



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 09:37 AM UTC

PDB ID : 7ZMN / pdb_00007zmn
Title : Crystal structure of human RECQL5 helicase APO form in complex with engineered nanobody (Gluebody) G3-048
Authors : Ye, M.; Makola, M.; Newman, J.A.; Fairhead, M.; MacLean, E.; Krojer, T.; Aitkenhead, H.; Bountra, C.; Gileadi, O.; von Delft, F.
Deposited on : 2022-04-19
Resolution : 3.20 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

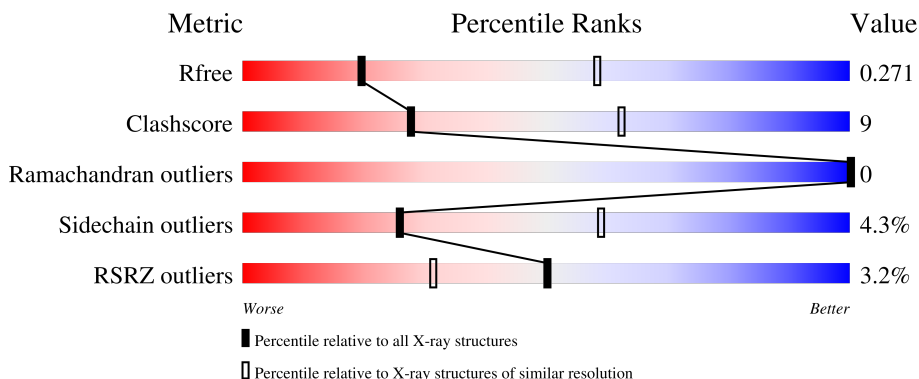
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	445	4% 76% 22% ..
1	B	445	4% 76% 22% ..
2	C	127	74% 24% .
2	K	127	69% 28% ..

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8800 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 DNA helicase Q5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	440	Total	C	N	O	S	0	4	0
			3405	2153	605	624	23			
1	B	440	Total	C	N	O	S	0	4	0
			3447	2177	620	627	23			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	9	SER	-	expression tag	UNP O94762
A	10	MET	-	expression tag	UNP O94762
B	9	SER	-	expression tag	UNP O94762
B	10	MET	-	expression tag	UNP O94762

- Molecule 2 is a protein called Gluebody G3-048.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	K	124	Total	C	N	O	S	0	2	0
			954	592	164	193	5			
2	C	124	Total	C	N	O	S	0	2	0
			950	590	163	192	5			

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		
3	B	1	Total	Zn	0	0
			1	1		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

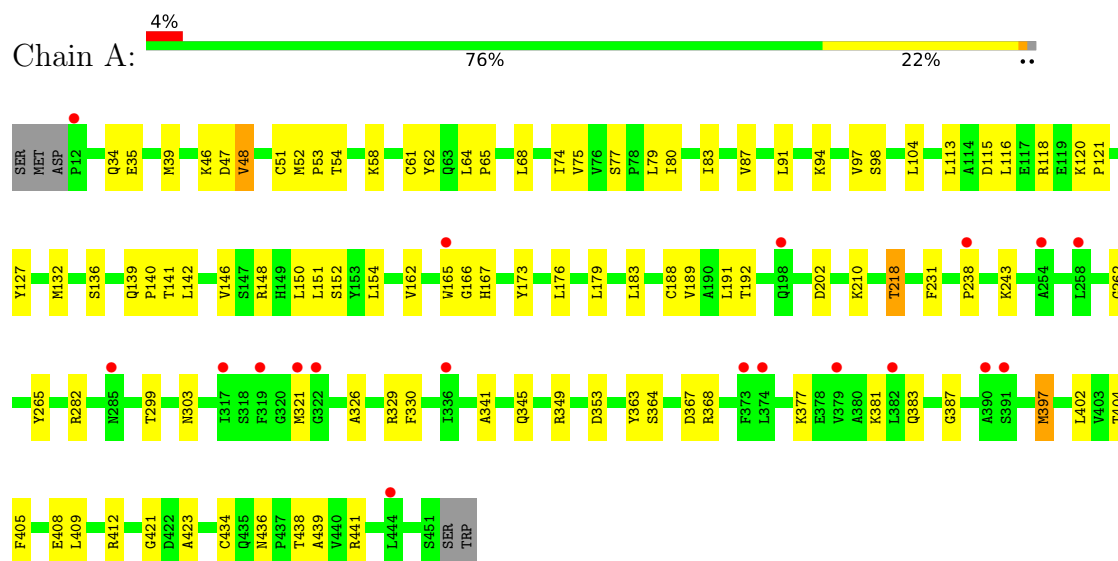
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	7	Total	O	0	0
			7	7		
5	B	9	Total	O	0	0
			9	9		
5	K	6	Total	O	0	0
			6	6		
5	C	5	Total	O	0	0
			5	5		

