



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 7, 2020 – 02:37 PM BST

PDB ID : 6RUU
Title : Pseudokinase domain of human IRAK3
Authors : Lange, S.M.; Kulathu, Y.; Cohen, P.
Deposited on : 2019-05-29
Resolution : 2.95 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.14.2
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) : 2.14.2

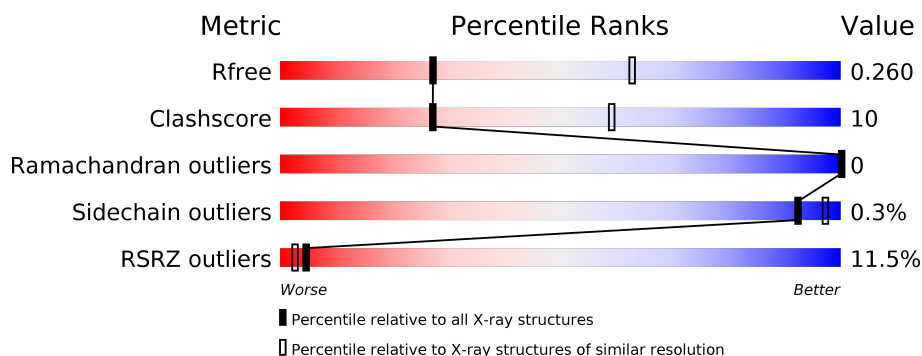
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.95 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	3104 (3.00-2.92)
Clashscore	141614	3462 (3.00-2.92)
Ramachandran outliers	138981	3340 (3.00-2.92)
Sidechain outliers	138945	3343 (3.00-2.92)
RSRZ outliers	127900	2986 (3.00-2.92)

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 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	315	<div> <div>3%</div> <div> <div></div> <div>73%</div> <div>17%</div> <div>10%</div> </div> </div>
1	B	315	<div> <div>20%</div> <div> <div></div> <div>69%</div> <div>22%</div> <div>9%</div> </div> </div>
1	C	315	<div> <div>6%</div> <div> <div></div> <div>60%</div> <div>17%</div> <div>22%</div> </div> </div>

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	SO4	C	508	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6252 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 Interleukin-1 receptor-associated kinase 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	285	Total	C	N	O	S	0	0	0
			2223	1438	376	392	17			
1	B	286	Total	C	N	O	S	0	0	0
			2102	1356	351	379	16			
1	C	245	Total	C	N	O	S	0	0	0
			1861	1205	311	330	15			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	140	GLY	-	expression tag	UNP Q9Y616
A	141	PRO	-	expression tag	UNP Q9Y616
A	142	LEU	-	expression tag	UNP Q9Y616
A	143	GLY	-	expression tag	UNP Q9Y616
A	144	SER	-	expression tag	UNP Q9Y616
A	147	VAL	ILE	variant	UNP Q9Y616
B	140	GLY	-	expression tag	UNP Q9Y616
B	141	PRO	-	expression tag	UNP Q9Y616
B	142	LEU	-	expression tag	UNP Q9Y616
B	143	GLY	-	expression tag	UNP Q9Y616
B	144	SER	-	expression tag	UNP Q9Y616
B	147	VAL	ILE	variant	UNP Q9Y616
C	140	GLY	-	expression tag	UNP Q9Y616
C	141	PRO	-	expression tag	UNP Q9Y616
C	142	LEU	-	expression tag	UNP Q9Y616
C	143	GLY	-	expression tag	UNP Q9Y616
C	144	SER	-	expression tag	UNP Q9Y616
C	147	VAL	ILE	variant	UNP Q9Y616

- Molecule 2 is MERCURY (II) ION (three-letter code: HG) (formula: Hg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Hg 1 1	0	0
2	A	2	Total Hg 2 2	0	0
2	C	5	Total Hg 5 5	0	0

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0
3	C	1	Total C O 6 3 3	0	0

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O_4S).

