Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:19 PM GMT

PDB ID	:	4OK2
Title	:	Crystal Structure of Alg17c Mutant Y258A
Authors	:	Nair, S.K.; Park, D.S.
Deposited on	:	2014-01-21
Resolution	:	2.45 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
http://wwpdb.org/validation/2016/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:

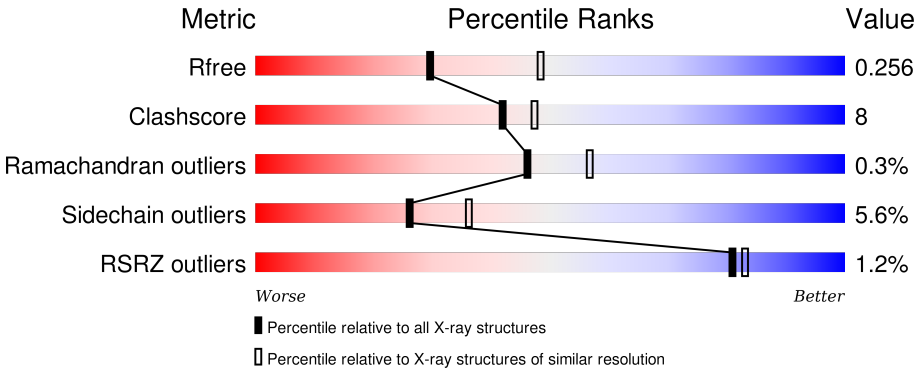
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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}	91344	4776 (2.50-2.42)
Clashscore	102246	1030 (2.48-2.44)
Ramachandran outliers	100387	1024 (2.48-2.44)
Sidechain outliers	100360	1024 (2.48-2.44)
RSRZ outliers	91569	4787 (2.50-2.42)

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. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	736	% 79% 15%
1	B	736	% 78% 16%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11521 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 Putative alginate lyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	706	Total	C	N	O	S	0	0	0
			5549	3531	938	1067	13			
1	B	705	Total	C	N	O	S	0	0	0
			5539	3525	935	1066	13			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	258	ALA	TYR	ENGINEERED MUTATION	UNP Q21FJ0
B	258	ALA	TYR	ENGINEERED MUTATION	UNP Q21FJ0

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 3 is water.

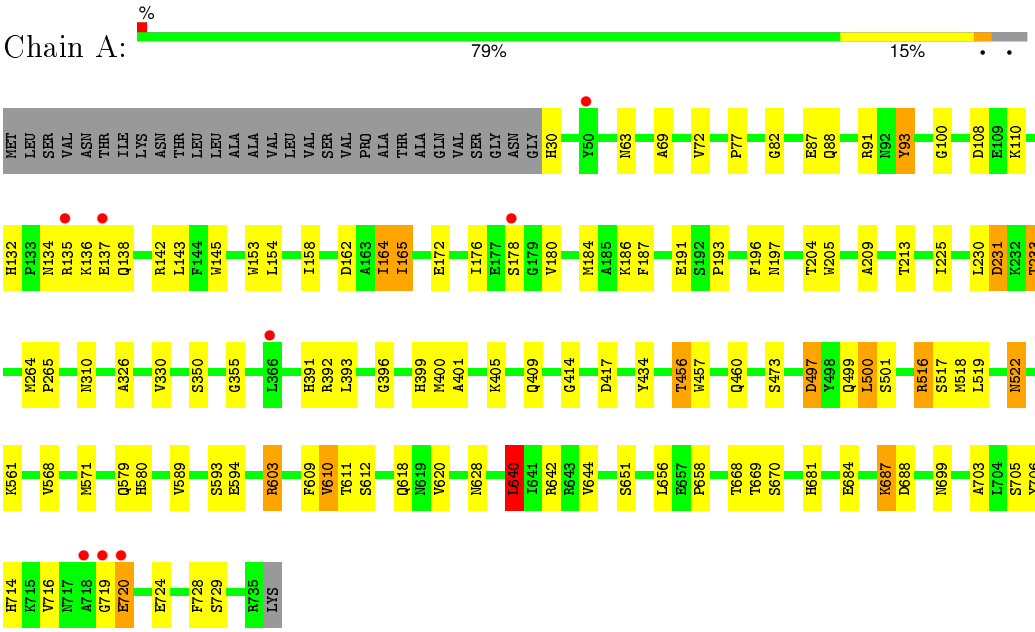
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	245	Total	O	0	0
			245	245		
3	B	186	Total	O	0	0
			186	186		

3 Residue-property plots (i)

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.

