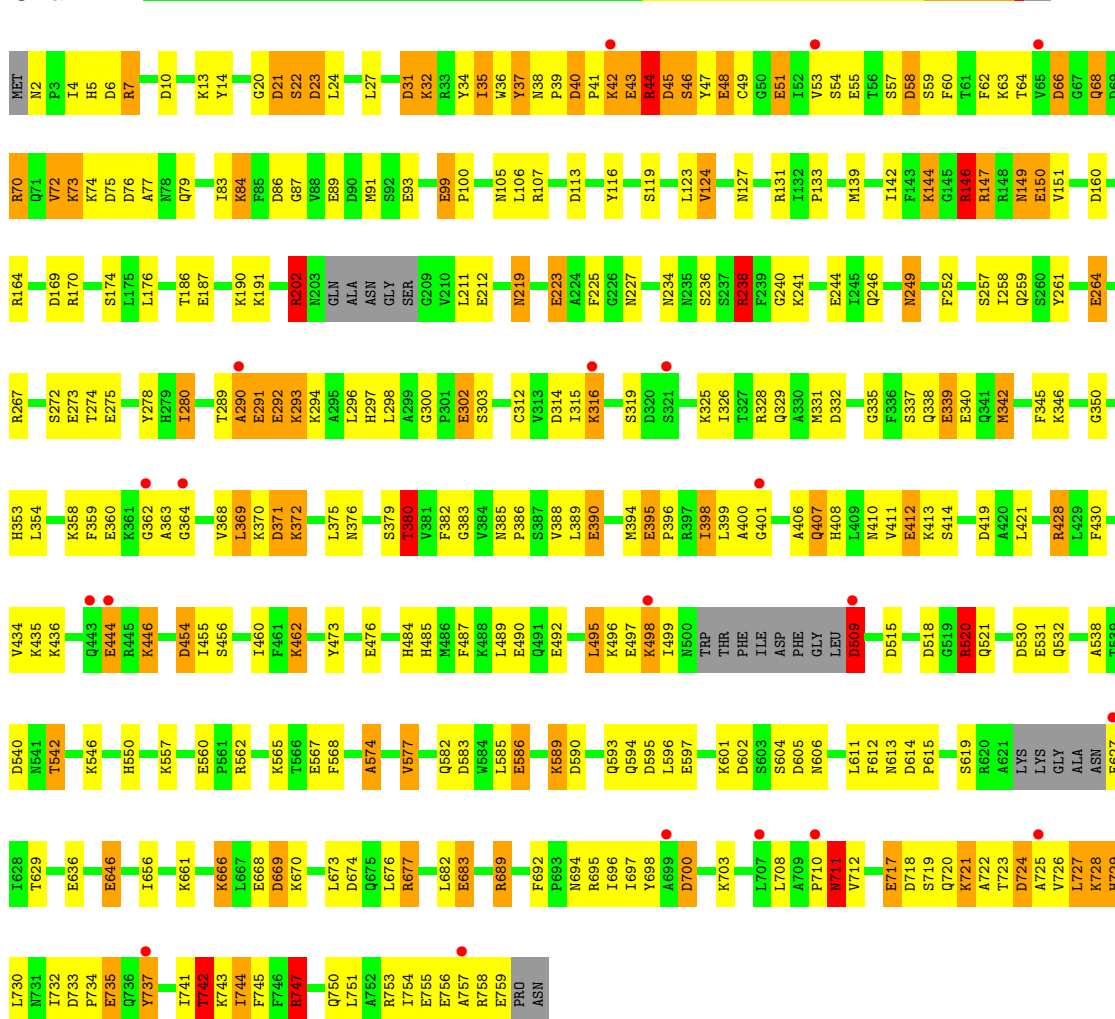
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	730	Total	O	0	0
			730	730		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: MYOSIN S1DC MOTOR DOMAIN

Chain A:



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	104.00Å 180.50Å 54.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 2.00 19.67 – 2.00	Depositor EDS
% Data completeness (in resolution range)	95.2 (25.00-2.00) 94.9 (19.67-2.00)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.98 (at 2.01Å)	Xtriage
Refinement program	TNT V. 5-E	Depositor
R, R_{free}	0.187 , (Not available) 0.197 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	19.7	Xtriage
Anisotropy	0.339	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.45 , 114.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 66024 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6643	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 3.82% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MNQ, MG

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	1.14	40/5999 (0.7%)	1.65	103/8097 (1.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	1	0