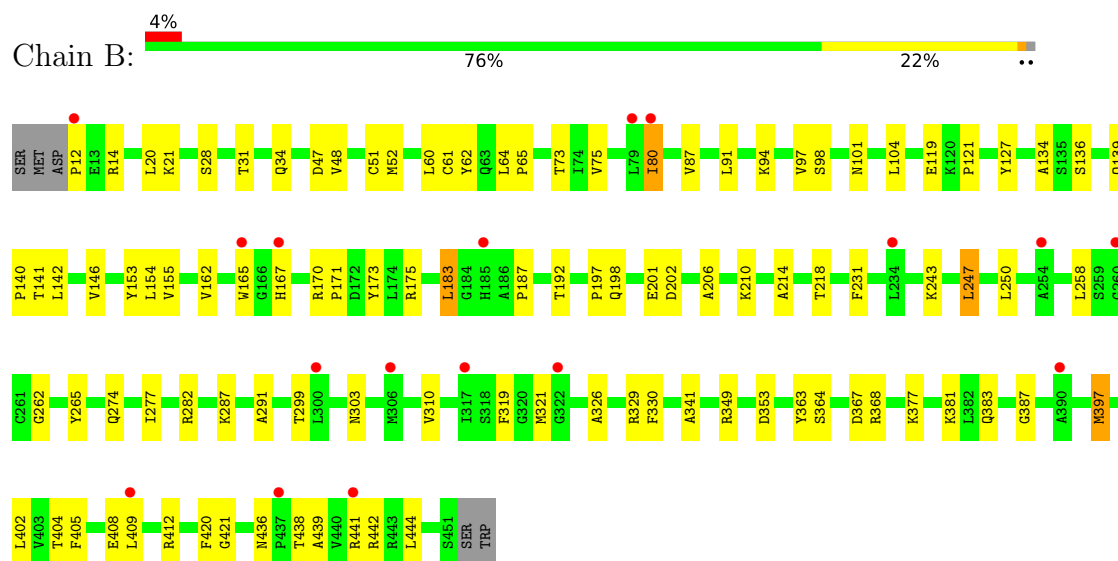
3 Residue-property plots [i](#)

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 DNA helicase Q5



• Molecule 1: ATP-dependent DNA helicase Q5



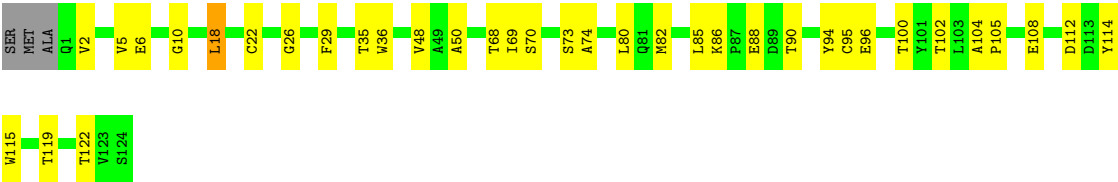
• Molecule 2: Gluebody G3-048

Chain K:

69%

28%

..



● Molecule 2: Gluebody G3-048

Chain C:

74%

24%

.