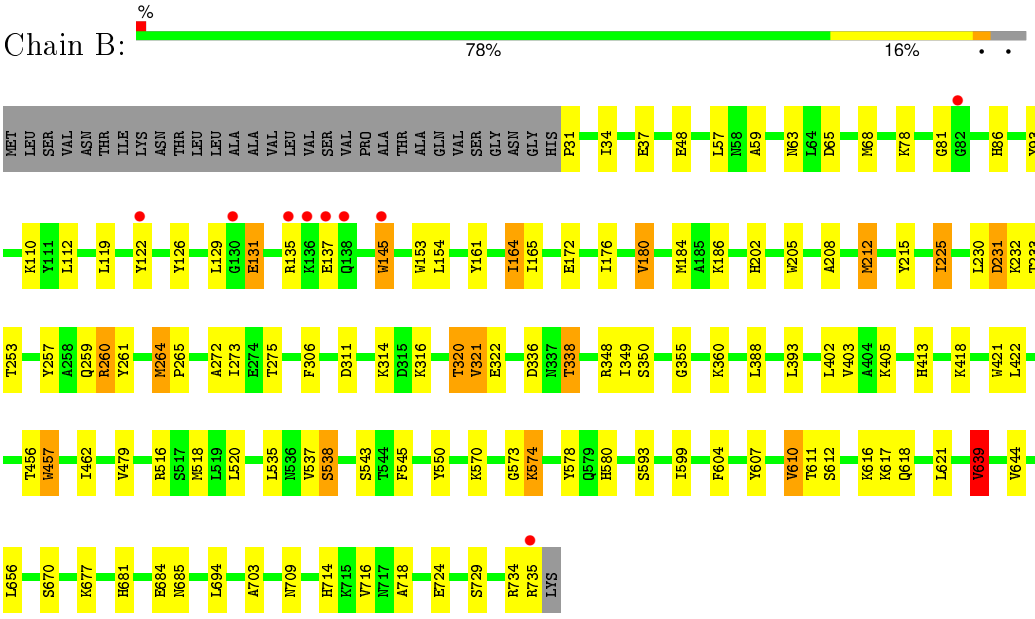
- Molecule 1: Putative alginate lyase

Chain A:



- Molecule 1: Putative alginate lyase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	84.95Å 136.63Å 88.86Å 90.00° 107.11° 90.00°	Depositor
Resolution (Å)	25.00 – 2.45 48.84 – 2.44	Depositor EDS
% Data completeness (in resolution range)	93.2 (25.00-2.45) 92.8 (48.84-2.44)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.57 (at 2.45Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.193 , 0.256 0.191 , 0.256	Depositor DCC
R_{free} test set	3316 reflections (5.27%)	DCC
Wilson B-factor (Å^2)	35.5	Xtriage
Anisotropy	0.070	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.29 , 29.5	EDS
Estimated twinning fraction	0.022 for l,-k,h	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Outliers	1 of 66619 reflections (0.002%)	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	11521	wwPDB-VP
Average B, all atoms (Å^2)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 14.29% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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

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	0.48	0/5677	0.65	2/7705 (0.0%)
1	B	0.48	0/5666	0.65	1/7689 (0.0%)
All	All	0.48	0/11343	0.65	3/15394 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	639	VAL	CB-CA-C	-5.88	100.23	111.40
1	A	535	LEU	CA-CB-CG	5.69	128.38	115.30
1	A	640	LEU	CA-CB-CG	5.04	126.89	115.30

There are no chirality outliers.

