



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 01:38 AM UTC

PDB ID : 5ZNY / pdb_00005zny
Title : Structure of mDR3_DD-C363G with MBP tag
Authors : Yin, X.; Jin, T.
Deposited on : 2018-04-11
Resolution : 2.74 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

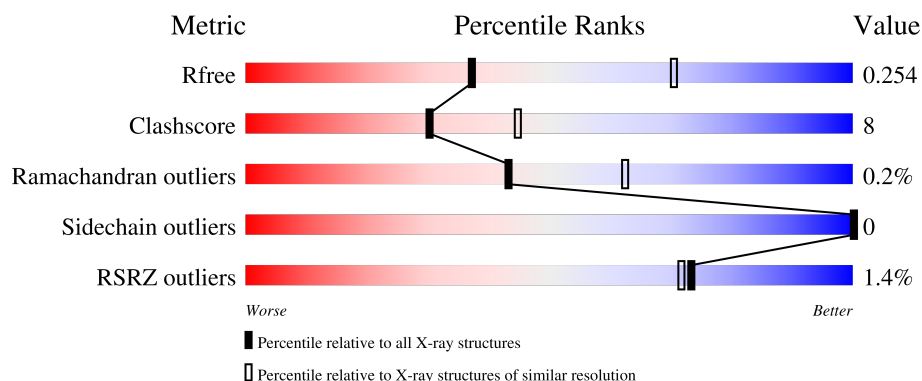
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

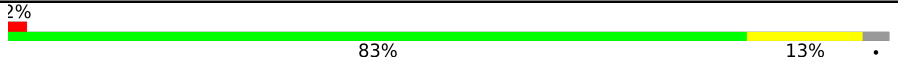
The reported resolution of this entry is 2.74 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1819 (2.76-2.72)
Clashscore	190562	1866 (2.76-2.72)
Ramachandran outliers	187476	1830 (2.76-2.72)
Sidechain outliers	187428	1831 (2.76-2.72)
RSRZ outliers	180081	1819 (2.76-2.72)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	464	
1	B	464	
1	C	464	
1	D	464	

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
2	SO4	D	505	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 14098 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 Maltose-binding periplasmic protein,Tumor necrosis factor receptor superfamily, member 25.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	448	Total	C	N	O	S	0	0	0
			3470	2225	579	656	10			
1	B	450	Total	C	N	O	S	0	0	0
			3487	2234	584	659	10			
1	C	450	Total	C	N	O	S	0	0	0
			3487	2234	584	659	10			
1	D	453	Total	C	N	O	S	0	0	0
			3515	2252	591	662	10			

There are 100 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP P0AEX9
A	83	ALA	ASP	engineered mutation	UNP P0AEX9
A	84	ALA	LYS	engineered mutation	UNP P0AEX9
A	173	ALA	GLU	engineered mutation	UNP P0AEX9
A	174	ALA	ASN	engineered mutation	UNP P0AEX9
A	240	ALA	LYS	engineered mutation	UNP P0AEX9
A	360	ALA	GLU	engineered mutation	UNP P0AEX9
A	363	ALA	LYS	engineered mutation	UNP P0AEX9
A	364	ALA	ASP	engineered mutation	UNP P0AEX9
A	368	ASN	-	linker	UNP P0AEX9
A	369	ALA	-	linker	UNP P0AEX9
A	370	ALA	-	linker	UNP P0AEX9
A	371	ARG	-	linker	UNP P0AEX9
A	372	ALA	-	linker	UNP P0AEX9
A	373	ALA	-	linker	UNP P0AEX9
A	374	ALA	-	linker	UNP P0AEX9
A	410	GLY	CYS	engineered mutation	UNP B1AWN9
A	457	LEU	-	expression tag	UNP B1AWN9
A	458	GLU	-	expression tag	UNP B1AWN9
A	459	HIS	-	expression tag	UNP B1AWN9

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Chain	Residue	Modelled	Actual	Comment	Reference
A	460	HIS	-	expression tag	UNP B1AWN9
A	461	HIS	-	expression tag	UNP B1AWN9
A	462	HIS	-	expression tag	UNP B1AWN9
A	463	HIS	-	expression tag	UNP B1AWN9
A	464	HIS	-	expression tag	UNP B1AWN9
B	1	MET	-	expression tag	UNP P0AEX9
B	83	ALA	ASP	engineered mutation	UNP P0AEX9
B	84	ALA	LYS	engineered mutation	UNP P0AEX9
B	173	ALA	GLU	engineered mutation	UNP P0AEX9
B	174	ALA	ASN	engineered mutation	UNP P0AEX9
B	240	ALA	LYS	engineered mutation	UNP P0AEX9
B	360	ALA	GLU	engineered mutation	UNP P0AEX9
B	363	ALA	LYS	engineered mutation	UNP P0AEX9
B	364	ALA	ASP	engineered mutation	UNP P0AEX9
B	368	ASN	-	linker	UNP P0AEX9
B	369	ALA	-	linker	UNP P0AEX9
B	370	ALA	-	linker	UNP P0AEX9
B	371	ARG	-	linker	UNP P0AEX9
B	372	ALA	-	linker	UNP P0AEX9
B	373	ALA	-	linker	UNP P0AEX9
B	374	ALA	-	linker	UNP P0AEX9
B	410	GLY	CYS	engineered mutation	UNP B1AWN9
B	457	LEU	-	expression tag	UNP B1AWN9
B	458	GLU	-	expression tag	UNP B1AWN9
B	459	HIS	-	expression tag	UNP B1AWN9
B	460	HIS	-	expression tag	UNP B1AWN9
B	461	HIS	-	expression tag	UNP B1AWN9
B	462	HIS	-	expression tag	UNP B1AWN9
B	463	HIS	-	expression tag	UNP B1AWN9
B	464	HIS	-	expression tag	UNP B1AWN9
C	1	MET	-	expression tag	UNP P0AEX9
C	83	ALA	ASP	engineered mutation	UNP P0AEX9
C	84	ALA	LYS	engineered mutation	UNP P0AEX9
C	173	ALA	GLU	engineered mutation	UNP P0AEX9
C	174	ALA	ASN	engineered mutation	UNP P0AEX9
C	240	ALA	LYS	engineered mutation	UNP P0AEX9
C	360	ALA	GLU	engineered mutation	UNP P0AEX9
C	363	ALA	LYS	engineered mutation	UNP P0AEX9
C	364	ALA	ASP	engineered mutation	UNP P0AEX9
C	368	ASN	-	linker	UNP P0AEX9
C	369	ALA	-	linker	UNP P0AEX9
C	370	ALA	-	linker	UNP P0AEX9