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	395	GLU	CD-OE2	8.51	1.35	1.25
1	A	93	GLU	CD-OE2	7.99	1.34	1.25
1	A	490	GLU	CD-OE2	7.98	1.34	1.25
1	A	531	GLU	CD-OE2	7.64	1.34	1.25
1	A	560	GLU	CD-OE1	-7.53	1.17	1.25
1	A	668	GLU	CD-OE2	7.52	1.33	1.25
1	A	187	GLU	CD-OE2	7.29	1.33	1.25
1	A	264	GLU	CD-OE1	-7.27	1.17	1.25
1	A	412	GLU	CD-OE2	7.17	1.33	1.25
1	A	444	GLU	CD-OE2	7.17	1.33	1.25
1	A	275	GLU	CD-OE2	7.09	1.33	1.25
1	A	150	GLU	CD-OE2	7.09	1.33	1.25
1	A	755	GLU	CD-OE2	6.96	1.33	1.25
1	A	717	GLU	CD-OE2	6.83	1.33	1.25
1	A	683	GLU	CD-OE2	6.79	1.33	1.25
1	A	646	GLU	CD-OE2	6.69	1.33	1.25
1	A	244	GLU	CD-OE2	6.61	1.32	1.25
1	A	340	GLU	CD-OE2	6.58	1.32	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	492	GLU	CD-OE2	6.57	1.32	1.25
1	A	756	GLU	CD-OE2	6.53	1.32	1.25
1	A	735	GLU	CD-OE2	6.49	1.32	1.25
1	A	390	GLU	CD-OE2	6.45	1.32	1.25
1	A	476	GLU	CD-OE2	6.36	1.32	1.25
1	A	292	GLU	CD-OE2	6.23	1.32	1.25
1	A	291	GLU	CD-OE2	6.22	1.32	1.25
1	A	89	GLU	CD-OE2	6.19	1.32	1.25
1	A	55	GLU	CD-OE2	6.17	1.32	1.25
1	A	636	GLU	CD-OE2	6.15	1.32	1.25
1	A	212	GLU	CD-OE2	6.05	1.32	1.25
1	A	360	GLU	CD-OE2	6.03	1.32	1.25
1	A	586	GLU	CD-OE2	6.02	1.32	1.25
1	A	174	SER	CB-OG	5.71	1.49	1.42
1	A	476	GLU	CD-OE1	-5.65	1.19	1.25
1	A	51	GLU	CD-OE2	5.54	1.31	1.25
1	A	302	GLU	CD-OE2	5.45	1.31	1.25
1	A	43	GLU	CD-OE2	5.32	1.31	1.25
1	A	567	GLU	CD-OE2	5.28	1.31	1.25
1	A	223	GLU	CD-OE2	5.26	1.31	1.25
1	A	339	GLU	CD-OE2	5.13	1.31	1.25
1	A	48	GLU	CD-OE2	5.05	1.31	1.25