There are no planarity outliers.

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	5549	0	5388	93	0
1	B	5539	0	5382	91	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	245	0	0	4	0
3	B	186	0	0	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	11521	0	10770	183	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 (183) 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:A:603:ARG:HH11	1:A:603:ARG:HG3	1.01	1.15
1:A:456:THR:CG2	1:A:580:HIS:NE2	2.17	1.08
1:B:612:SER:OG	1:B:681:HIS:HD2	1.50	0.92
1:A:393:LEU:HD21	1:A:522:ASN:HD21	1.33	0.91
1:A:603:ARG:NH1	1:A:603:ARG:HG3	1.81	0.88
1:B:421:TRP:HB3	1:B:518:MET:HE3	1.58	0.84
1:B:100:GLY:HA2	1:B:164:ILE:HD13	1.59	0.82
1:B:260:ARG:HG3	1:B:261:TYR:N	1.95	0.81
1:A:456:THR:HG23	1:A:580:HIS:NE2	1.98	0.78
1:B:208:ALA:O	1:B:212:MET:CG	2.34	0.76
1:A:393:LEU:HD21	1:A:522:ASN:ND2	2.00	0.75
1:A:612:SER:OG	1:A:681:HIS:HD2	1.69	0.75
1:A:571:MET:CE	1:A:628:ASN:OD1	2.33	0.75
1:B:456:THR:HG22	1:B:580:HIS:NE2	2.03	0.74
1:A:132:HIS:CD2	1:A:134:ASN:H	2.07	0.73
1:B:208:ALA:O	1:B:212:MET:HG2	1.91	0.71
1:B:418:LYS:HE2	3:B:986:HOH:O	1.90	0.70
1:A:392:ARG:HH11	1:A:399:HIS:HE1	1.38	0.70
1:A:603:ARG:HH11	1:A:603:ARG:CG	1.92	0.69
1:B:336:ASP:OD1	1:B:338:THR:HB	1.91	0.69
1:B:316:LYS:HZ3	1:B:320:THR:HG23	1.56	0.69
1:A:456:THR:HG22	1:A:580:HIS:NE2	2.06	0.68
1:A:589:VAL:HG22	1:A:594:GLU:HG3	1.75	0.67
1:B:421:TRP:HB3	1:B:518:MET:CE	2.24	0.67
1:B:253:THR:HG22	3:B:916:HOH:O	1.94	0.66
1:A:100:GLY:HA2	1:A:164:ILE:HD13	1.78	0.65
1:B:208:ALA:O	1:B:212:MET:HG3	1.96	0.65
1:A:400:MET:HE1	1:A:658:PRO:HG2	1.79	0.65
1:B:612:SER:OG	1:B:681:HIS:CD2	2.42	0.65
1:B:31:PRO:HD2	1:B:37:GLU:HG3	1.80	0.64
1:A:571:MET:HE1	1:A:628:ASN:OD1	1.97	0.64
1:B:100:GLY:HA2	1:B:164:ILE:CD1	2.28	0.64
1:B:338:THR:HG21	3:B:963:HOH:O	2.00	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:434:TYR:CD2	1:A:457:TRP:HZ3	2.18	0.62
1:B:516:ARG:HD3	1:B:518:MET:HE2	1.82	0.62
1:B:316:LYS:NZ	1:B:320:THR:HG23	2.13	0.62
1:B:112:LEU:HD12	1:B:164:ILE:HG13	1.81	0.61
1:B:456:THR:CG2	1:B:580:HIS:NE2	2.63	0.61
1:A:154:LEU:HD13	1:A:184:MET:CE	2.29	0.61