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	79.50Å 89.57Å 259.06Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	129.53 – 3.20 129.53 – 3.20	Depositor EDS
% Data completeness (in resolution range)	99.5 (129.53-3.20) 99.5 (129.53-3.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.78 (at 3.19Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.242 , 0.275 0.241 , 0.271	Depositor DCC
R_{free} test set	1559 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	61.8	Xtriage
Anisotropy	0.119	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 65.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	8800	wwPDB-VP
Average B, all atoms (Å ²)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.14% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, 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	1.06	0/3496	1.57	7/4728 (0.1%)
1	B	1.04	0/3520	1.57	2/4756 (0.0%)
2	C	1.04	0/979	1.46	3/1328 (0.2%)
2	K	1.04	1/979 (0.1%)	1.42	4/1328 (0.3%)
All	All	1.05	1/8974 (0.0%)	1.54	16/12140 (0.1%)

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	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	K	70	SER	CA-CB	-5.26	1.44	1.53

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	167[A]	HIS	CB-CA-C	6.74	120.10	111.40
1	A	167[B]	HIS	CB-CA-C	6.74	120.10	111.40
2	K	26	GLY	CA-C-O	-6.56	117.34	122.52
1	A	166	GLY	CA-C-O	-5.70	116.17	121.18
1	A	238	PRO	CA-C-N	5.48	127.63	120.28
1	A	238	PRO	C-N-CA	5.48	127.63	120.28
1	B	187	PRO	CB-CA-C	-5.39	104.51	111.46
1	A	341	ALA	CA-C-N	5.36	125.89	119.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	341	ALA	C-N-CA	5.36	125.89	119.94
1	B	121	PRO	CB-CA-C	-5.23	104.13	110.98
2	C	105	PRO	N-CA-C	-5.21	107.34	113.86
2	C	73	SER	CA-C-N	5.19	127.50	120.65
2	C	73	SER	C-N-CA	5.19	127.50	120.65
2	K	73	SER	CA-C-N	5.02	127.27	120.65
2	K	73	SER	C-N-CA	5.02	127.27	120.65
2	K	10	GLY	CA-C-O	-5.01	117.40	121.76

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	77	SER	Peptide

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	3405	0	3398	60	0
1	B	3447	0	3473	63	0
2	C	950	0	905	20	0
2	K	954	0	909	18	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	10	0	0	1	0
4	B	5	0	0	0	0
5	A	7	0	0	3	0
5	B	9	0	0	3	0
5	C	5	0	0	2	0
5	K	6	0	0	1	0
All	All	8800	0	8685	158	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 9.