All (103) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	238[A]	ARG	NE-CZ-NH2	-17.07	111.76	120.30
1	A	238[A]	ARG	NE-CZ-NH1	15.85	128.23	120.30
1	A	44[A]	ARG	NE-CZ-NH1	13.43	127.02	120.30
1	A	677[A]	ARG	NE-CZ-NH1	13.14	126.87	120.30
1	A	747[A]	ARG	NE-CZ-NH1	-10.93	114.83	120.30
1	A	23	ASP	CB-CG-OD2	-10.17	109.15	118.30
1	A	6	ASP	CB-CG-OD2	-9.62	109.64	118.30
1	A	700	ASP	CB-CG-OD2	-9.62	109.64	118.30
1	A	23	ASP	CB-CG-OD1	9.24	126.61	118.30
1	A	590	ASP	CB-CG-OD2	-8.99	110.21	118.30
1	A	76	ASP	CB-CG-OD2	-8.62	110.54	118.30
1	A	700	ASP	CB-CG-OD1	8.59	126.03	118.30
1	A	530	ASP	CB-CG-OD1	8.40	125.86	118.30
1	A	21	ASP	CB-CG-OD2	-8.23	110.89	118.30
1	A	202[A]	ARG	NE-CZ-NH1	8.15	124.38	120.30
1	A	520[A]	ARG	NE-CZ-NH1	8.09	124.34	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	160	ASP	CB-CG-OD1	8.03	125.53	118.30
1	A	219	ASN	CB-CA-C	8.01	126.43	110.40
1	A	45	ASP	CB-CG-OD2	-8.01	111.09	118.30
1	A	515	ASP	CB-CG-OD1	8.01	125.50	118.30
1	A	520[A]	ARG	NE-CZ-NH2	-7.87	116.37	120.30
1	A	238[A]	ARG	CD-NE-CZ	7.71	134.39	123.60
1	A	7[A]	ARG	NE-CZ-NH1	7.69	124.14	120.30
1	A	164[A]	ARG	NE-CZ-NH1	7.60	124.10	120.30
1	A	45	ASP	CB-CG-OD1	7.59	125.13	118.30
1	A	169	ASP	CB-CG-OD1	7.58	125.12	118.30
1	A	509	ASP	CB-CG-OD2	-7.52	111.53	118.30
1	A	454	ASP	CB-CG-OD1	7.49	125.04	118.30
1	A	509	ASP	CB-CG-OD1	7.45	125.01	118.30
1	A	689[A]	ARG	NE-CZ-NH1	7.42	124.01	120.30
1	A	419	ASP	CB-CG-OD2	-7.34	111.69	118.30
1	A	58	ASP	CB-CG-OD1	7.33	124.90	118.30
1	A	119	SER	N-CA-CB	7.14	121.22	110.50
1	A	44[A]	ARG	CD-NE-CZ	7.12	133.57	123.60
1	A	518	ASP	CB-CG-OD1	7.07	124.66	118.30
1	A	170[A]	ARG	NE-CZ-NH1	7.07	123.83	120.30
1	A	473	TYR	CB-CG-CD2	-7.07	116.76	121.00
1	A	267[A]	ARG	NE-CZ-NH1	7.02	123.81	120.30
1	A	76	ASP	CB-CG-OD1	6.96	124.56	118.30
1	A	419	ASP	CB-CG-OD1	6.92	124.53	118.30
1	A	605	ASP	CB-CG-OD2	-6.89	112.10	118.30
1	A	747[A]	ARG	CD-NE-CZ	-6.81	114.06	123.60
1	A	314	ASP	CB-CG-OD2	-6.81	112.17	118.30
1	A	86	ASP	CB-CG-OD1	6.80	124.42	118.30
1	A	562[A]	ARG	NE-CZ-NH2	-6.65	116.97	120.30
1	A	6	ASP	CB-CG-OD1	6.65	124.29	118.30
1	A	454	ASP	CB-CG-OD2	-6.65	112.31	118.30
1	A	332	ASP	CB-CG-OD2	-6.64	112.32	118.30
1	A	724	ASP	CB-CG-OD2	-6.64	112.32	118.30
1	A	22	SER	N-CA-CB	6.50	120.25	110.50
1	A	10	ASP	CB-CG-OD2	-6.45	112.50	118.30
1	A	583	ASP	CB-CG-OD1	6.39	124.05	118.30
1	A	674	ASP	CB-CG-OD1	6.38	124.05	118.30
1	A	170[A]	ARG	NE-CZ-NH2	-6.37	117.11	120.30
1	A	37	TYR	CB-CG-CD1	6.37	124.82	121.00
1	A	733	ASP	CB-CG-OD2	-6.34	112.59	118.30
1	A	147[A]	ARG	NE-CZ-NH1	6.31	123.46	120.30
1	A	75	ASP	CB-CG-OD2	-6.31	112.62	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	574	ALA	N-CA-CB	-6.18	101.45	110.10
1	A	58	ASP	CB-CG-OD2	-6.17	112.75	118.30
1	A	249	ASN	CA-CB-CG	-6.11	99.95	113.40
1	A	595	ASP	CB-CG-OD1	6.11	123.80	118.30
1	A	590	ASP	CB-CG-OD1	6.08	123.77	118.30
1	A	124	VAL	CA-CB-CG2	-6.08	101.79	110.90
1	A	146[A]	ARG	NE-CZ-NH1	6.07	123.33	120.30
1	A	66	ASP	CB-CG-OD2	-6.00	112.90	118.30
1	A	695[A]	ARG	NE-CZ-NH1	5.99	123.30	120.30
1	A	41	PRO	N-CA-CB	5.97	110.46	103.30
1	A	677[A]	ARG	NH1-CZ-NH2	-5.92	112.89	119.40
1	A	733	ASP	CB-CG-OD1	5.74	123.47	118.30
1	A	131[A]	ARG	NE-CZ-NH1	5.69	123.15	120.30
1	A	577	VAL	CG1-CB-CG2	-5.67	101.82	110.90
1	A	371	ASP	CB-CG-OD2	-5.66	113.20	118.30
1	A	21	ASP	CB-CG-OD1	5.59	123.33	118.30
1	A	147[A]	ARG	NE-CZ-NH2	-5.57	117.51	120.30
1	A	41	PRO	C-N-CA	5.55	135.59	121.70
1	A	86	ASP	CB-CG-OD2	-5.55	113.31	118.30
1	A	107[A]	ARG	CG-CD-NE	-5.52	100.20	111.80
1	A	113	ASP	CB-CG-OD1	5.51	123.26	118.30
1	A	380	THR	N-CA-CB	5.50	120.74	110.30
1	A	99	GLU	CG-CD-OE1	5.49	129.27	118.30
1	A	669	ASP	CB-CG-OD2	-5.48	113.37	118.30
1	A	99	GLU	CG-CD-OE2	-5.47	107.35	118.30
1	A	202[A]	ARG	NE-CZ-NH2	-5.43	117.59	120.30
1	A	538	ALA	N-CA-CB	5.42	117.68	110.10
1	A	602	ASP	CB-CG-OD2	-5.42	113.43	118.30
1	A	428[A]	ARG	NE-CZ-NH2	-5.41	117.59	120.30
1	A	742	THR	CA-CB-CG2	-5.41	104.83	112.40
1	A	606	ASN	CB-CA-C	5.39	121.17	110.40
1	A	677[A]	ARG	O-C-N	-5.29	114.24	122.70
1	A	629	THR	CA-CB-CG2	-5.29	105.00	112.40
1	A	737	TYR	O-C-N	5.24	131.09	122.70
1	A	44[A]	ARG	NH1-CZ-NH2	-5.20	113.68	119.40
1	A	518	ASP	CB-CG-OD2	-5.14	113.67	118.30
1	A	113	ASP	CB-CG-OD2	-5.14	113.67	118.30
1	A	332	ASP	CB-CG-OD1	5.13	122.92	118.30
1	A	568	PHE	CZ-CE2-CD2	-5.10	113.98	120.10
1	A	395	GLU	CG-CD-OE2	-5.06	108.18	118.30
1	A	530	ASP	CB-CG-OD2	-5.06	113.75	118.30
1	A	44[A]	ARG	CG-CD-NE	5.05	122.41	111.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	729	HIS	CA-CB-CG	-5.05	105.02	113.60
1	A	712	VAL	N-CA-C	5.05	124.63	111.00
1	A	302	GLU	CG-CD-OE2	-5.01	108.28	118.30

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	606	ASN	CA

There are no planarity outliers.

5.2 Close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5887	0	5597	272	0
2	A	1	0	0	0	0
3	A	25	0	9	2	0
4	A	730	0	0	39	0
All	All	6643	0	5606	275	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 24.