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Chain	Residue	Modelled	Actual	Comment	Reference
C	371	ARG	-	linker	UNP P0AEX9
C	372	ALA	-	linker	UNP P0AEX9
C	373	ALA	-	linker	UNP P0AEX9
C	374	ALA	-	linker	UNP P0AEX9
C	410	GLY	CYS	engineered mutation	UNP B1AWN9
C	457	LEU	-	expression tag	UNP B1AWN9
C	458	GLU	-	expression tag	UNP B1AWN9
C	459	HIS	-	expression tag	UNP B1AWN9
C	460	HIS	-	expression tag	UNP B1AWN9
C	461	HIS	-	expression tag	UNP B1AWN9
C	462	HIS	-	expression tag	UNP B1AWN9
C	463	HIS	-	expression tag	UNP B1AWN9
C	464	HIS	-	expression tag	UNP B1AWN9
D	1	MET	-	expression tag	UNP P0AEX9
D	83	ALA	ASP	engineered mutation	UNP P0AEX9
D	84	ALA	LYS	engineered mutation	UNP P0AEX9
D	173	ALA	GLU	engineered mutation	UNP P0AEX9
D	174	ALA	ASN	engineered mutation	UNP P0AEX9
D	240	ALA	LYS	engineered mutation	UNP P0AEX9
D	360	ALA	GLU	engineered mutation	UNP P0AEX9
D	363	ALA	LYS	engineered mutation	UNP P0AEX9
D	364	ALA	ASP	engineered mutation	UNP P0AEX9
D	368	ASN	-	linker	UNP P0AEX9
D	369	ALA	-	linker	UNP P0AEX9
D	370	ALA	-	linker	UNP P0AEX9
D	371	ARG	-	linker	UNP P0AEX9
D	372	ALA	-	linker	UNP P0AEX9
D	373	ALA	-	linker	UNP P0AEX9
D	374	ALA	-	linker	UNP P0AEX9
D	410	GLY	CYS	engineered mutation	UNP B1AWN9
D	457	LEU	-	expression tag	UNP B1AWN9
D	458	GLU	-	expression tag	UNP B1AWN9
D	459	HIS	-	expression tag	UNP B1AWN9
D	460	HIS	-	expression tag	UNP B1AWN9
D	461	HIS	-	expression tag	UNP B1AWN9
D	462	HIS	-	expression tag	UNP B1AWN9
D	463	HIS	-	expression tag	UNP B1AWN9
D	464	HIS	-	expression tag	UNP B1AWN9

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

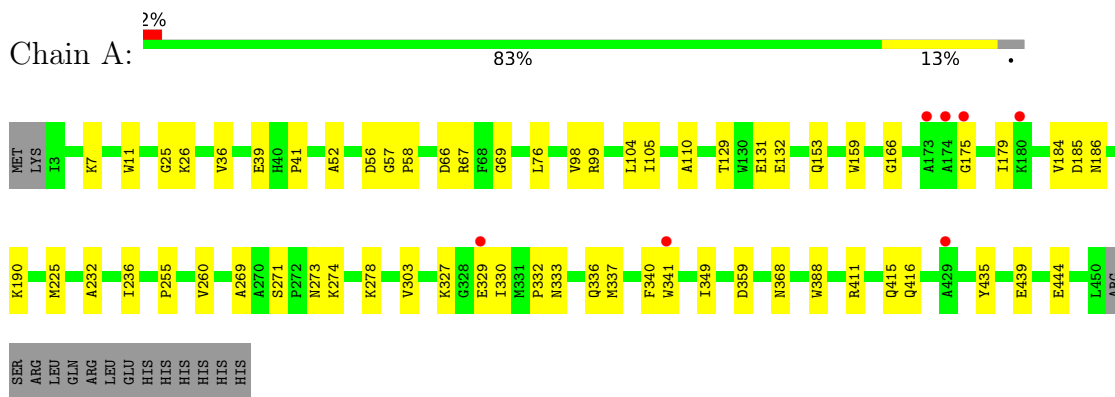
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	O	0	0
			1	1		
3	B	2	Total	O	0	0
			2	2		
3	C	1	Total	O	0	0
			1	1		

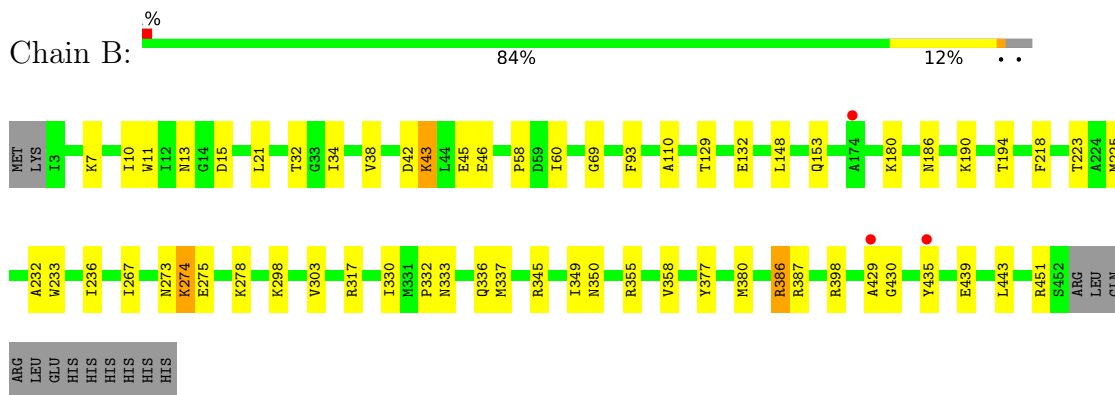
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