1:B:65:ASP:HA	1:B:68:MET:HE2	1.80	0.61
1:A:522:ASN:HD22	1:A:522:ASN:N	1.98	0.61
1:A:571:MET:HE2	1:A:628:ASN:OD1	2.01	0.61
1:A:119:LEU:HB3	1:A:176:ILE:HD12	1.83	0.60
1:A:30:HIS:N	3:A:953:HOH:O	2.34	0.60
1:A:310:ASN:OD1	1:A:405:LYS:NZ	2.34	0.60
1:A:687:LYS:HG3	1:A:705:SER:HA	1.83	0.60
1:A:620:VAL:HG13	1:A:640:LEU:HD13	1.83	0.60
1:A:136:LYS:HB2	1:A:138:GLN:HG2	1.84	0.60
1:A:456:THR:HG21	1:A:580:HIS:NE2	2.13	0.59
1:A:205:TRP:CZ2	1:A:265:PRO:HD3	2.37	0.59
1:A:392:ARG:HH11	1:A:399:HIS:CE1	2.20	0.58
1:A:132:HIS:HD2	1:A:134:ASN:H	1.49	0.58
1:B:119:LEU:HB3	1:B:176:ILE:HD12	1.85	0.58
1:B:714:HIS:O	1:B:724:GLU:HA	2.03	0.57
1:A:593:SER:HA	1:A:610:VAL:HG13	1.86	0.57
1:A:400:MET:CE	1:A:658:PRO:HG2	2.34	0.57
1:A:350:SER:O	1:A:355:GLY:HA3	2.04	0.57
1:B:456:THR:HG22	1:B:580:HIS:CD2	2.40	0.56
1:A:178:SER:HB2	3:A:1120:HOH:O	2.04	0.56
1:B:231:ASP:O	1:B:232:LYS:HB2	2.05	0.56
1:A:204:THR:HG22	1:A:265:PRO:HB2	1.88	0.56
1:B:574:LYS:HG2	3:B:959:HOH:O	2.06	0.56
1:B:593:SER:HA	1:B:610:VAL:HG13	1.86	0.55
1:A:460:GLN:NE2	1:A:473:SER:OG	2.40	0.55
1:B:316:LYS:NZ	1:B:320:THR:CG2	2.69	0.55
1:B:86:HIS:CE1	1:B:145:TRP:HZ3	2.23	0.55
1:B:422:LEU:HD11	1:B:433:ASP:HB2	1.89	0.55
1:B:321:VAL:HB	3:B:934:HOH:O	2.06	0.54
1:A:500:LEU:HG	1:A:501:SER:N	2.21	0.54
1:A:687:LYS:NZ	1:A:714:HIS:HD2	2.05	0.54
1:B:57:LEU:HD11	1:B:102:LEU:HD21	1.89	0.54
1:A:252:TYR:H	1:A:259:GLN:NE2	2.06	0.54
1:A:172:GLU:O	1:A:176:ILE:HG12	2.07	0.53
1:A:568:VAL:HG23	1:B:570:LYS:HG2	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:405:LYS:HE2	1:A:417:ASP:OD1	2.08	0.53
1:A:714:HIS:O	1:A:724:GLU:HA	2.08	0.53
1:B:402:LEU:HD13	1:B:520:LEU:HD11	1.91	0.53
1:B:434:TYR:CD2	1:B:457:TRP:HZ3	2.27	0.53
1:A:225:ILE:HD12	1:A:230:LEU:HA	1.91	0.52
1:A:205:TRP:CH2	1:A:264:MET:HG2	2.45	0.52
1:A:714:HIS:HE1	3:A:923:HOH:O	1.92	0.51
1:A:193:PRO:O	1:A:197:ASN:HB2	2.11	0.51
1:A:516:ARG:HD2	1:A:533:ASP:OD1	2.11	0.51
1:B:257:TYR:O	1:B:260:ARG:HG2	2.10	0.51
1:A:618:GLN:HG2	1:A:644:VAL:HG22	1.92	0.51
1:A:396:GLY:O	1:A:399:HIS:HD2	1.94	0.51
1:B:426:ASN:ND2	3:B:1001:HOH:O	2.44	0.51
1:A:681:HIS:HE1	1:A:688:ASP:OD1	1.95	0.50
1:A:82:GLY:HA2	1:A:87:GLU:OE2	2.11	0.50