All (275) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:4:ILE:HD11	1:A:142:ILE:HG23	1.24	1.17
1:A:45:ASP:HB2	1:A:670:LYS:HD3	1.17	1.15
1:A:45:ASP:HB2	1:A:670:LYS:CD	1.76	1.13
1:A:35:ILE:HD11	1:A:77:ALA:HB1	1.24	1.09
1:A:398:ILE:HG12	1:A:407:GLN:HG3	1.31	1.08
1:A:40:ASP:HB3	1:A:42:LYS:HB2	1.45	0.98
1:A:485:HIS:NE2	1:A:489:LEU:HD11	1.84	0.93
1:A:62:PHE:HE2	1:A:72:VAL:HG22	1.34	0.93
1:A:35:ILE:HD11	1:A:77:ALA:CB	1.98	0.92

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:45:ASP:CG	1:A:677[A]:ARG:HH22	1.73	0.91
1:A:398:ILE:HG12	1:A:407:GLN:CG	2.01	0.91
1:A:697:ILE:HD13	1:A:743:LYS:HG2	1.53	0.90
1:A:45:ASP:CB	1:A:670:LYS:HD3	2.01	0.89
1:A:296:LEU:HB2	1:A:298:LEU:HD11	1.56	0.88
1:A:342:MET:HG3	1:A:346:LYS:HE3	1.55	0.87
1:A:127:ASN:HD21	3:A:999:MNQ:HA21	1.37	0.86
1:A:4:ILE:HD11	1:A:142:ILE:CG2	2.05	0.86
1:A:741:ILE:HG22	1:A:742:THR:HG22	1.58	0.84
1:A:53:VAL:HG11	1:A:63:LYS:HD2	1.58	0.84
1:A:249:ASN:ND2	4:A:1086:HOH:O	2.04	0.82
1:A:149:ASN:HD22	1:A:150:GLU:N	1.78	0.82
1:A:2:ASN:HB3	1:A:5:HIS:HD2	1.45	0.81
1:A:385:ASN:HB3	1:A:388:VAL:CG2	2.10	0.81
1:A:64:THR:HG23	1:A:68:GLN:O	1.81	0.81
1:A:296:LEU:HB2	1:A:298:LEU:CD1	2.11	0.80
1:A:249:ASN:H	1:A:249:ASN:ND2	1.79	0.77
1:A:280:ILE:O	1:A:280:ILE:HG13	1.85	0.76
1:A:698:TYR:CZ	1:A:720:GLN:HG3	2.21	0.75
1:A:698:TYR:CE1	1:A:720:GLN:HG3	2.22	0.74
1:A:289:THR:HG23	1:A:292:GLU:OE2	1.87	0.74
1:A:99:GLU:OE2	4:A:1670:HOH:O	2.05	0.74
1:A:62:PHE:CE2	1:A:72:VAL:HG22	2.19	0.74
1:A:149:ASN:C	1:A:149:ASN:HD22	1.90	0.73
1:A:399:LEU:HD12	1:A:400:ALA:N	2.03	0.73
1:A:342:MET:HE3	1:A:342:MET:HA	1.70	0.73
1:A:385:ASN:HB3	1:A:388:VAL:HG23	1.71	0.72
1:A:40:ASP:C	1:A:42:LYS:H	1.90	0.72
1:A:219:ASN:HB3	4:A:1724:HOH:O	1.89	0.72
1:A:45:ASP:HB2	1:A:670:LYS:HD2	1.67	0.72
1:A:498:LYS:HG2	1:A:498:LYS:O	1.88	0.72
1:A:521:GLN:OE1	4:A:1657:HOH:O	2.08	0.72
1:A:40:ASP:CB	1:A:42:LYS:HB2	2.19	0.71
1:A:495:LEU:HD12	4:A:1654:HOH:O	1.90	0.71
1:A:509:ASP:OD2	1:A:557:LYS:NZ	2.22	0.71
1:A:290:ALA:HA	1:A:293:LYS:HD2	1.73	0.70
1:A:342:MET:HG3	1:A:346:LYS:CE	2.20	0.70
1:A:2:ASN:HB3	1:A:5:HIS:CD2	2.27	0.70
1:A:718:ASP:CG	1:A:721:LYS:HB2	2.12	0.70
1:A:87:GLY:H	1:A:105:ASN:ND2	1.89	0.69
1:A:259:GLN:HG2	1:A:261:TYR:CZ	2.28	0.68
1:A:372:LYS:O	1:A:376:ASN:ND2	2.27	0.68