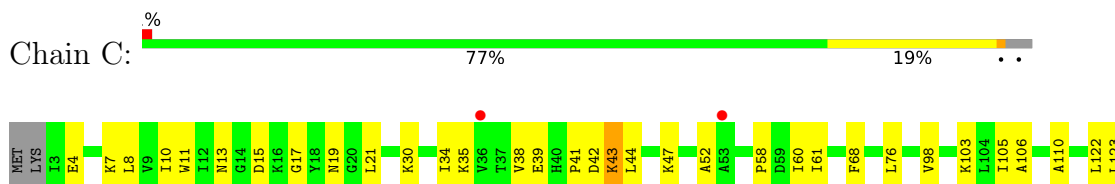
- Molecule 1: Maltose-binding periplasmic protein,Tumor necrosis factor receptor superfamily, member 25

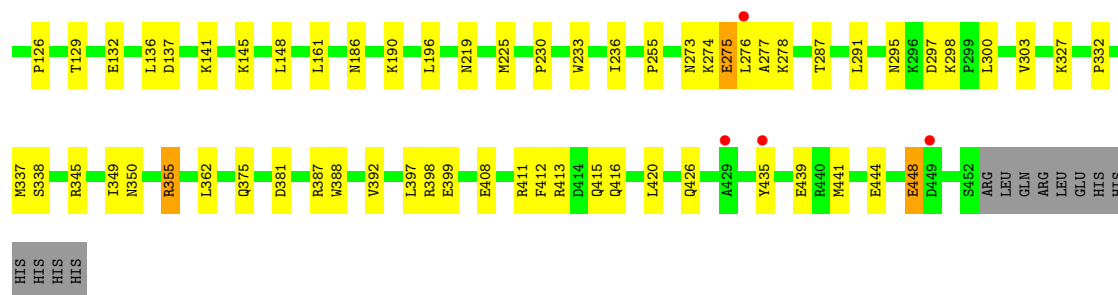


- Molecule 1: Maltose-binding periplasmic protein,Tumor necrosis factor receptor superfamily, member 25

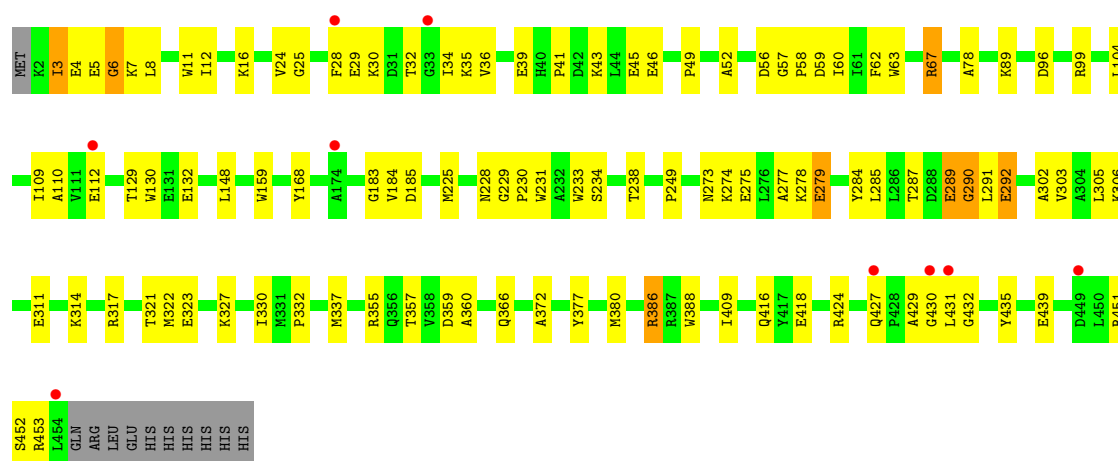
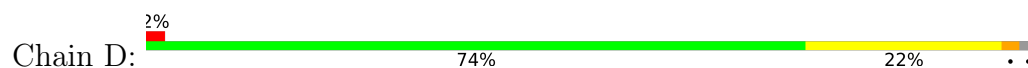


- Molecule 1: Maltose-binding periplasmic protein,Tumor necrosis factor receptor superfamily, member 25





- Molecule 1: Maltose-binding periplasmic protein,Tumor necrosis factor receptor superfamily, member 25



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	90.32Å 145.15Å 178.52Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.78 – 2.74 47.78 – 2.74	Depositor EDS
% Data completeness (in resolution range)	99.7 (47.78-2.74) 99.7 (47.78-2.74)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.18 (at 2.73Å)	Xtriage
Refinement program	PHENIX 1.12-2829	Depositor
R, R_{free}	0.210 , 0.248 0.216 , 0.254	Depositor DCC
R_{free} test set	3112 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	76.3	Xtriage
Anisotropy	0.471	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 46.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	14098	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 21.93 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 6.3250e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SO4

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	0/3550	0.62	1/4819 (0.0%)
1	B	0.40	1/3567 (0.0%)	0.79	13/4841 (0.3%)
1	C	0.39	0/3567	0.80	9/4841 (0.2%)
1	D	0.46	1/3595 (0.0%)	0.90	24/4877 (0.5%)
All	All	0.40	2/14279 (0.0%)	0.78	47/19378 (0.2%)

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	B	0	1
1	D	0	1
All	All	0	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	386	ARG	CD-NE	-6.17	1.37	1.46
1	B	180	LYS	CD-CE	5.14	1.67	1.52