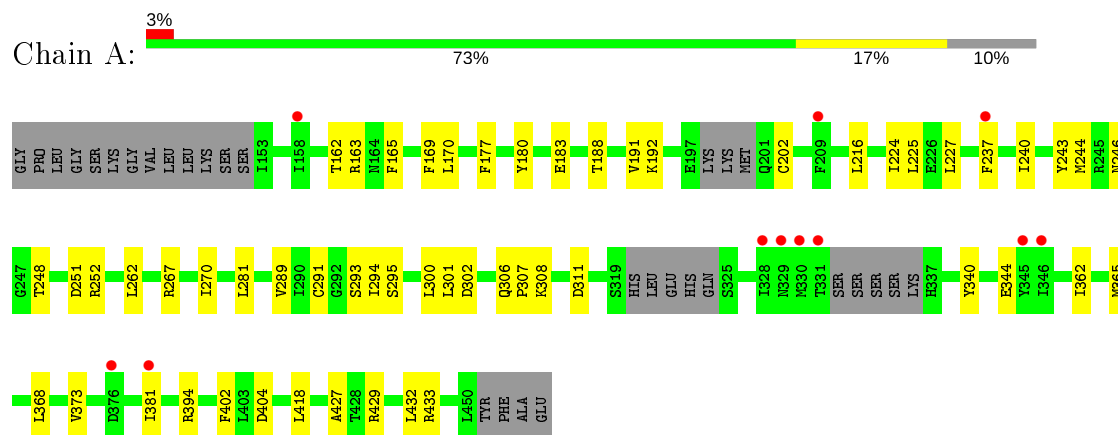


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		

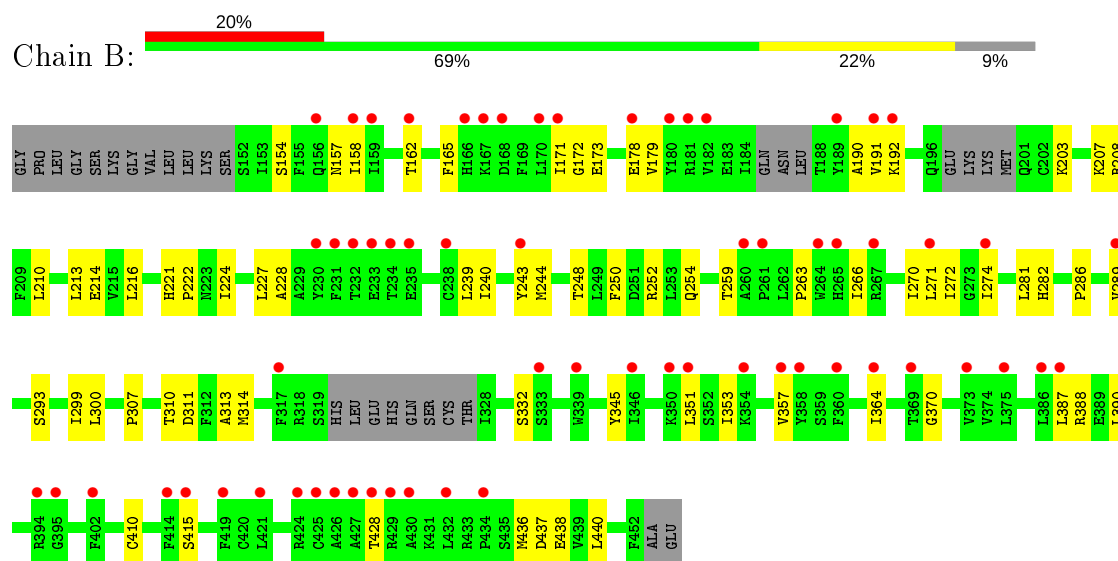
3 Residue-property plots [i](#)

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: Interleukin-1 receptor-associated kinase 3