All (158) 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)
2:C:71[B]:ARG:HD2	5:C:205:HOH:O	1.72	0.88
1:A:136:SER:HA	1:A:139:GLN:HG2	1.57	0.86
1:A:183:LEU:O	1:A:183:LEU:HD12	1.73	0.86
1:B:12:PRO:HB2	5:B:607:HOH:O	1.85	0.76
1:B:136:SER:HA	1:B:139:GLN:HG2	1.67	0.75
1:B:243:LYS:HD3	1:B:282:ARG:HB3	1.68	0.74
1:A:243:LYS:HD3	1:A:282:ARG:HB3	1.69	0.72
1:B:198[A]:GLN:HA	1:B:198[A]:GLN:OE1	1.90	0.70
1:B:167[A]:HIS:CD2	1:B:167[A]:HIS:O	2.46	0.68
1:A:349:ARG:NH2	5:A:601:HOH:O	2.23	0.68
1:B:438:THR:HG22	1:B:441[A]:ARG:HH21	1.59	0.67
1:B:438:THR:HG22	1:B:441[A]:ARG:NH2	2.10	0.67
1:A:151:LEU:CD2	1:A:183:LEU:HD22	2.26	0.66
1:B:167[A]:HIS:O	1:B:167[A]:HIS:HD2	1.79	0.66
1:B:154:LEU:HD22	1:B:183:LEU:HD12	1.78	0.66
1:B:162:VAL:HG22	1:B:173:TYR:HB3	1.78	0.65
1:A:162:VAL:HG12	1:A:173:TYR:HB3	1.80	0.63
2:C:71[B]:ARG:HH11	2:C:71[B]:ARG:HB2	1.63	0.62
1:B:397:MET:HA	1:B:397:MET:HE2	1.81	0.62
1:A:151:LEU:HD23	1:A:183:LEU:HD22	1.80	0.61
1:A:183:LEU:O	1:A:183:LEU:CD1	2.47	0.61
2:K:69:ILE:CG2	2:K:80:LEU:HD13	2.30	0.61
2:C:69:ILE:CG2	2:C:80:LEU:HD13	2.31	0.61
2:C:35:THR:O	2:C:95:CYS:HA	2.02	0.59
2:K:35:THR:O	2:K:95:CYS:HA	2.02	0.59
2:C:71[B]:ARG:CD	5:C:205:HOH:O	2.42	0.58
1:A:321:MET:HB2	1:A:349:ARG:NE	2.19	0.58
1:B:162:VAL:HG11	1:B:206:ALA:HB3	1.86	0.58
1:A:397:MET:HA	1:A:397:MET:HE2	1.84	0.57
1:B:134:ALA:HB1	1:B:175[A]:ARG:HD2	1.85	0.57
1:B:287:LYS:HG2	1:B:310:VAL:HG11	1.85	0.57
1:B:438:THR:CG2	1:B:441[A]:ARG:HH21	2.16	0.57
1:B:134:ALA:CB	1:B:175[A]:ARG:HD2	2.35	0.57
1:B:48:VAL:HG22	1:B:214:ALA:HB3	1.87	0.57
1:B:64:LEU:HB3	1:B:65:PRO:HD3	1.88	0.56
1:B:438:THR:CG2	1:B:441[A]:ARG:NH2	2.68	0.56
2:C:6:GLU:OE2	2:C:94:TYR:HA	2.04	0.56
2:C:90:THR:HG23	2:C:122:THR:HA	1.87	0.56
1:A:326:ALA:HB1	1:A:353:ASP:HB3	1.88	0.56
2:K:90:THR:HG23	2:K:122:THR:HA	1.88	0.56
1:A:64:LEU:HB3	1:A:65:PRO:HD3	1.89	0.55
1:B:247:LEU:HD23	1:B:258:LEU:HD11	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:K:2:VAL:CG2	2:K:102:THR:HG21	2.36	0.55
2:K:69:ILE:HG23	2:K:80:LEU:HD13	1.87	0.55
1:B:20:LEU:HD13	1:B:60:LEU:HD21	1.89	0.55
2:K:6:GLU:OE2	2:K:94:TYR:HA	2.06	0.54
2:C:71[B]:ARG:HH11	2:C:71[B]:ARG:CB	2.20	0.54
1:B:14:ARG:HG2	5:B:607:HOH:O	2.07	0.54
1:B:34:GLN:HA	1:B:52:MET:HE1	1.89	0.53
1:B:326:ALA:HB1	1:B:353:ASP:HB3	1.89	0.53
2:C:71[B]:ARG:CB	2:C:71[B]:ARG:NH1	2.72	0.53
2:K:74:ALA:HB3	5:K:202:HOH:O	2.09	0.53