1:A:706:TYR:HA	1:A:728:PHE:CE2	2.46	0.50
1:B:537:VAL:HG22	1:B:538:SER:N	2.27	0.49
1:B:86:HIS:CE1	1:B:145:TRP:CZ3	3.00	0.49
1:A:703:ALA:O	1:A:729:SER:HA	2.12	0.49
1:A:231:ASP:OD2	1:A:233:THR:OG1	2.31	0.49
1:B:154:LEU:HB2	1:B:184:MET:HE1	1.94	0.49
1:B:607:TYR:HE1	1:B:654:SER:HB3	1.78	0.49
1:A:154:LEU:HB3	1:A:209:ALA:HB1	1.95	0.49
1:A:196:PHE:CZ	1:A:225:ILE:HG22	2.48	0.48
1:A:497:ASP:OD2	1:A:497:ASP:N	2.43	0.48
1:B:212:MET:HE3	1:B:272:ALA:HB2	1.95	0.48
1:A:609:PHE:HB2	1:A:642:ARG:CZ	2.44	0.48
1:B:709:ASN:HB2	3:B:1033:HOH:O	2.12	0.48
1:B:252:TYR:H	1:B:259:GLN:NE2	2.12	0.48
1:B:122:TYR:HB3	1:B:180:VAL:HG11	1.93	0.48
1:B:348:ARG:HG3	3:B:1031:HOH:O	2.13	0.48
1:A:593:SER:CA	1:A:610:VAL:HG13	2.43	0.48
1:B:231:ASP:HB2	1:B:233:THR:H	1.77	0.47
1:B:68:MET:HE1	1:B:110:LYS:HD3	1.97	0.47
1:B:611:THR:HA	1:B:651:SER:O	2.15	0.47
1:A:326:ALA:O	1:A:330:VAL:HG23	2.14	0.47
1:B:607:TYR:CE1	1:B:654:SER:HB3	2.50	0.47
1:B:611:THR:HB	1:B:618:GLN:OE1	2.15	0.47
1:B:172:GLU:O	1:B:176:ILE:HG12	2.14	0.46
1:A:88:GLN:OE1	1:A:91:ARG:NH1	2.48	0.46
1:B:684:GLU:HB2	1:B:716:VAL:HG13	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:550:TYR:CD1	1:B:639:VAL:HG13	2.49	0.46
1:B:161:TYR:CZ	1:B:165:ILE:HD11	2.50	0.46
1:B:520:LEU:N	1:B:520:LEU:HD12	2.31	0.46
1:A:88:GLN:HA	1:A:91:ARG:NH1	2.31	0.46
1:A:158:ILE:HG12	1:A:213:THR:HA	1.97	0.46
1:B:264:MET:HB3	1:B:265:PRO:CD	2.46	0.46
1:A:154:LEU:HD13	1:A:184:MET:HE3	1.97	0.46
1:B:306:PHE:HD2	1:B:316:LYS:HG3	1.81	0.46
1:B:264:MET:HB3	1:B:265:PRO:HD3	1.97	0.46
1:A:186:LYS:HE3	1:A:191:GLU:OE2	2.16	0.45
1:B:543:SER:HB2	1:B:545:PHE:CE1	2.51	0.45
1:B:403:VAL:HB	1:B:422:LEU:HB2	1.99	0.45
1:B:215:TYR:HE1	1:B:273:ILE:HD13	1.82	0.45
1:B:703:ALA:O	1:B:729:SER:HA	2.16	0.45
1:A:687:LYS:HZ3	1:A:714:HIS:HD2	1.65	0.45
1:A:612:SER:OG	1:A:681:HIS:CD2	2.60	0.45
1:A:603:ARG:NH1	1:A:603:ARG:CG	2.61	0.45
1:B:31:PRO:N	3:B:908:HOH:O	2.49	0.44
1:B:231:ASP:O	1:B:232:LYS:CB	2.66	0.44
1:A:589:VAL:CG2	1:A:594:GLU:HG3	2.42	0.44
1:B:516:ARG:CD	1:B:518:MET:HE2	2.46	0.44
1:A:72:VAL:HG13	1:A:118:LEU:HD23	1.99	0.44
1:A:93:TYR:HB3	1:A:153:TRP:CD1	2.52	0.44