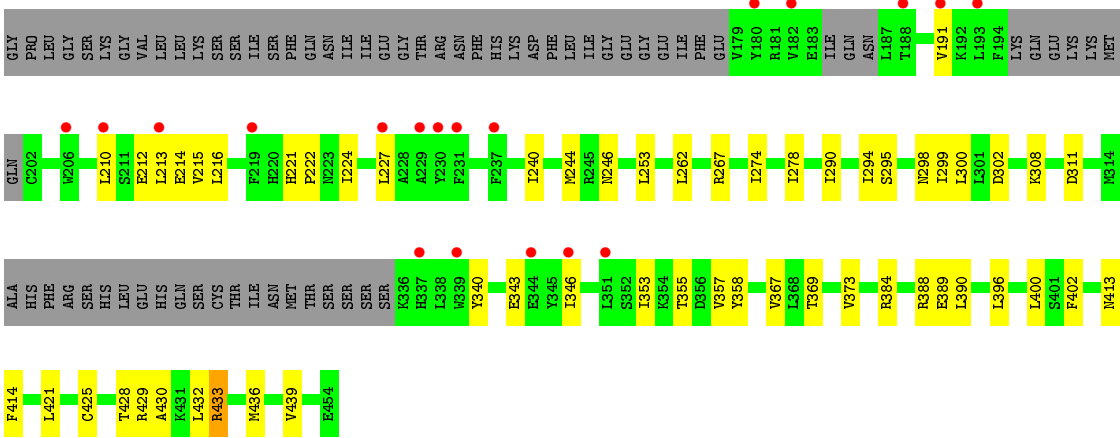


- Molecule 1: Interleukin-1 receptor-associated kinase 3



- Molecule 1: Interleukin-1 receptor-associated kinase 3





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	53.55Å 167.16Å 179.51Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.65 – 2.95 49.06 – 2.95	Depositor EDS
% Data completeness (in resolution range)	61.6 (48.65-2.95) 57.4 (49.06-2.95)	Depositor EDS
R_{merge}	0.28	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.16 (at 2.96Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.228 , 0.260 0.228 , 0.260	Depositor DCC
R_{free} test set	1119 reflections (5.20%)	wwPDB-VP
Wilson B-factor (Å ²)	70.4	Xtriage
Anisotropy	0.115	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 73.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	6252	wwPDB-VP
Average B, all atoms (Å ²)	90.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.61% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.27	0/2272	0.47	0/3076
1	B	0.25	0/2147	0.44	0/2927
1	C	0.27	0/1902	0.46	0/2588
All	All	0.26	0/6321	0.46	0/8591

There are no bond length outliers.

There are no bond angle outliers.

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	2223	0	2162	43	0
1	B	2102	0	1935	43	0
1	C	1861	0	1786	43	0
2	A	2	0	0	0	0
2	B	1	0	0	0	0
2	C	5	0	0	0	0
3	A	12	0	16	0	0
3	C	6	0	8	0	0
4	A	20	0	0	0	0
4	B	10	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	10	0	0	3	0
All	All	6252	0	5907	123	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 10.