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	355	ARG	CB-CG-CD	-16.41	73.56	111.30
1	C	274	LYS	CA-C-N	-13.13	101.64	122.65
1	C	274	LYS	C-N-CA	-13.13	101.64	122.65
1	D	386	ARG	CD-NE-CZ	10.98	139.77	124.40
1	B	386	ARG	NE-CZ-NH2	-10.14	110.07	119.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	5	GLU	N-CA-CB	9.51	124.75	110.29
1	B	386	ARG	CD-NE-CZ	9.40	137.56	124.40
1	B	386	ARG	CG-CD-NE	9.27	132.40	112.00
1	B	43	LYS	CG-CD-CE	-8.78	91.10	111.30
1	D	278	LYS	CA-C-N	8.11	131.81	120.29
1	D	278	LYS	C-N-CA	8.11	131.81	120.29
1	D	4	GLU	CA-CB-CG	7.87	129.83	114.10
1	D	6	GLY	N-CA-C	-7.75	101.58	114.76
1	C	30	LYS	CD-CE-NZ	-7.52	87.83	111.90
1	D	386	ARG	NE-CZ-NH2	-7.28	112.64	119.20
1	D	43	LYS	CB-CG-CD	-7.28	94.55	111.30
1	D	43	LYS	CA-CB-CG	7.14	128.38	114.10
1	B	274	LYS	CA-C-N	-6.94	109.12	120.72
1	B	274	LYS	C-N-CA	-6.94	109.12	120.72
1	D	5	GLU	CB-CA-C	-6.83	96.90	109.46
1	C	30	LYS	CG-CD-CE	6.80	126.93	111.30
1	B	180	LYS	CG-CD-CE	6.76	126.84	111.30
1	D	43	LYS	CG-CD-CE	6.73	126.77	111.30
1	C	275	GLU	N-CA-CB	6.67	120.30	110.56
1	D	279	GLU	N-CA-CB	-6.48	100.56	110.16
1	D	289	GLU	CA-CB-CG	6.32	126.75	114.10
1	D	3	ILE	CA-C-N	6.20	129.63	120.95
1	D	3	ILE	C-N-CA	6.20	129.63	120.95
1	B	43	LYS	CB-CG-CD	6.18	125.52	111.30
1	B	275	GLU	N-CA-CB	6.12	120.55	110.39
1	C	103	LYS	CB-CG-CD	6.10	125.32	111.30
1	D	289	GLU	CB-CA-C	-6.06	101.37	110.88
1	B	386	ARG	CB-CG-CD	6.01	125.12	111.30
1	D	67	ARG	N-CA-CB	-5.97	100.48	110.39
1	B	386	ARG	NE-CZ-NH1	5.91	127.41	121.50
1	A	329	GLU	CA-CB-CG	5.84	125.78	114.10
1	B	180	LYS	CB-CG-CD	-5.66	98.29	111.30
1	D	290	GLY	N-CA-C	-5.64	105.81	113.37
1	C	35	LYS	CB-CG-CD	-5.61	98.39	111.30
1	B	398	ARG	CG-CD-NE	5.58	124.28	112.00
1	C	355	ARG	CB-CG-CD	-5.54	98.57	111.30
1	D	427	GLN	CA-CB-CG	5.50	125.10	114.10
1	D	67	ARG	CB-CG-CD	5.46	123.86	111.30
1	D	292	GLU	CB-CG-CD	5.43	121.84	112.60
1	D	30	LYS	CB-CA-C	5.28	120.45	109.95
1	C	43	LYS	CB-CA-C	5.23	119.44	111.80
1	D	30	LYS	N-CA-CB	-5.17	101.72	110.41

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	386	ARG	Sidechain
1	D	386	ARG	Sidechain

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3470	0	3432	44	0
1	B	3487	0	3449	38	0
1	C	3487	0	3450	68	0
1	D	3515	0	3486	85	0
2	A	25	0	0	1	0
2	B	35	0	0	0	0
2	C	35	0	0	1	0
2	D	40	0	0	3	0
3	A	1	0	0	0	0
3	B	2	0	0	0	0
3	C	1	0	0	0	0
All	All	14098	0	13817	234	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 8.