1:A:421:GLY:HA2	2:C:29:PHE:CZ	2.43	0.52
1:A:423:ALA:HB2	2:C:103:LEU:O	2.08	0.52
2:C:50:ALA:C	2:C:69:ILE:HD11	2.34	0.52
1:A:349:ARG:NH1	5:A:601:HOH:O	2.42	0.52
1:B:262:GLY:HA3	1:B:330:PHE:CE2	2.44	0.52
2:C:100:THR:HB	2:C:114:TYR:CE2	2.45	0.52
1:A:34:GLN:HA	1:A:52:MET:HE1	1.92	0.52
1:A:265:TYR:OH	1:A:349:ARG:NH1	2.44	0.51
2:C:69:ILE:HG23	2:C:80:LEU:HD13	1.92	0.51
1:B:404:THR:O	1:B:408:GLU:HG2	2.11	0.50
1:A:404:THR:O	1:A:408:GLU:HG2	2.12	0.50
1:B:377:LYS:O	1:B:381:LYS:HG2	2.11	0.50
1:B:139:GLN:N	1:B:140:PRO:HD2	2.27	0.49
2:K:104:ALA:HB1	2:K:105:PRO:HD2	1.94	0.49
1:A:139:GLN:N	1:A:140:PRO:HD2	2.27	0.49
2:K:50:ALA:C	2:K:69:ILE:HD11	2.36	0.49
1:A:377:LYS:O	1:A:381:LYS:HG2	2.11	0.49
1:B:73:THR:HG23	1:B:153:TYR:HB2	1.94	0.49
1:A:115:ASP:O	1:A:118:ARG:HG2	2.13	0.48
1:B:402:LEU:O	1:B:405:PHE:HB3	2.13	0.48
1:A:61:CYS:O	1:A:65:PRO:HG2	2.13	0.48
1:A:402:LEU:O	1:A:405:PHE:HB3	2.13	0.48
1:B:231:PHE:HA	1:B:363:TYR:O	2.13	0.48
1:A:363:TYR:OH	1:A:368:ARG:HD3	2.14	0.48
2:K:18:LEU:HD12	2:K:85:LEU:CD1	2.43	0.48
2:K:100:THR:HB	2:K:114:TYR:CE2	2.48	0.48
1:B:62:TYR:C	1:B:65:PRO:HD2	2.39	0.48
1:B:87:VAL:HG23	1:B:97:VAL:HG22	1.96	0.48
1:A:136:SER:HA	1:A:139:GLN:CG	2.37	0.48
1:A:62:TYR:C	1:A:65:PRO:HD2	2.39	0.48
1:B:136:SER:HA	1:B:139:GLN:CG	2.42	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:262:GLY:HA3	1:A:330:PHE:CE2	2.49	0.48
1:B:363:TYR:OH	1:B:368:ARG:HD3	2.13	0.47
2:K:112:ASP:HB3	2:K:114:TYR:CZ	2.49	0.47
1:A:47:ASP:OD1	1:A:210:LYS:HE2	2.14	0.47
1:A:64:LEU:C	1:A:64:LEU:HD23	2.40	0.47
2:K:36:TRP:O	2:K:48:VAL:HB	2.14	0.47
1:B:61:CYS:O	1:B:65:PRO:HG2	2.14	0.47
1:A:436:ASN:OD1	1:A:439:ALA:HB2	2.14	0.47
2:C:18:LEU:HD12	2:C:85:LEU:CD1	2.44	0.47
2:C:4:LEU:HB3	2:C:116:GLY:HA2	1.97	0.47
1:B:91:LEU:O	1:B:94:LYS:N	2.41	0.47
1:B:162:VAL:HG11	1:B:206:ALA:CB	2.45	0.46
1:A:91:LEU:O	1:A:94:LYS:N	2.40	0.46
1:A:83:ILE:HG23	1:A:127:TYR:HB3	1.97	0.46
1:A:51:CYS:HA	1:A:192:THR:O	2.16	0.46
1:A:299:THR:O	1:A:303:ASN:HB2	2.16	0.46
1:B:421:GLY:HA2	2:K:29:PHE:CZ	2.50	0.46
1:B:436:ASN:OD1	1:B:439:ALA:HB2	2.14	0.46
2:C:36:TRP:O	2:C:48:VAL:HB	2.16	0.46
1:A:405:PHE:CE2	1:A:412:ARG:HB3	2.51	0.46
1:B:341:ALA:HA	1:B:420:PHE:HZ	1.80	0.46
1:A:74:ILE:CD1	1:A:151:LEU:HD11	2.46	0.45
1:B:64:LEU:C	1:B:64:LEU:HD23	2.41	0.45
1:A:176:LEU:O	1:A:179:LEU:HB3	2.17	0.45
1:A:46:LYS:HA	1:A:210:LYS:HE3	1.99	0.45
1:A:116:LEU:O	1:A:148:ARG:NH2	2.50	0.44