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:273:GLU:O	1:A:274:THR:OG1	2.12	0.68
1:A:249:ASN:HD22	1:A:249:ASN:H	1.41	0.67
1:A:574:ALA:HA	4:A:1583:HOH:O	1.93	0.67
1:A:710:PRO:O	1:A:711:ASN:ND2	2.27	0.67
1:A:302:GLU:H	1:A:302:GLU:CD	1.97	0.67
1:A:697:ILE:HD13	1:A:743:LYS:CG	2.23	0.67
1:A:497:GLU:OE1	1:A:742:THR:HG23	1.95	0.67
1:A:542:THR:HG23	4:A:1661:HOH:O	1.94	0.67
1:A:219:ASN:ND2	4:A:1274:HOH:O	2.27	0.66
1:A:62:PHE:HE2	1:A:72:VAL:CG2	2.06	0.66
1:A:289:THR:O	1:A:291:GLU:N	2.27	0.66
1:A:697:ILE:HB	1:A:700:ASP:OD1	1.96	0.66
1:A:225:PHE:CZ	1:A:280:ILE:HD13	2.31	0.66
1:A:238[A]:ARG:HD3	1:A:264:GLU:OE2	1.95	0.65
1:A:395:GLU:HA	1:A:407:GLN:O	1.96	0.65
1:A:42:LYS:O	1:A:43:GLU:HG3	1.96	0.65
1:A:31:ASP:OD2	1:A:31:ASP:N	2.21	0.65
1:A:485:HIS:NE2	1:A:489:LEU:CD1	2.58	0.65
1:A:710:PRO:HD3	1:A:729:HIS:ND1	2.11	0.65
1:A:296:LEU:HB2	1:A:298:LEU:CG	2.28	0.64
1:A:139:MET:HA	1:A:142:ILE:HD12	1.81	0.63
1:A:399:LEU:HD11	1:A:401:GLY:C	2.19	0.63
1:A:435:LYS:HG3	4:A:1505:HOH:O	1.99	0.63
1:A:241:LYS:NZ	1:A:454:ASP:OD1	2.29	0.63
1:A:710:PRO:HD3	1:A:729:HIS:CG	2.34	0.62
1:A:646:GLU:HG3	4:A:1653:HOH:O	1.98	0.62
1:A:290:ALA:HA	1:A:293:LYS:CD	2.30	0.62
1:A:62:PHE:CE2	1:A:72:VAL:CG2	2.82	0.62
1:A:34:TYR:CE1	1:A:51:GLU:HB2	2.36	0.61
1:A:399:LEU:HD11	1:A:401:GLY:O	2.00	0.61
1:A:710:PRO:HG3	1:A:729:HIS:CE1	2.35	0.61
1:A:718:ASP:OD1	1:A:721:LYS:HB2	2.01	0.61
1:A:520[A]:ARG:NE	4:A:1702:HOH:O	2.34	0.61
1:A:399:LEU:HD12	1:A:400:ALA:H	1.66	0.60
1:A:385:ASN:OD1	1:A:386:PRO:HD2	2.02	0.60
1:A:40:ASP:OD2	1:A:42:LYS:HB2	2.02	0.59
1:A:176:LEU:HD12	1:A:176:LEU:N	2.16	0.59
1:A:328[A]:ARG:HD2	4:A:1207:HOH:O	2.02	0.59
1:A:45:ASP:H	1:A:670:LYS:HE2	1.67	0.59
1:A:45:ASP:CB	1:A:670:LYS:CD	2.68	0.59
1:A:312:CYS:HB2	4:A:1638:HOH:O	2.02	0.59
1:A:542:THR:CB	4:A:1661:HOH:O	2.50	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:296:LEU:CB	1:A:298:LEU:HD11	2.29	0.59
1:A:722:ALA:O	1:A:725:ALA:HB3	2.02	0.59
1:A:4:ILE:CD1	1:A:142:ILE:HG23	2.16	0.59
1:A:289:THR:O	1:A:292:GLU:N	2.36	0.58
1:A:296:LEU:O	1:A:297:HIS:HB2	2.04	0.58
1:A:410:ASN:OD1	1:A:410:ASN:C	2.42	0.58
1:A:290:ALA:CA	1:A:293:LYS:HD2	2.34	0.57
1:A:683:GLU:HG3	4:A:1365:HOH:O	2.04	0.57
1:A:601:LYS:HG2	1:A:613:ASN:ND2	2.19	0.57
1:A:296:LEU:HB2	1:A:298:LEU:HG	1.86	0.57
1:A:710:PRO:HD3	1:A:729:HIS:CE1	2.40	0.57
1:A:741:ILE:HG22	1:A:742:THR:CG2	2.34	0.56
1:A:726:VAL:O	1:A:730:LEU:HG	2.06	0.56
1:A:59:SER:HB2	1:A:72:VAL:O	2.06	0.56
1:A:698:TYR:CZ	1:A:720:GLN:CG	2.89	0.56
1:A:91:MET:CE	1:A:106:LEU:HD13	2.36	0.56
1:A:542:THR:CG2	4:A:1661:HOH:O	2.54	0.55
1:A:371:ASP:OD2	1:A:372:LYS:N	2.39	0.55
1:A:24:LEU:HD23	1:A:24:LEU:N	2.22	0.55
1:A:38:ASN:C	1:A:40:ASP:H	2.10	0.55
1:A:58:ASP:OD2	1:A:58:ASP:N	2.40	0.55
1:A:732:ILE:HD12	1:A:750:GLN:NE2	2.22	0.54
1:A:396:PRO:HG2	1:A:398:ILE:HD11	1.88	0.54