All (234) 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:B:10:ILE:HD11	1:B:38:VAL:HG22	1.41	1.00
1:D:7:LYS:HA	1:D:34:ILE:HD11	1.45	0.96
1:D:275:GLU:OE2	1:D:279:GLU:HG3	1.67	0.93
1:D:228:ASN:HD22	1:D:229:GLY:H	1.06	0.93
1:D:24:VAL:HG11	1:D:290:GLY:HA2	1.56	0.88
1:D:289:GLU:O	1:D:292:GLU:HG3	1.75	0.87
1:C:4:GLU:HB3	1:C:7:LYS:HE2	1.57	0.86
1:A:340:PHE:HD2	1:A:341:TRP:HD1	1.28	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:32:THR:HB	1:B:34:ILE:HD13	1.61	0.81
1:B:10:ILE:HG22	1:B:60:ILE:HB	1.64	0.80
1:D:29:GLU:HA	1:D:32:THR:HG22	1.60	0.80
1:D:311:GLU:OE2	1:D:314:LYS:NZ	2.15	0.79
1:C:439:GLU:OE2	1:C:439:GLU:N	2.17	0.78
1:D:32:THR:HG23	1:D:34:ILE:H	1.47	0.78
1:C:349:ILE:HG22	1:C:355:ARG:HH12	1.47	0.77
1:C:17:GLY:HA2	1:C:298:LYS:HD2	1.65	0.77
1:A:332:PRO:HB2	1:A:337:MET:HE3	1.66	0.77
1:D:25:GLY:HA3	1:D:36:VAL:HG21	1.67	0.76
1:D:228:ASN:HD22	1:D:229:GLY:N	1.83	0.76
1:A:368:ASN:HB3	1:C:408:GLU:OE2	1.87	0.74
1:A:25:GLY:HA3	1:A:36:VAL:HG21	1.71	0.73
1:D:285:LEU:O	1:D:291:LEU:HD13	1.88	0.73
1:C:43:LYS:HD2	1:C:435:TYR:OH	1.90	0.72
1:C:123:LEU:HD21	1:C:126:PRO:HA	1.72	0.72
1:B:332:PRO:HB2	1:B:337:MET:HE3	1.72	0.72
1:D:7:LYS:HA	1:D:34:ILE:CD1	2.20	0.71
1:B:93:PHE:HD2	1:B:330:ILE:HD11	1.55	0.71
1:C:4:GLU:HB3	1:C:7:LYS:CE	2.20	0.71
1:D:62:PHE:HE2	1:D:109:ILE:HD12	1.56	0.71
1:D:357:THR:OG1	1:D:359:ASP:OD1	2.04	0.70
1:A:444:GLU:OE1	1:A:444:GLU:N	2.23	0.69
1:D:8:LEU:HB2	1:D:36:VAL:HG12	1.75	0.69
1:D:60:ILE:HD11	1:D:277:ALA:HB1	1.74	0.69
1:D:7:LYS:HG2	1:D:8:LEU:N	2.08	0.69
1:B:190:LYS:O	1:B:194:THR:HG23	1.93	0.68
1:D:63:TRP:NE1	1:D:67:ARG:HG3	2.09	0.67
1:B:153:GLN:O	1:B:345:ARG:NH1	2.28	0.67
1:B:45:GLU:HG2	1:B:46:GLU:HG2	1.77	0.67
1:C:350:ASN:OD1	1:C:355:ARG:NH1	2.28	0.66
1:A:260:VAL:HG13	1:A:330:ILE:HD13	1.76	0.66
1:D:228:ASN:ND2	1:D:229:GLY:H	1.88	0.66
1:C:412:PHE:N	2:C:505:SO4:O1	2.28	0.66
1:A:52:ALA:HB3	1:A:76:LEU:HD23	1.77	0.65
1:D:63:TRP:HE1	1:D:67:ARG:HG3	1.61	0.65
1:B:435:TYR:CE1	1:B:451:ARG:HG2	2.32	0.65
1:A:274:LYS:O	1:A:278:LYS:HG3	1.98	0.64
1:C:10:ILE:HB	1:C:38:VAL:HG13	1.79	0.64
1:D:29:GLU:HA	1:D:32:THR:CG2	2.29	0.63
1:A:340:PHE:CD2	1:A:341:TRP:HD1	2.14	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:336:GLN:N	1:A:336:GLN:OE1	2.31	0.61
1:B:129:THR:OG1	1:B:132:GLU:HG3	2.01	0.61
1:C:34:ILE:HG13	1:C:276:LEU:HD21	1.81	0.61
1:C:444:GLU:OE1	1:C:444:GLU:N	2.26	0.60
1:B:60:ILE:HG12	1:B:267:ILE:HG22	1.82	0.60
1:D:34:ILE:HD12	1:D:35:LYS:N	2.17	0.60
1:D:29:GLU:OE1	1:D:35:LYS:HD2	2.01	0.60
1:B:233:TRP:HB3	1:B:317:ARG:HH21	1.65	0.59
1:D:129:THR:OG1	1:D:132:GLU:HG3	2.02	0.59
1:A:166:GLY:O	1:A:186:ASN:ND2	2.36	0.59
1:C:332:PRO:HB2	1:C:337:MET:HE3	1.85	0.59
1:D:435:TYR:O	1:D:439:GLU:HG3	2.01	0.59
1:C:338:SER:HB2	1:C:375:GLN:HE22	1.67	0.59
1:D:148:LEU:HB2	1:D:225:MET:HE3	1.85	0.59
1:A:340:PHE:HD2	1:A:341:TRP:CD1	2.15	0.59
1:D:25:GLY:CA	1:D:36:VAL:HG21	2.33	0.58
1:D:28:PHE:O	1:D:32:THR:HG22	2.04	0.58
1:B:10:ILE:CD1	1:B:38:VAL:HG22	2.26	0.58
1:B:13:ASN:ND2	1:B:15:ASP:OD1	2.37	0.57
1:A:269:ALA:O	1:A:274:LYS:NZ	2.33	0.57
1:C:129:THR:OG1	1:C:132:GLU:HG3	2.05	0.57
1:B:336:GLN:OE1	1:B:336:GLN:N	2.29	0.57
1:C:295:ASN:ND2	1:C:300:LEU:H	2.02	0.57
1:A:25:GLY:CA	1:A:36:VAL:HG21	2.35	0.57
1:C:41:PRO:HD2	1:C:44:LEU:HD12	1.85	0.57
1:D:12:ILE:HD11	1:D:16:LYS:HG3	1.85	0.57
1:D:112:GLU:OE2	1:D:231:TRP:HE3	1.87	0.57
1:B:153:GLN:HA	1:B:349:ILE:HD11	1.87	0.57
1:D:228:ASN:HD21	1:D:231:TRP:NE1	2.04	0.56
1:C:137:ASP:OD2	1:C:141:LYS:HE2	2.05	0.56