2:K:82:MET:HE2	2:K:85:LEU:HD21	1.99	0.44
2:C:112:ASP:HB3	2:C:114:TYR:CZ	2.51	0.44
1:A:231:PHE:HA	1:A:363:TYR:O	2.18	0.44
1:B:383:GLN:O	1:B:387:GLY:HA2	2.17	0.44
1:A:183:LEU:HD12	1:A:183:LEU:C	2.42	0.44
1:A:329:ARG:HH11	1:A:329:ARG:HA	1.83	0.44
1:B:170:ARG:O	1:B:171:PRO:C	2.61	0.44
1:A:53:PRO:HD3	1:A:218:THR:HG21	2.00	0.44
1:B:247:LEU:O	1:B:250:LEU:O	2.36	0.44
2:C:96:GLU:HB2	2:C:115:TRP:CZ3	2.53	0.43
1:A:154:LEU:O	1:A:188:CYS:HA	2.17	0.43
1:B:258:LEU:HD12	1:B:258:LEU:HA	1.87	0.43
1:A:58:LYS:N	4:A:502:SO4:O4	2.51	0.43
1:A:142:LEU:O	1:A:146:VAL:HG23	2.18	0.43
1:B:299:THR:O	1:B:303:ASN:HB2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:142:LEU:O	1:B:146:VAL:HG23	2.18	0.43
1:A:349:ARG:CZ	5:A:601:HOH:O	2.65	0.43
1:B:265:TYR:OH	1:B:349:ARG:NH1	2.44	0.43
1:A:383:GLN:O	1:A:387:GLY:HA2	2.18	0.43
1:B:51:CYS:HA	1:B:192:THR:O	2.18	0.43
1:B:75:VAL:HA	1:B:155:VAL:HB	2.01	0.42
1:A:75:VAL:HG11	1:A:127:TYR:CE1	2.54	0.42
1:A:438:THR:HG22	1:A:441[A]:ARG:HH21	1.84	0.42
1:B:80:ILE:HD12	1:B:101:ASN:HB2	2.01	0.42
1:B:329:ARG:HH11	1:B:329:ARG:HA	1.84	0.42
1:A:151:LEU:HD21	1:A:183:LEU:HD22	1.99	0.42
1:A:165:TRP:HE3	1:A:202:ASP:OD2	2.02	0.42
1:B:165:TRP:HE3	1:B:202:ASP:OD2	2.03	0.42
1:B:197:PRO:O	1:B:201:GLU:HG2	2.20	0.42
1:B:442:ARG:NH2	5:B:601:HOH:O	2.53	0.42
1:A:120:LYS:N	1:A:121:PRO:HD3	2.35	0.42
2:K:86:LYS:HB3	2:K:88:GLU:OE1	2.20	0.41
1:A:87:VAL:HG23	1:A:97:VAL:HG22	2.02	0.41
1:B:291:ALA:HB2	1:B:319:PHE:CD1	2.56	0.41
2:K:2:VAL:HG21	2:K:102:THR:HG21	2.03	0.41
1:A:48:VAL:HG13	1:A:189:VAL:HG22	2.03	0.41
1:A:364:SER:O	1:A:367:ASP:HB2	2.19	0.41
1:A:189:VAL:HG12	1:A:191:LEU:HD11	2.03	0.41
1:A:35:GLU:O	1:A:39:MET:HG2	2.21	0.41
1:B:321:MET:HE3	1:B:321:MET:HB3	1.98	0.41
1:A:68:LEU:HD23	1:A:68:LEU:HA	1.96	0.41
1:B:47:ASP:OD1	1:B:210:LYS:HE2	2.20	0.41
1:B:274:GLN:O	1:B:277:ILE:HB	2.21	0.41
1:B:364:SER:O	1:B:367:ASP:HB2	2.20	0.41
2:C:100:THR:HB	2:C:114:TYR:HE2	1.84	0.41
1:B:405:PHE:CE2	1:B:412:ARG:HB3	2.56	0.40
1:B:75:VAL:HG11	1:B:127:TYR:CE1	2.57	0.40
2:K:96:GLU:HB2	2:K:115:TRP:CZ3	2.55	0.40
1:A:148:ARG:HE	1:A:150:LEU:HD12	1.87	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	442/445 (99%)	415 (94%)	27 (6%)	0	100	100
1	B	442/445 (99%)	414 (94%)	28 (6%)	0	100	100
2	C	124/127 (98%)	120 (97%)	4 (3%)	0	100	100
2	K	124/127 (98%)	119 (96%)	5 (4%)	0	100	100
All	All	1132/1144 (99%)	1068 (94%)	64 (6%)	0	100	100