1:A:79:GLN:HB2	4:A:1089:HOH:O	2.06	0.54
1:A:540:ASP:OD2	4:A:1511:HOH:O	2.19	0.54
1:A:289:THR:C	1:A:291:GLU:N	2.62	0.53
1:A:582:GLN:HG3	4:A:1325:HOH:O	2.07	0.53
1:A:697:ILE:O	1:A:700:ASP:HB2	2.08	0.53
1:A:124:VAL:HG13	1:A:656:ILE:HD12	1.90	0.53
1:A:293:LYS:HA	1:A:298:LEU:HD12	1.91	0.53
1:A:51:GLU:O	1:A:53:VAL:HG13	2.09	0.52
1:A:149:ASN:C	1:A:149:ASN:ND2	2.59	0.52
1:A:297:HIS:HB2	1:A:353:HIS:HE2	1.75	0.51
1:A:191:LYS:NZ	4:A:1424:HOH:O	2.31	0.51
1:A:45:ASP:HB3	1:A:673:LEU:HD12	1.92	0.51
1:A:190:LYS:HE2	1:A:223:GLU:OE2	2.11	0.51
1:A:362:GLY:O	1:A:364:GLY:N	2.44	0.51
1:A:37:TYR:O	1:A:47:TYR:HA	2.11	0.51
1:A:127:ASN:ND2	3:A:999:MNQ:HA21	2.18	0.50
1:A:79:GLN:HG3	4:A:1260:HOH:O	2.10	0.50
1:A:619:SER:HB3	1:A:627:PHE:CE2	2.45	0.50
1:A:290:ALA:N	1:A:293:LYS:HD2	2.27	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:359:PHE:HB3	1:A:411:VAL:HG22	1.94	0.50
1:A:383:GLY:O	1:A:604:SER:N	2.37	0.50
1:A:289:THR:C	1:A:291:GLU:H	2.14	0.50
1:A:737:TYR:HB2	1:A:744:ILE:HD11	1.94	0.50
1:A:297:HIS:HB2	1:A:353:HIS:NE2	2.27	0.49
1:A:87:GLY:H	1:A:105:ASN:HD21	1.57	0.49
1:A:708:LEU:HD23	1:A:759:GLU:HA	1.93	0.49
1:A:45:ASP:H	1:A:670:LYS:CE	2.25	0.49
1:A:498:LYS:C	1:A:499:ILE:HG13	2.33	0.49
1:A:593:GLN:HB2	1:A:596:LEU:HD12	1.94	0.49
1:A:219:ASN:CB	4:A:1724:HOH:O	2.54	0.49
1:A:734:PRO:HA	1:A:737:TYR:CE2	2.48	0.49
1:A:498:LYS:O	1:A:499:ILE:HG13	2.12	0.49
1:A:550:HIS:HD2	4:A:1303:HOH:O	1.95	0.49
1:A:532:GLN:OE1	1:A:542:THR:HB	2.12	0.48
1:A:34:TYR:HB3	1:A:49:CYS:SG	2.53	0.48
1:A:40:ASP:OD2	1:A:42:LYS:HG3	2.15	0.47
1:A:149:ASN:ND2	1:A:150:GLU:HG3	2.29	0.47
1:A:227:ASN:HA	1:A:236:SER:O	2.14	0.47
1:A:354:LEU:HD13	1:A:421:LEU:HD23	1.97	0.47
1:A:331:MET:HE3	1:A:345:PHE:HZ	1.79	0.47
1:A:342:MET:CG	1:A:346:LYS:CE	2.93	0.47
1:A:40:ASP:C	1:A:42:LYS:N	2.64	0.46
1:A:38:ASN:ND2	1:A:46:SER:O	2.35	0.46
1:A:350:GLY:HA3	1:A:382:PHE:CZ	2.51	0.46
1:A:32:LYS:HB2	1:A:34:TYR:CE2	2.51	0.46
1:A:335:GLY:HA3	4:A:1640:HOH:O	2.16	0.46
1:A:35:ILE:HD13	4:A:1087:HOH:O	2.15	0.45
1:A:40:ASP:HB3	1:A:42:LYS:CB	2.30	0.45
1:A:359:PHE:CB	1:A:411:VAL:HG22	2.46	0.45
1:A:585:LEU:O	1:A:589:LYS:HD2	2.16	0.45
1:A:73:LYS:HA	1:A:73:LYS:HD2	1.44	0.45
1:A:39:PRO:HG3	1:A:48:GLU:HG3	1.98	0.45
1:A:99:GLU:HB2	1:A:100:PRO:HD3	1.97	0.45
1:A:462:LYS:HD2	1:A:462:LYS:HA	1.57	0.45
1:A:83:ILE:HD11	4:A:1044:HOH:O	2.16	0.45
1:A:249:ASN:N	1:A:249:ASN:ND2	2.49	0.45
1:A:698:TYR:CE2	1:A:720:GLN:HG3	2.51	0.45
1:A:614:ASP:OD1	1:A:615:PRO:HD2	2.17	0.45
1:A:455:ILE:O	1:A:455:ILE:HG13	2.15	0.45
1:A:338:GLN:NE2	4:A:1289:HOH:O	2.50	0.45
1:A:14:TYR:CE2	1:A:133:PRO:HG2	2.52	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:331:MET:CE	1:A:345:PHE:HZ	2.31	0.44
1:A:484:HIS:O	1:A:487:PHE:HB3	2.17	0.44
1:A:38:ASN:C	1:A:40:ASP:N	2.71	0.44
1:A:720:GLN:O	1:A:723:THR:HB	2.18	0.44