1:B:34:ILE:HD12	1:B:34:ILE:N	2.21	0.55
1:C:388:TRP:HB3	1:C:416:GLN:CD	2.31	0.55
1:D:62:PHE:CE2	1:D:109:ILE:HD12	2.38	0.55
1:D:63:TRP:CD1	1:D:67:ARG:HG3	2.41	0.55
1:C:388:TRP:CE3	1:C:416:GLN:HG3	2.40	0.55
1:C:381:ASP:OD1	1:C:413:ARG:NH2	2.40	0.55
1:C:13:ASN:ND2	1:C:15:ASP:OD1	2.39	0.55
1:A:260:VAL:CG1	1:A:330:ILE:HD13	2.37	0.55
1:D:34:ILE:HD12	1:D:35:LYS:H	1.72	0.55
1:C:110:ALA:HA	1:C:303:VAL:HA	1.88	0.55
1:D:332:PRO:HB2	1:D:337:MET:HE3	1.87	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:397:LEU:HD12	1:C:426:GLN:HE22	1.73	0.54
1:D:45:GLU:HG2	1:D:46:GLU:HG3	1.89	0.54
1:D:112:GLU:OE2	1:D:231:TRP:CE3	2.60	0.54
1:B:7:LYS:O	1:B:273:ASN:ND2	2.41	0.54
1:D:377:TYR:HA	1:D:380:MET:HG3	1.89	0.53
1:C:148:LEU:HB2	1:C:225:MET:HE3	1.89	0.53
1:D:56:ASP:CG	1:D:57:GLY:H	2.15	0.53
1:D:159:TRP:HH2	1:D:184:VAL:HG12	1.73	0.53
1:C:122:LEU:HD21	1:C:145:LYS:HE3	1.89	0.53
1:C:435:TYR:O	1:C:439:GLU:OE2	2.27	0.53
1:D:388:TRP:CE2	1:D:416:GLN:HG2	2.44	0.53
1:C:52:ALA:CB	1:C:76:LEU:HD22	2.39	0.53
1:C:255:PRO:HB3	1:C:327:LYS:HD3	1.91	0.53
1:B:10:ILE:HD12	1:B:21:LEU:HD21	1.91	0.52
1:C:350:ASN:HA	1:C:355:ARG:NH1	2.24	0.52
1:D:7:LYS:O	1:D:273:ASN:ND2	2.40	0.52
1:D:424:ARG:CZ	1:D:429:ALA:HB2	2.40	0.52
1:A:26:LYS:O	1:A:26:LYS:HG3	2.09	0.52
1:C:98:VAL:CG2	1:C:106:ALA:HB3	2.40	0.52
1:C:21:LEU:HD23	1:C:38:VAL:HG11	1.91	0.52
1:A:7:LYS:O	1:A:273:ASN:ND2	2.42	0.51
1:B:377:TYR:HA	1:B:380:MET:HG3	1.92	0.51
1:C:275:GLU:HB2	1:C:278:LYS:HG2	1.91	0.51
1:D:230:PRO:HA	1:D:233:TRP:CE2	2.45	0.51
1:B:42:ASP:OD2	1:B:43:LYS:HG2	2.10	0.51
1:A:153:GLN:HA	1:A:349:ILE:HD11	1.91	0.51
1:D:89:LYS:O	1:D:306:LYS:HG3	2.11	0.51
1:B:439:GLU:OE2	1:B:451:ARG:NH2	2.43	0.51
1:D:359:ASP:OD1	1:D:359:ASP:N	2.43	0.51
1:C:11:TRP:CD2	1:C:58:PRO:HG3	2.46	0.51
1:A:232:ALA:O	1:A:236:ILE:HG13	2.11	0.51
1:B:232:ALA:O	1:B:236:ILE:HG13	2.12	0.50
1:D:359:ASP:OD1	1:D:360:ALA:N	2.43	0.50
1:D:3:ILE:HG13	1:D:56:ASP:HA	1.92	0.50
1:B:274:LYS:O	1:B:278:LYS:HG3	2.12	0.50
1:D:424:ARG:NH2	2:D:505:SO4:O4	2.44	0.50
1:D:49:PRO:HB2	1:D:431:LEU:HB3	1.93	0.50
1:D:45:GLU:HG2	1:D:46:GLU:N	2.25	0.50
1:A:185:ASP:O	1:A:190:LYS:HE2	2.12	0.50
1:A:255:PRO:HB3	1:A:327:LYS:HD3	1.94	0.50
1:D:3:ILE:CG1	1:D:56:ASP:HA	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:78:ALA:HB2	1:D:274:LYS:HE2	1.94	0.49
1:A:411:ARG:O	1:A:415:GLN:HG3	2.12	0.49
1:B:194:THR:HG22	1:B:358:VAL:HG11	1.94	0.49
1:A:271:SER:O	1:A:274:LYS:HE3	2.12	0.49
1:B:110:ALA:HA	1:B:303:VAL:HA	1.95	0.49
1:A:388:TRP:HB3	1:A:416:GLN:CD	2.37	0.48
1:D:60:ILE:HD11	1:D:277:ALA:CB	2.41	0.48
1:D:317:ARG:O	1:D:321:THR:HG23	2.13	0.48
1:D:409:ILE:HD13	1:D:418:GLU:HG3	1.94	0.48
1:C:388:TRP:CD2	1:C:416:GLN:HG3	2.48	0.48
1:C:398:ARG:HG3	1:C:399:GLU:H	1.79	0.48
1:D:372:ALA:N	2:D:503:SO4:O2	2.43	0.48
1:C:7:LYS:O	1:C:273:ASN:ND2	2.47	0.48
1:D:6:GLY:O	1:D:34:ILE:HD13	2.12	0.48
1:D:234:SER:O	1:D:238:THR:HG23	2.13	0.48
1:C:44:LEU:HD21	1:C:61:ILE:HD11	1.95	0.48
1:C:398:ARG:HG3	1:C:399:GLU:N	2.29	0.48
1:D:323:GLU:O	1:D:327:LYS:HG3	2.13	0.48
1:C:123:LEU:HD11	1:C:136:LEU:HD21	1.96	0.47
1:B:11:TRP:CD2	1:B:58:PRO:HG3	2.49	0.47
1:D:35:LYS:HG3	1:D:36:VAL:N	2.30	0.47
1:D:287:THR:O	1:D:291:LEU:HB2	2.15	0.47
1:C:7:LYS:H	1:C:273:ASN:ND2	2.12	0.47
1:D:96:ASP:HB2	1:D:330:ILE:HD13	1.96	0.47
1:C:10:ILE:HG13	1:C:38:VAL:HG22	1.96	0.47
1:D:185:ASP:HB2	1:D:366:GLN:HB2	1.97	0.47
1:C:42:ASP:OD2	1:C:43:LYS:N	2.47	0.47
1:A:129:THR:HB	1:A:132:GLU:HG3	1.97	0.46