All (123) 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:162:THR:HG23	1:B:165:PHE:H	1.43	0.83
1:B:252:ARG:HE	1:B:259:THR:HG21	1.48	0.79
1:C:191:VAL:HG12	1:C:240:ILE:HG22	1.65	0.78
1:C:430:ALA:HA	1:C:433:ARG:HE	1.49	0.76
1:A:295:SER:HA	1:A:340:TYR:HE2	1.49	0.75
1:B:179:VAL:HG12	1:B:192:LYS:HG2	1.67	0.74
1:A:362:ILE:HA	1:A:365:MET:HE3	1.70	0.73
1:A:381:ILE:HD11	1:C:389:GLU:HG3	1.71	0.71
1:B:263:PRO:HD2	1:B:266:ILE:HD12	1.73	0.70
1:A:295:SER:HA	1:A:340:TYR:CE2	2.27	0.70
1:B:228:ALA:HB3	1:B:240:ILE:HG13	1.75	0.69
1:A:246:ASN:ND2	1:A:302:ASP:O	2.25	0.68
1:C:246:ASN:ND2	1:C:302:ASP:O	2.29	0.66
1:A:162:THR:HG23	1:A:165:PHE:H	1.59	0.66
1:C:244:MET:HE3	1:C:308:LYS:HG3	1.78	0.66
1:A:163:ARG:HD3	1:A:169:PHE:HE2	1.60	0.66
1:B:171:ILE:HB	1:B:179:VAL:HG23	1.77	0.65
1:C:221:HIS:CD2	1:C:222:PRO:HD2	2.32	0.64
1:A:402:PHE:HE2	1:C:390:LEU:HD23	1.63	0.63
1:A:427:ALA:O	1:A:433:ARG:NH1	2.32	0.62
1:C:421:LEU:HD21	1:C:439:VAL:HG23	1.82	0.61
1:A:191:VAL:HG12	1:A:240:ILE:HG22	1.82	0.61
1:A:394:ARG:NH2	1:C:369:THR:O	2.34	0.60
1:B:270:ILE:HG22	1:B:307:PRO:HD3	1.83	0.60
1:C:430:ALA:HA	1:C:433:ARG:NE	2.18	0.59
1:A:163:ARG:HD3	1:A:169:PHE:CE2	2.38	0.59
1:B:410:CYS:SG	1:B:415:SER:OG	2.61	0.58
1:A:293:SER:OG	1:A:311:ASP:O	2.22	0.58
1:B:274:ILE:HD11	1:B:307:PRO:HB2	1.86	0.58
1:A:252:ARG:HB3	1:A:262:LEU:HD11	1.86	0.58
1:B:224:ILE:HD11	1:B:281:LEU:HD21	1.85	0.58
1:A:237:PHE:CE2	1:B:286:PRO:HG3	2.40	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:213:LEU:HD12	1:B:239:LEU:HD21	1.88	0.56
1:A:270:ILE:HG23	1:A:307:PRO:HB3	1.89	0.55
1:B:388:ARG:HH22	1:B:428:THR:HG23	1.72	0.54
1:B:173:GLU:HA	1:B:178:GLU:HA	1.89	0.54
1:B:191:VAL:HG12	1:B:240:ILE:HG22	1.90	0.54
1:A:402:PHE:CE2	1:C:390:LEU:HD23	2.43	0.54
1:A:248:THR:HG23	1:A:251:ASP:H	1.73	0.53
1:B:208:ARG:HB3	1:B:313:ALA:HB1	1.89	0.52
1:B:158:ILE:O	1:B:162:THR:HG22	2.09	0.52
1:B:437:ASP:OD1	1:B:438:GLU:N	2.42	0.52
1:C:253:LEU:HD21	1:C:367:VAL:HG23	1.91	0.52
1:C:358:TYR:HB2	1:C:433:ARG:NH2	2.25	0.52
1:A:225:LEU:HD22	1:A:300:LEU:HD12	1.92	0.52
1:B:282:HIS:HD2	1:B:353:ILE:HG23	1.74	0.51
1:B:387:LEU:HD23	1:B:390:LEU:HD12	1.91	0.51
1:A:281:LEU:HB3	1:A:289:VAL:HG13	1.92	0.51
1:A:344:GLU:OE2	1:A:433:ARG:NH2	2.43	0.51
1:C:295:SER:HA	1:C:340:TYR:CE2	2.44	0.51
1:B:271:LEU:HD11	1:B:364:ILE:HG12	1.93	0.51
1:A:183:GLU:HA	1:A:188:THR:HA	1.92	0.50
1:C:210:LEU:O	1:C:214:GLU:HG3	2.11	0.50
1:B:208:ARG:NH2	1:B:332:SER:O	2.41	0.50
1:A:188:THR:HG23	1:A:243:TYR:HB3	1.93	0.50
1:A:394:ARG:HG3	1:C:402:PHE:HA	1.93	0.50
1:A:267:ARG:NH2	1:A:404:ASP:OD2	2.45	0.49
1:A:224:ILE:HD11	1:A:281:LEU:HD11	1.94	0.49
1:C:295:SER:HA	1:C:340:TYR:HE2	1.77	0.49
1:C:357:VAL:HG22	1:C:436:MET:HE2	1.95	0.49
1:B:244:MET:HG3	1:B:300:LEU:HB3	1.95	0.49
1:A:162:THR:HA	1:A:169:PHE:HD2	1.77	0.49
1:C:413:ASN:OD1	1:C:414:PHE:N	2.46	0.49
1:A:244:MET:HE3	1:A:308:LYS:HG3	1.95	0.49
1:C:221:HIS:HB3	1:C:224:ILE:HG12	1.93	0.48
1:C:358:TYR:HA	1:C:425:CYS:O	2.13	0.48