1:A:84:LYS:HB3	4:A:1261:HOH:O	2.17	0.44
1:A:462:LYS:HG2	4:A:1649:HOH:O	2.18	0.44
1:A:316:LYS:HB2	1:A:316:LYS:HE3	1.66	0.44
1:A:22:SER:OG	1:A:23:ASP:N	2.51	0.44
1:A:223:GLU:O	1:A:227:ASN:HB2	2.18	0.44
1:A:696:ILE:O	1:A:744:ILE:N	2.44	0.43
1:A:144:LYS:HD3	4:A:1619:HOH:O	2.16	0.43
1:A:293:LYS:HG3	1:A:293:LYS:H	1.63	0.43
1:A:293:LYS:O	1:A:297:HIS:N	2.51	0.43
1:A:589:LYS:NZ	4:A:1255:HOH:O	2.50	0.43
1:A:259:GLN:HG2	1:A:261:TYR:OH	2.17	0.43
1:A:656:ILE:HD13	1:A:676:LEU:HD21	2.01	0.43
1:A:240:GLY:HA3	1:A:455:ILE:HG22	1.99	0.43
1:A:724:ASP:O	1:A:728:LYS:HB2	2.18	0.43
1:A:698:TYR:HB3	1:A:719:SER:HB3	2.00	0.43
1:A:53:VAL:HG22	1:A:62:PHE:HA	2.01	0.43
1:A:331:MET:HE3	1:A:345:PHE:CZ	2.54	0.43
1:A:727:LEU:HD12	1:A:727:LEU:HA	1.69	0.43
1:A:116:TYR:HB3	1:A:123:LEU:HD11	2.01	0.43
1:A:296:LEU:HA	1:A:296:LEU:HD23	1.89	0.43
1:A:300:GLY:HA3	1:A:302:GLU:OE2	2.18	0.43
1:A:246:GLN:HG2	1:A:446:LYS:HD3	2.00	0.43
1:A:44[A]:ARG:HD3	1:A:45:ASP:OD2	2.19	0.42
1:A:60:PHE:CE1	1:A:74:LYS:HA	2.54	0.42
1:A:754:ILE:O	1:A:757:ALA:HB3	2.19	0.42
1:A:586:GLU:HB2	4:A:1514:HOH:O	2.19	0.42
1:A:597:GLU:OE1	1:A:612:PHE:CD2	2.73	0.42
1:A:410:ASN:OD1	1:A:413:LYS:N	2.46	0.42
1:A:546:LYS:NZ	4:A:1450:HOH:O	2.30	0.42
1:A:331:MET:CE	1:A:345:PHE:CZ	3.03	0.42
1:A:289:THR:O	1:A:289:THR:OG1	2.37	0.42
1:A:408:HIS:ND1	1:A:408:HIS:C	2.73	0.42
1:A:593:GLN:NE2	4:A:1572:HOH:O	2.42	0.41
1:A:369:LEU:HA	1:A:369:LEU:HD12	1.78	0.41
1:A:708:LEU:CD2	1:A:758:ARG:O	2.68	0.41
1:A:40:ASP:CG	1:A:42:LYS:HB2	2.39	0.41
1:A:485:HIS:CD2	1:A:489:LEU:CD1	3.03	0.41
1:A:376:ASN:O	1:A:380:THR:HG22	2.21	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:258:ILE:HD13	1:A:258:ILE:HG21	1.82	0.41
1:A:614:ASP:OD1	1:A:614:ASP:C	2.58	0.41
1:A:60:PHE:N	1:A:72:VAL:O	2.49	0.41
1:A:23:ASP:N	1:A:23:ASP:OD1	2.52	0.41
1:A:186:THR:HG22	1:A:190:LYS:HE3	2.03	0.41
1:A:362:GLY:C	1:A:364:GLY:N	2.74	0.41
1:A:428[A]:ARG:HD3	1:A:611:LEU:O	2.21	0.41
1:A:698:TYR:CD2	1:A:720:GLN:HA	2.56	0.41
1:A:146[A]:ARG:HD3	1:A:151:VAL:CG1	2.51	0.41
1:A:146[A]:ARG:HD3	1:A:151:VAL:HG13	2.01	0.41
1:A:747[A]:ARG:HH11	1:A:747[A]:ARG:HD3	1.54	0.41
1:A:430:PHE:O	1:A:434:VAL:HG23	2.21	0.41
1:A:698:TYR:CD1	1:A:720:GLN:HG3	2.56	0.40
1:A:99:GLU:N	1:A:100:PRO:HD2	2.36	0.40
1:A:21:ASP:OD2	1:A:21:ASP:N	2.54	0.40
1:A:375:LEU:HD12	1:A:389:LEU:HD23	2.03	0.40
1:A:460:ILE:HG12	1:A:577:VAL:HG22	2.02	0.40
1:A:234:ASN:OD1	1:A:315:ILE:HG23	2.20	0.40
1:A:20:GLY:N	4:A:1385:HOH:O	2.27	0.40
1:A:142:ILE:CG2	1:A:142:ILE:O	2.69	0.40
1:A:34:TYR:CD1	1:A:51:GLU:HB2	2.56	0.40
1:A:386:PRO:O	1:A:390:GLU:HB2	2.22	0.40
1:A:372:LYS:HB2	1:A:372:LYS:HE3	1.52	0.40
1:A:710:PRO:CG	1:A:729:HIS:CE1	3.05	0.40
1:A:39:PRO:HG3	1:A:48:GLU:CG	2.52	0.40
1:A:496:LYS:HB3	1:A:496:LYS:NZ	2.34	0.40
1:A:666:LYS:HE2	4:A:1217:HOH:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