1:B:186:ASN:O	1:B:190:LYS:HG3	2.14	0.46
1:D:99:ARG:CZ	1:D:104:LEU:HD21	2.44	0.46
1:D:24:VAL:HG11	1:D:290:GLY:CA	2.36	0.46
1:B:350:ASN:OD1	1:B:355:ARG:NH1	2.48	0.46
1:D:451:ARG:C	1:D:453:ARG:H	2.23	0.46
1:A:260:VAL:HG13	1:A:330:ILE:HA	1.97	0.46
1:A:435:TYR:CZ	1:A:439:GLU:OE2	2.68	0.46
1:B:435:TYR:HB3	1:B:451:ARG:HH12	1.80	0.46
1:C:39:GLU:C	1:C:41:PRO:HD3	2.40	0.46
1:D:453:ARG:HA	1:D:453:ARG:HD3	1.83	0.46
1:D:429:ALA:HA	1:D:430:GLY:HA2	1.77	0.46
1:C:411:ARG:O	1:C:415:GLN:HG3	2.16	0.45
1:B:69:GLY:HA3	1:B:333:ASN:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:148:LEU:HB2	1:B:225:MET:HE3	1.99	0.45
1:A:179:ILE:CD1	1:A:336:GLN:HE21	2.30	0.45
1:D:302:ALA:HB1	1:D:322:MET:HE2	1.99	0.45
1:C:338:SER:CB	1:C:375:GLN:HE22	2.27	0.45
1:C:420:LEU:HD23	1:C:420:LEU:HA	1.83	0.45
1:A:186:ASN:O	1:A:190:LYS:HG3	2.16	0.45
1:B:298:LYS:HB3	1:B:298:LYS:HE2	1.52	0.45
1:D:59:ASP:C	1:D:60:ILE:HD13	2.42	0.44
1:A:11:TRP:CD2	1:A:58:PRO:HG3	2.52	0.44
1:A:110:ALA:HA	1:A:303:VAL:HA	1.99	0.44
1:B:387:ARG:HH11	1:B:443:LEU:HD21	1.82	0.44
1:D:39:GLU:C	1:D:41:PRO:HD3	2.42	0.44
1:C:287:THR:O	1:C:291:LEU:HB2	2.17	0.44
1:B:42:ASP:O	1:B:43:LYS:HB2	2.17	0.44
1:C:98:VAL:HG21	1:C:106:ALA:HB3	1.98	0.44
1:D:159:TRP:CH2	1:D:184:VAL:HG12	2.53	0.44
1:A:99:ARG:NH2	1:A:104:LEU:HD21	2.33	0.44
1:A:190:LYS:HD3	1:A:359:ASP:OD1	2.19	0.43
1:A:129:THR:HG23	1:A:131:GLU:OE2	2.19	0.43
1:D:130:TRP:CD1	1:D:249:PRO:HB2	2.54	0.43
1:D:432:GLY:N	2:D:505:SO4:O3	2.49	0.43
1:A:159:TRP:HH2	1:A:184:VAL:HG12	1.84	0.43
1:A:66:ASP:OD1	1:A:67:ARG:N	2.51	0.43
1:A:98:VAL:O	1:A:105:ILE:HG13	2.18	0.43
1:C:230:PRO:HA	1:C:233:TRP:CE2	2.53	0.43
1:C:345:ARG:O	1:C:349:ILE:HG12	2.18	0.43
1:A:388:TRP:HB3	1:A:416:GLN:OE1	2.19	0.43
1:B:429:ALA:HA	1:B:430:GLY:HA2	1.75	0.43
1:A:56:ASP:CG	1:A:57:GLY:H	2.27	0.43
1:D:110:ALA:HA	1:D:303:VAL:HA	2.00	0.43
1:D:225:MET:HE3	1:D:225:MET:HB2	1.89	0.43
1:C:19:ASN:HB2	1:C:297:ASP:OD2	2.18	0.43
1:C:42:ASP:O	1:C:47:LYS:HE3	2.19	0.43
1:A:411:ARG:HB3	2:A:504:SO4:O1	2.19	0.42
1:C:444:GLU:O	1:C:448:GLU:HB2	2.18	0.42
1:D:52:ALA:HA	1:D:56:ASP:O	2.19	0.42
1:D:89:LYS:O	1:D:305:LEU:HD12	2.18	0.42
1:C:186:ASN:O	1:C:190:LYS:HG3	2.20	0.42
1:C:362:LEU:HD23	1:C:362:LEU:HA	1.80	0.42
1:C:392:VAL:HG22	1:C:397:LEU:HB3	2.02	0.42
1:C:161:LEU:HD23	1:C:196:LEU:HB2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:34:ILE:N	1:B:34:ILE:CD1	2.83	0.42
1:A:39:GLU:C	1:A:41:PRO:HD3	2.45	0.41
1:C:8:LEU:HD11	1:C:277:ALA:HA	2.02	0.41
1:A:129:THR:HG22	1:A:132:GLU:H	1.85	0.41
1:D:451:ARG:C	1:D:453:ARG:N	2.79	0.41
1:C:44:LEU:HD21	1:C:61:ILE:CD1	2.51	0.41
1:C:68:PHE:HB3	1:C:105:ILE:HD12	2.01	0.41
1:A:225:MET:HE3	1:A:225:MET:HB2	1.91	0.41
1:C:387:ARG:NH2	1:C:441:MET:HB3	2.36	0.41
1:B:218:PHE:HA	1:B:223:THR:HG22	2.03	0.41
1:D:28:PHE:HA	1:D:284:TYR:CE2	2.56	0.41
1:D:168:TYR:CE1	1:D:183:GLY:HA3	2.55	0.41
1:C:219:ASN:OD1	1:C:236:ILE:HG12	2.20	0.41
1:C:255:PRO:CB	1:C:327:LYS:HD3	2.51	0.40
1:C:60:ILE:HD11	1:C:277:ALA:HB1	2.03	0.40
1:D:11:TRP:CD2	1:D:58:PRO:HG3	2.56	0.40
1:D:25:GLY:O	1:D:36:VAL:HG22	2.21	0.40
1:D:56:ASP:OD1	1:D:57:GLY:N	2.50	0.40
1:A:69:GLY:HA3	1:A:333:ASN:O	2.21	0.40
1:C:52:ALA:HB1	1:C:76:LEU:HD22	2.04	0.40
1:C:225:MET:HE3	1:C:225:MET:HB2	1.98	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	446/464 (96%)	431 (97%)	14 (3%)	1 (0%)	43	62
1	B	448/464 (97%)	439 (98%)	9 (2%)	0	100	100
1	C	448/464 (97%)	438 (98%)	9 (2%)	1 (0%)	43	62
1	D	451/464 (97%)	440 (98%)	10 (2%)	1 (0%)	43	62