1:B:260:ARG:CD	1:B:316:LYS:HD2	2.46	0.44
1:A:108:ASP:OD1	1:A:110:LYS:HG2	2.17	0.44
1:A:231:ASP:OD1	1:A:231:ASP:N	2.39	0.44
1:B:320:THR:HB	1:B:322:GLU:OE1	2.18	0.44
1:B:34:ILE:HD11	1:B:212:MET:HE1	2.00	0.44
1:B:126:TYR:HA	1:B:129:LEU:HD12	2.00	0.44
1:A:72:VAL:HG21	1:A:117:ASP:HB3	2.00	0.43
1:B:734:ARG:HB3	1:B:735:ARG:H	1.72	0.43
1:A:719:GLY:O	1:A:720:GLU:HB2	2.18	0.43
1:B:311:ASP:HB2	1:B:413:HIS:HB3	2.00	0.43
1:A:77:PRO:HG3	1:A:143:LEU:HA	1.99	0.43
1:A:162:ASP:HA	1:A:165:ILE:HD11	1.99	0.43
1:A:391:HIS:O	1:A:401:ALA:HA	2.19	0.43
1:A:571:MET:HE2	1:A:579:GLN:O	2.19	0.43
1:B:205:TRP:CZ2	1:B:265:PRO:HD3	2.54	0.43
1:A:142:ARG:HG3	1:A:187:PHE:CE2	2.54	0.43
1:B:350:SER:O	1:B:355:GLY:HA3	2.19	0.43
1:B:611:THR:HG21	1:B:644:VAL:HG21	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:516:ARG:HD3	1:A:535:LEU:HB2	2.01	0.42
1:B:154:LEU:HB2	1:B:184:MET:CE	2.48	0.42
1:B:599:ILE:HD13	1:B:604:PHE:CE1	2.54	0.42
1:B:225:ILE:HG13	1:B:230:LEU:HA	2.01	0.42
1:A:400:MET:HE3	1:A:400:MET:HB2	1.91	0.42
1:B:93:TYR:CD2	1:B:153:TRP:HA	2.54	0.42
1:B:349:ILE:HD12	3:B:936:HOH:O	2.19	0.42
1:B:402:LEU:HD13	1:B:520:LEU:CD1	2.49	0.42
1:A:93:TYR:CD1	1:A:153:TRP:HA	2.54	0.42
1:A:409:GLN:O	1:A:414:GLY:HA3	2.19	0.42
1:A:684:GLU:O	1:A:687:LYS:HB2	2.19	0.42
1:A:63:ASN:HB3	3:A:1063:HOH:O	2.19	0.41
1:A:393:LEU:HD13	1:A:499:GLN:HE21	1.85	0.41
1:B:573:GLY:HA3	1:B:578:TYR:HB3	2.01	0.41
1:B:537:VAL:CG2	1:B:538:SER:N	2.83	0.41
1:B:131:GLU:H	1:B:131:GLU:HG2	1.64	0.41
1:A:611:THR:HA	1:A:651:SER:O	2.19	0.41
1:A:669:THR:HG22	1:A:670:SER:N	2.34	0.41
1:B:612:SER:HG	1:B:681:HIS:HD2	1.64	0.41
1:B:59:ALA:O	1:B:63:ASN:ND2	2.54	0.41
1:B:215:TYR:CE1	1:B:273:ILE:HD13	2.54	0.41
1:A:500:LEU:HD12	1:A:519:LEU:HG	2.03	0.41
1:B:388:LEU:HD12	1:B:405:LYS:HA	2.03	0.41
1:B:421:TRP:CB	1:B:518:MET:HE3	2.38	0.41
1:A:684:GLU:HB3	1:A:716:VAL:HG13	2.02	0.41
1:B:68:MET:CE	1:B:110:LYS:HD3	2.51	0.40
1:A:516:ARG:HG3	1:A:517:SER:N	2.36	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	704/736 (96%)	671 (95%)	31 (4%)	2 (0%)	46 57
1	B	703/736 (96%)	659 (94%)	42 (6%)	2 (0%)	46 57
All	All	1407/1472 (96%)	1330 (94%)	73 (5%)	4 (0%)	46 57