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	732/761 (96%)	698 (95%)	29 (4%)	5 (1%)	30	20

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	44[A]	ARG
1	A	290	ALA
1	A	363	ALA
1	A	711	ASN
1	A	7[A]	ARG

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	632/665 (95%)	555 (88%)	77 (12%)	7 4

All (77) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	LYS
1	A	27	LEU
1	A	31	ASP
1	A	32	LYS
1	A	35	ILE
1	A	40	ASP
1	A	42	LYS
1	A	44[A]	ARG
1	A	46	SER
1	A	54	SER
1	A	57	SER
1	A	66	ASP
1	A	68	GLN
1	A	70[A]	ARG
1	A	72	VAL
1	A	73	LYS
1	A	84	LYS
1	A	144	LYS
1	A	146[A]	ARG
1	A	147[A]	ARG
1	A	149	ASN
1	A	202[A]	ARG

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Mol	Chain	Res	Type
1	A	211	LEU
1	A	238[A]	ARG
1	A	257	SER
1	A	272	SER
1	A	280	ILE
1	A	293	LYS
1	A	294	LYS
1	A	303	SER
1	A	316	LYS
1	A	319	SER
1	A	325	LYS
1	A	326	ILE
1	A	329	GLN
1	A	337	SER
1	A	339	GLU
1	A	342	MET
1	A	358	LYS
1	A	368	VAL
1	A	369	LEU
1	A	370	LYS
1	A	372	LYS
1	A	379	SER
1	A	380	THR
1	A	394	MET
1	A	398	ILE
1	A	407	GLN
1	A	412	GLU
1	A	414	SER
1	A	436	LYS
1	A	444	GLU
1	A	446	LYS
1	A	456	SER
1	A	462	LYS
1	A	495	LEU
1	A	498	LYS
1	A	509	ASP
1	A	520[A]	ARG
1	A	542	THR
1	A	565	LYS
1	A	589	LYS
1	A	594	GLN
1	A	661	LYS

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Mol	Chain	Res	Type
1	A	666	LYS
1	A	669	ASP
1	A	689[A]	ARG
1	A	703	LYS
1	A	711	ASN
1	A	717	GLU
1	A	721	LYS
1	A	727	LEU
1	A	728	LYS
1	A	742	THR
1	A	744	ILE
1	A	747[A]	ARG
1	A	751	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	HIS
1	A	68	GLN
1	A	105	ASN
1	A	149	ASN
1	A	219	ASN
1	A	249	ASN
1	A	259	GLN
1	A	283	GLN
1	A	329	GLN
1	A	376	ASN
1	A	439	ASN
1	A	479	GLN
1	A	491	GLN
1	A	521	GLN
1	A	594	GLN
1	A	616	ASN
1	A	662	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 for Mogul analysis.

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	MNQ	A	999	2	24,25,25	2.16	6 (25%)	30,37,37	3.12	10 (33%)

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
3	MNQ	A	999	2	-	0/18/24/24	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	999	MNQ	O3B-N3	6.16	1.33	1.23
3	A	999	MNQ	PB-OB3	4.78	1.61	1.51
3	A	999	MNQ	C5-C6	3.56	1.47	1.39
3	A	999	MNQ	C4-C3	2.66	1.44	1.38
3	A	999	MNQ	C6-C1	2.44	1.43	1.39
3	A	999	MNQ	CA2-NA3	2.24	1.51	1.46

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	999	MNQ	C4-C3-N3	9.75	127.98	118.96
3	A	999	MNQ	C2-C3-N3	-9.13	111.49	118.71
3	A	999	MNQ	CA2-NA3-C1	-7.23	104.19	122.77
3	A	999	MNQ	C2-C1-NA3	-3.57	114.00	120.80
3	A	999	MNQ	C3-C2-C1	3.35	122.13	118.42
3	A	999	MNQ	O3A-N3-C3	2.77	119.74	114.42
3	A	999	MNQ	C5-C6-C1	-2.42	116.61	119.73
3	A	999	MNQ	C6-C1-NA3	2.23	125.57	121.02
3	A	999	MNQ	OA3-PB-OB3	-2.22	103.47	108.73
3	A	999	MNQ	CA1-CA2-NA3	2.14	116.31	111.09

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	740/761 (97%)	-0.24	20 (2%) 52 52	12, 32, 74, 100	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	627	PHE	6.3
1	A	65	VAL	5.5
1	A	42	LYS	5.4
1	A	362	GLY	4.1
1	A	710	PRO	4.1
1	A	737	TYR	4.0
1	A	707	LEU	3.9
1	A	364	GLY	3.3
1	A	321	SER	3.2
1	A	290	ALA	3.2
1	A	401	GLY	2.9
1	A	725	ALA	2.8
1	A	53	VAL	2.8
1	A	443	GLN	2.7
1	A	444	GLU	2.5
1	A	316	LYS	2.4
1	A	498	LYS	2.3
1	A	509	ASP	2.2
1	A	699	ALA	2.1
1	A	757	ALA	2.1

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

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

6.3 Carbohydrates

There are no carbohydrates in this entry.

6.4 Ligands

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	MNQ	A	999	25/25	0.09	0.82	12,22,98,100	0
2	MG	A	998	1/1	0.03	-2.63	20,20,20,20	0

6.5 Other polymers

There are no such residues in this entry.