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1793/1856 (97%)	1748 (98%)	42 (2%)	3 (0%)	43 62

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	448	GLU
1	D	452	SER
1	A	175	GLY

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	347/363 (96%)	347 (100%)	0	100 100
1	B	349/363 (96%)	349 (100%)	0	100 100
1	C	349/363 (96%)	349 (100%)	0	100 100
1	D	352/363 (97%)	352 (100%)	0	100 100
All	All	1397/1452 (96%)	1397 (100%)	0	100 100

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

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	40	HIS
1	A	375	GLN
1	A	425	GLN
1	A	426	GLN
1	A	427	GLN
1	B	242	ASN
1	C	50	GLN
1	C	254	GLN
1	C	283	ASN
1	C	295	ASN

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Mol	Chain	Res	Type
1	C	375	GLN
1	C	416	GLN
1	D	228	ASN
1	D	326	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

27 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	$\# Z > 2$
2	SO4	C	501	-	4,4,4	0.26	0	6,6,6	0.06	0
2	SO4	D	503	-	4,4,4	0.23	0	6,6,6	0.05	0
2	SO4	C	507	-	4,4,4	0.24	0	6,6,6	0.14	0
2	SO4	A	504	-	4,4,4	0.25	0	6,6,6	0.19	0
2	SO4	B	503	-	4,4,4	0.28	0	6,6,6	0.19	0
2	SO4	D	502	-	4,4,4	0.36	0	6,6,6	0.65	0
2	SO4	B	506	-	4,4,4	0.34	0	6,6,6	0.41	0
2	SO4	C	503	-	4,4,4	0.23	0	6,6,6	0.11	0
2	SO4	D	501	-	4,4,4	0.24	0	6,6,6	0.13	0
2	SO4	D	504	-	4,4,4	0.27	0	6,6,6	0.09	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	B	502	-	4,4,4	0.23	0	6,6,6	0.06	0
2	SO4	B	505	-	4,4,4	0.24	0	6,6,6	0.11	0
2	SO4	A	502	-	4,4,4	0.21	0	6,6,6	0.13	0
2	SO4	D	506	-	4,4,4	0.26	0	6,6,6	0.14	0
2	SO4	C	506	-	4,4,4	0.24	0	6,6,6	0.12	0
2	SO4	C	504	-	4,4,4	0.21	0	6,6,6	0.35	0
2	SO4	B	507	-	4,4,4	0.23	0	6,6,6	0.08	0
2	SO4	C	505	-	4,4,4	0.25	0	6,6,6	0.25	0
2	SO4	B	501	-	4,4,4	0.26	0	6,6,6	0.11	0
2	SO4	A	501	-	4,4,4	0.24	0	6,6,6	0.08	0
2	SO4	C	502	-	4,4,4	0.23	0	6,6,6	0.10	0
2	SO4	A	503	-	4,4,4	0.23	0	6,6,6	0.41	0
2	SO4	A	505	-	4,4,4	0.27	0	6,6,6	0.18	0
2	SO4	D	508	-	4,4,4	0.20	0	6,6,6	0.22	0
2	SO4	B	504	-	4,4,4	0.23	0	6,6,6	0.11	0
2	SO4	D	505	-	4,4,4	0.48	0	6,6,6	0.29	0
2	SO4	D	507	-	4,4,4	0.24	0	6,6,6	0.09	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	503	SO4	1	0
2	A	504	SO4	1	0
2	C	505	SO4	1	0
2	D	505	SO4	2	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

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	448/464 (96%)	0.05	7 (1%) 70 68	50, 75, 110, 155	0
1	B	450/464 (96%)	0.03	3 (0%) 84 84	49, 73, 108, 153	0
1	C	450/464 (96%)	0.11	6 (1%) 75 74	49, 76, 123, 159	0
1	D	453/464 (97%)	0.21	9 (1%) 65 62	44, 83, 129, 158	0
All	All	1801/1856 (97%)	0.10	25 (1%) 73 72	44, 76, 118, 159	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	329	GLU	3.7
1	A	175	GLY	3.4
1	D	454	LEU	3.4
1	B	435	TYR	3.2
1	D	28	PHE	2.9
1	D	33	GLY	2.8
1	A	429	ALA	2.8
1	D	431	LEU	2.7
1	A	341	TRP	2.6
1	C	449	ASP	2.6
1	D	112	GLU	2.5
1	D	174	ALA	2.4
1	C	435	TYR	2.3
1	D	430	GLY	2.3
1	A	174	ALA	2.2
1	C	429	ALA	2.2
1	A	173	ALA	2.2
1	A	180	LYS	2.2
1	C	276	LEU	2.2
1	C	53	ALA	2.1
1	D	449	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	174	ALA	2.1
1	B	429	ALA	2.1
1	D	427	GLN	2.0
1	C	36	VAL	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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($\text{\AA}^2$)	Q<0.9
2	SO4	A	505	5/5	0.58	0.12	166,168,168,168	0
2	SO4	C	507	5/5	0.64	0.13	150,150,150,152	0
2	SO4	C	503	5/5	0.68	0.08	162,162,163,163	0
2	SO4	C	506	5/5	0.68	0.12	136,137,138,138	0
2	SO4	B	507	5/5	0.68	0.11	146,146,147,147	0
2	SO4	C	502	5/5	0.69	0.10	128,128,128,129	0
2	SO4	B	502	5/5	0.69	0.15	163,163,164,165	0
2	SO4	D	508	5/5	0.70	0.14	157,157,158,158	0
2	SO4	D	504	5/5	0.71	0.12	131,131,131,133	0
2	SO4	D	507	5/5	0.73	0.08	151,151,151,152	0
2	SO4	D	506	5/5	0.74	0.07	142,144,145,145	0
2	SO4	B	504	5/5	0.75	0.09	133,134,135,135	0
2	SO4	A	502	5/5	0.77	0.09	132,132,133,133	0
2	SO4	B	501	5/5	0.78	0.08	117,119,120,120	0
2	SO4	D	501	5/5	0.78	0.08	132,133,134,134	0
2	SO4	C	501	5/5	0.80	0.11	127,129,130,130	0
2	SO4	B	505	5/5	0.82	0.12	113,113,114,114	0
2	SO4	D	505	5/5	0.83	0.12	116,117,117,119	0
2	SO4	D	503	5/5	0.85	0.12	103,104,105,106	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	A	501	5/5	0.86	0.08	124,125,125,125	0
2	SO4	B	506	5/5	0.89	0.10	88,89,91,93	0
2	SO4	C	504	5/5	0.93	0.07	69,69,71,72	0
2	SO4	A	503	5/5	0.94	0.07	69,71,72,78	0
2	SO4	B	503	5/5	0.96	0.07	67,68,71,73	0
2	SO4	C	505	5/5	0.97	0.08	69,71,73,74	0
2	SO4	A	504	5/5	0.97	0.07	69,70,71,78	0
2	SO4	D	502	5/5	0.98	0.10	67,69,79,79	0

6.5 Other polymers [i](#)

There are no such residues in this entry.