1:C:384:ARG:O	1:C:388:ARG:HG2	2.14	0.48
1:C:388:ARG:NH1	1:C:428:THR:HG23	2.29	0.48
1:A:252:ARG:HB3	1:A:262:LEU:CD1	2.43	0.48
1:C:212:GLU:OE1	1:C:311:ASP:HB3	2.13	0.48
1:A:177:PHE:CD2	1:A:192:LYS:HG2	2.50	0.47
1:C:340:TYR:CE1	1:C:373:VAL:HG21	2.49	0.47
1:B:216:LEU:HB3	1:B:227:LEU:HB2	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:278:ILE:HG21	1:C:294:ILE:HD12	1.95	0.47
1:B:357:VAL:HG12	1:B:436:MET:SD	2.55	0.47
1:B:248:THR:HG22	1:B:250:PHE:H	1.79	0.47
1:B:254:GLN:NE2	1:B:370:GLY:O	2.48	0.47
1:A:340:TYR:CE1	1:A:373:VAL:HG21	2.50	0.46
1:A:429:ARG:HG3	1:A:432:LEU:HD23	1.98	0.46
1:C:429:ARG:HG3	1:C:432:LEU:HD13	1.98	0.46
1:A:216:LEU:HD13	1:A:227:LEU:HB2	1.98	0.46
1:B:154:SER:OG	1:B:157:ASN:OD1	2.23	0.46
1:B:274:ILE:HG13	1:B:299:ILE:HD13	1.97	0.46
1:C:290:ILE:HG21	1:C:353:ILE:HA	1.98	0.46
1:C:300:LEU:HB2	1:C:308:LYS:HB2	1.98	0.45
1:B:274:ILE:CD1	1:B:307:PRO:HB2	2.46	0.45
1:C:210:LEU:O	1:C:214:GLU:N	2.45	0.45
1:C:388:ARG:NH2	4:C:508:SO4:O1	2.28	0.45
1:C:213:LEU:HD12	1:C:227:LEU:HD13	1.98	0.45
1:A:180:TYR:HB2	1:A:191:VAL:CG2	2.47	0.44
1:B:248:THR:HG22	1:B:250:PHE:N	2.32	0.44
1:C:429:ARG:NE	4:C:508:SO4:O2	2.50	0.44
1:A:340:TYR:HE1	1:A:373:VAL:HG11	1.83	0.44
1:B:210:LEU:O	1:B:214:GLU:HG3	2.18	0.44
1:B:221:HIS:CG	1:B:222:PRO:HD2	2.52	0.44
1:A:302:ASP:OD1	1:A:306:GLN:N	2.42	0.44
1:B:203:LYS:O	1:B:207:LYS:HD3	2.17	0.44
1:B:208:ARG:HE	1:B:314:MET:HE3	1.83	0.44
1:C:262:LEU:HB2	1:C:267:ARG:HE	1.83	0.44
1:B:281:LEU:HD12	1:B:289:VAL:HG11	1.99	0.44
1:C:396:LEU:O	1:C:400:LEU:HG	2.17	0.43
1:B:345:TYR:CE1	1:B:351:LEU:HD12	2.54	0.43
1:C:215:VAL:HG13	1:C:216:LEU:HD12	2.00	0.43
1:C:343:GLU:HA	1:C:346:ILE:HG12	2.00	0.43
1:A:362:ILE:HD13	1:A:365:MET:HE3	2.01	0.42
1:A:170:LEU:HA	1:A:180:TYR:HA	2.01	0.42
1:A:368:LEU:HD21	1:A:418:LEU:HD22	2.00	0.42
1:B:310:THR:HG22	1:B:311:ASP:N	2.35	0.42
1:A:270:ILE:HA	1:A:307:PRO:HG3	2.02	0.42
1:B:172:GLY:O	1:B:179:VAL:N	2.52	0.42
1:B:216:LEU:HD13	1:B:227:LEU:HB2	2.02	0.42
1:B:190:ALA:HB2	1:B:243:TYR:HD1	1.83	0.42
1:A:291:CYS:HB2	1:A:294:ILE:HD11	2.02	0.42
1:A:252:ARG:HG3	1:A:301:LEU:HD12	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:272:ILE:HG23	1:B:440:LEU:HD11	2.02	0.41
1:A:368:LEU:HD11	1:A:418:LEU:HB3	2.03	0.41
1:C:353:ILE:H	1:C:353:ILE:HD12	1.86	0.41
1:C:222:PRO:O	1:C:308:LYS:NZ	2.36	0.41
1:C:355:THR:O	1:C:358:TYR:HB3	2.20	0.41
1:C:428:THR:N	4:C:508:SO4:O4	2.54	0.41
1:C:298:ASN:ND2	1:C:311:ASP:O	2.54	0.40
1:B:293:SER:OG	1:B:311:ASP:O	2.40	0.40
1:C:274:ILE:HD13	1:C:299:ILE:HD13	2.03	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	277/315 (88%)	264 (95%)	13 (5%)	0	100	100
1	B	278/315 (88%)	268 (96%)	10 (4%)	0	100	100
1	C	237/315 (75%)	228 (96%)	9 (4%)	0	100	100
All	All	792/945 (84%)	760 (96%)	32 (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	233/284 (82%)	232 (100%)	1 (0%)	91	96
1	B	204/284 (72%)	204 (100%)	0	100	100
1	C	191/284 (67%)	190 (100%)	1 (0%)	88	95
All	All	628/852 (74%)	626 (100%)	2 (0%)	92	97