There are no Ramachandran outliers to report.

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	363/373 (97%)	348 (96%)	15 (4%)	27	60
1	B	368/373 (99%)	354 (96%)	14 (4%)	29	62
2	C	100/100 (100%)	95 (95%)	5 (5%)	22	55
2	K	100/100 (100%)	94 (94%)	6 (6%)	17	50
All	All	931/946 (98%)	891 (96%)	40 (4%)	26	59

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	48	VAL
1	A	54	THR

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Mol	Chain	Res	Type
1	A	79	LEU
1	A	80	ILE
1	A	98	SER
1	A	104	LEU
1	A	113	LEU
1	A	132	MET
1	A	141	THR
1	A	152	SER
1	A	218	THR
1	A	345	GLN
1	A	397	MET
1	A	409	LEU
1	A	434	CYS
1	B	21	LYS
1	B	28	SER
1	B	31	THR
1	B	80	ILE
1	B	98	SER
1	B	104	LEU
1	B	119	GLU
1	B	141	THR
1	B	183	LEU
1	B	218	THR
1	B	247	LEU
1	B	397	MET
1	B	409	LEU
1	B	444	LEU
2	K	5	VAL
2	K	18	LEU
2	K	22	CYS
2	K	68	THR
2	K	108	GLU
2	K	119	THR
2	C	61	ASP
2	C	68	THR
2	C	108	GLU
2	C	110	GLU
2	C	119	THR

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

Mol	Chain	Res	Type
1	A	63	GLN
1	A	139	GLN
1	A	285	ASN
1	A	302	GLN
1	B	63	GLN
1	B	149	HIS
1	B	285	ASN
1	B	370	GLN
2	K	3	GLN
2	K	39	GLN
2	C	3	GLN
2	C	39	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](#)

Of 5 ligands modelled in this entry, 2 are monoatomic - leaving 3 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	$\# Z > 2$
4	SO4	A	502	-	4,4,4	0.35	0	6,6,6	0.13	0
4	SO4	A	503	-	4,4,4	0.56	0	6,6,6	0.20	0
4	SO4	B	502	-	4,4,4	0.30	0	6,6,6	0.16	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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	502	SO4	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

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	440/445 (98%)	0.35	19 (4%)	40	24	28, 78, 141, 198	2 (0%)
1	B	440/445 (98%)	0.34	17 (3%)	43	27	27, 74, 127, 162	4 (0%)
2	C	124/127 (97%)	-0.10	0	100	100	32, 49, 82, 122	2 (1%)
2	K	124/127 (97%)	-0.14	0	100	100	23, 48, 77, 106	2 (1%)
All	All	1128/1144 (98%)	0.24	36 (3%)	50	31	23, 68, 129, 198	10 (0%)

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	322	GLY	4.7
1	B	167[A]	HIS	4.6
1	A	12	PRO	4.5
1	B	317	ILE	3.2
1	A	254	ALA	3.1
1	B	300	LEU	3.1
1	A	390	ALA	2.9
1	B	12	PRO	2.9
1	A	165	TRP	2.9
1	A	382	LEU	2.7
1	A	285	ASN	2.7
1	A	336	ILE	2.6
1	A	373	PHE	2.6
1	B	234	LEU	2.6
1	A	317	ILE	2.5
1	A	374	LEU	2.5
1	A	319	PHE	2.5
1	A	379	VAL	2.5
1	B	390	ALA	2.4
1	A	391	SER	2.4
1	A	238	PRO	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	258	LEU	2.3
1	A	198[A]	GLN	2.3
1	B	409	LEU	2.3
1	B	79	LEU	2.3
1	A	444	LEU	2.2
1	B	254	ALA	2.2
1	B	306	MET	2.2
1	B	185	HIS	2.2
1	B	165	TRP	2.1
1	B	437	PRO	2.1
1	A	321	MET	2.1
1	B	260	GLY	2.1
1	B	441[A]	ARG	2.0
1	B	322	GLY	2.0
1	B	80	ILE	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
4	SO4	A	503	5/5	0.85	0.18	66,67,70,70	0
4	SO4	A	502	5/5	0.97	0.09	62,63,67,69	0
4	SO4	B	502	5/5	0.98	0.04	45,51,55,58	0
3	ZN	B	501	1/1	1.00	0.02	70,70,70,70	0
3	ZN	A	501	1/1	1.00	0.02	89,89,89,89	0

6.5 Other polymers [i](#)

There are no such residues in this entry.