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	202	CYS
1	C	433	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	221	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 19 ligands modelled in this entry, 8 are monoatomic - leaving 11 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	GOL	A	503	-	5,5,5	0.92	0	5,5,5	0.98	0
4	SO4	A	506	-	4,4,4	0.15	0	6,6,6	0.04	0
4	SO4	B	502	-	4,4,4	0.14	0	6,6,6	0.08	0
3	GOL	C	506	-	5,5,5	0.96	0	5,5,5	0.95	0
4	SO4	C	507	-	4,4,4	0.15	0	6,6,6	0.07	0
4	SO4	C	508	-	4,4,4	0.13	0	6,6,6	0.10	0
4	SO4	A	507	-	4,4,4	0.14	0	6,6,6	0.07	0
4	SO4	B	503	-	4,4,4	0.14	0	6,6,6	0.06	0
3	GOL	A	504	-	5,5,5	0.90	0	5,5,5	0.99	0
4	SO4	A	505	-	4,4,4	0.14	0	6,6,6	0.08	0
4	SO4	A	508	-	4,4,4	0.15	0	6,6,6	0.06	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
3	GOL	A	503	-	-	0/4/4/4	-
3	GOL	C	506	-	-	2/4/4/4	-
3	GOL	A	504	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	506	GOL	C1-C2-C3-O3
3	C	506	GOL	O2-C2-C3-O3

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	508	SO4	3	0

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	285/315 (90%)	0.27	11 (3%) 39 25	30, 69, 126, 171	0
1	B	286/315 (90%)	1.02	64 (22%) 0 0	83, 123, 167, 196	0
1	C	245/315 (77%)	0.48	19 (7%) 13 7	28, 60, 125, 150	0
All	All	816/945 (86%)	0.59	94 (11%) 4 3	28, 96, 151, 196	0

All (94) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	234	THR	9.0
1	C	229	ALA	6.9
1	B	425	CYS	5.8
1	C	230	TYR	5.6
1	A	329	ASN	5.5
1	B	264	TRP	5.5
1	B	191	VAL	5.3
1	B	421	LEU	5.1
1	B	267	ARG	5.0
1	B	394	ARG	5.0
1	B	162	THR	4.7
1	A	331	THR	4.7
1	B	415	SER	4.7
1	A	330	MET	4.5
1	B	181	ARG	4.5
1	B	351	LEU	4.0
1	B	180	TYR	3.9
1	B	360	PHE	3.9
1	C	337	HIS	3.8
1	B	170	LEU	3.8
1	A	237	PHE	3.8
1	B	354	LYS	3.7
1	B	168	ASP	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	369	THR	3.6
1	C	188	THR	3.6
1	C	206	TRP	3.5
1	C	227	LEU	3.5
1	C	193	LEU	3.5
1	B	414	PHE	3.5
1	A	376	ASP	3.5
1	B	158	ILE	3.4
1	B	238	CYS	3.3
1	C	191	VAL	3.2
1	B	430	ALA	3.2
1	A	328	ILE	3.1
1	B	419	PHE	3.1
1	B	171	ILE	3.1
1	B	289	VAL	3.1
1	C	219	PHE	3.0
1	C	210	LEU	3.0
1	B	167	LYS	3.0
1	B	357	VAL	3.0
1	B	271	LEU	3.0
1	C	346	ILE	2.9
1	B	189	TYR	2.8
1	B	192	LYS	2.8
1	C	213	LEU	2.8
1	B	432	LEU	2.8
1	B	424	ARG	2.8
1	B	373	VAL	2.8
1	A	158	ILE	2.7
1	B	182	VAL	2.7
1	B	231	PHE	2.7
1	B	426	ALA	2.7
1	A	346	ILE	2.6
1	B	232	THR	2.6
1	B	402	PHE	2.6
1	B	166	HIS	2.6
1	C	339	TRP	2.6
1	B	428	THR	2.6
1	B	387	LEU	2.6
1	B	159	ILE	2.6
1	B	178	GLU	2.5
1	B	233	GLU	2.5
1	B	230	TYR	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	358	TYR	2.4
1	C	180	TYR	2.4
1	B	350	LYS	2.3
1	A	209	PHE	2.3
1	C	182	VAL	2.3
1	C	344	GLU	2.3
1	B	375	LEU	2.3
1	B	395	GLY	2.3
1	B	434	PRO	2.3
1	B	427	ALA	2.2
1	B	346	ILE	2.2
1	B	265	HIS	2.2
1	B	364	ILE	2.2
1	C	237	PHE	2.2
1	B	261	PRO	2.2
1	A	381	ILE	2.2
1	B	339	TRP	2.2
1	B	260	ALA	2.2
1	C	351	LEU	2.2
1	B	274	ILE	2.2
1	C	231	PHE	2.1
1	A	345	TYR	2.1
1	B	156	GLN	2.1
1	B	235	GLU	2.1
1	B	429	ARG	2.1
1	B	243	TYR	2.0
1	B	317	PHE	2.0
1	B	333	SER	2.0
1	B	386	LEU	2.0

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 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	SO4	B	502	5/5	0.78	0.20	82,97,104,120	0
3	GOL	A	503	6/6	0.81	0.29	90,97,102,108	0
4	SO4	A	508	5/5	0.83	0.17	130,137,144,151	0
3	GOL	C	506	6/6	0.84	0.28	66,86,93,94	0
4	SO4	A	507	5/5	0.85	0.21	119,126,128,134	0
4	SO4	B	503	5/5	0.91	0.31	125,135,143,145	0
2	HG	C	504	1/1	0.91	0.15	162,162,162,162	1
3	GOL	A	504	6/6	0.91	0.16	73,75,77,85	0
2	HG	C	501	1/1	0.92	0.09	78,78,78,78	1
4	SO4	C	507	5/5	0.93	0.14	69,77,96,103	0
2	HG	A	501	1/1	0.95	0.14	100,100,100,100	1
2	HG	C	505	1/1	0.95	0.09	130,130,130,130	1
4	SO4	C	508	5/5	0.97	0.10	54,62,67,75	0
4	SO4	A	505	5/5	0.97	0.14	59,67,69,69	0
2	HG	C	503	1/1	0.98	0.12	114,114,114,114	0
4	SO4	A	506	5/5	0.98	0.14	62,67,71,73	0
2	HG	C	502	1/1	0.98	0.14	56,56,56,56	1
2	HG	B	501	1/1	0.98	0.07	74,74,74,74	1
2	HG	A	502	1/1	0.99	0.15	73,73,73,73	1

6.5 Other polymers [i](#)

There are no such residues in this entry.