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	720	GLU
1	B	81	GLY
1	B	718	ALA
1	A	69	ALA

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	586/610 (96%)	560 (96%)	26 (4%)	35 49
1	B	585/610 (96%)	546 (93%)	39 (7%)	20 27
All	All	1171/1220 (96%)	1106 (94%)	65 (6%)	26 36

All (65) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	93	TYR
1	A	135	ARG
1	A	137	GLU
1	A	145	TRP
1	A	164	ILE
1	A	165	ILE
1	A	180	VAL
1	A	231	ASP
1	A	233	THR
1	A	456	THR
1	A	497	ASP
1	A	500	LEU

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Mol	Chain	Res	Type
1	A	516	ARG
1	A	518	MET
1	A	522	ASN
1	A	535	LEU
1	A	561	LYS
1	A	603	ARG
1	A	610	VAL
1	A	640	LEU
1	A	656	LEU
1	A	668	THR
1	A	687	LYS
1	A	699	ASN
1	A	711	GLU
1	A	713	VAL
1	B	48	GLU
1	B	78	LYS
1	B	131	GLU
1	B	135	ARG
1	B	137	GLU
1	B	145	TRP
1	B	164	ILE
1	B	180	VAL
1	B	186	LYS
1	B	202	HIS
1	B	212	MET
1	B	225	ILE
1	B	231	ASP
1	B	240	LYS
1	B	260	ARG
1	B	264	MET
1	B	275	THR
1	B	314	LYS
1	B	320	THR
1	B	321	VAL
1	B	338	THR
1	B	360	LYS
1	B	393	LEU
1	B	457	TRP
1	B	462	ILE
1	B	479	VAL
1	B	535	LEU
1	B	538	SER

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Mol	Chain	Res	Type
1	B	574	LYS
1	B	610	VAL
1	B	616	LYS
1	B	617	LYS
1	B	621	LEU
1	B	639	VAL
1	B	656	LEU
1	B	670	SER
1	B	677	LYS
1	B	685	ASN
1	B	694	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	ASN
1	A	132	HIS
1	A	174	GLN
1	A	259	GLN
1	A	344	GLN
1	A	385	GLN
1	A	399	HIS
1	A	428	ASN
1	A	460	GLN
1	A	522	ASN
1	A	585	ASN
1	A	636	GLN
1	A	681	HIS
1	A	714	HIS
1	B	32	ASN
1	B	42	ASN
1	B	86	HIS
1	B	149	ASN
1	B	197	ASN
1	B	259	GLN
1	B	276	ASN
1	B	344	GLN
1	B	409	GLN
1	B	426	ASN
1	B	428	ASN
1	B	486	HIS
1	B	628	ASN

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Mol	Chain	Res	Type
1	B	681	HIS

5.3.3 RNA

There are no RNA molecules 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, 2 are monoatomic - leaving 0 for Mogul analysis.

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.

No monomer is involved in short contacts.

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, 95^{th} 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	706/736 (95%)	-0.30	8 (1%) 82 84	17, 37, 65, 120	0
1	B	705/736 (95%)	-0.16	9 (1%) 79 81	22, 40, 77, 134	0
All	All	1411/1472 (95%)	-0.23	17 (1%) 81 83	17, 39, 72, 134	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	137	GLU	4.9
1	A	720	GLU	4.4
1	B	137	GLU	4.1
1	B	135	ARG	3.4
1	B	136	LYS	3.3
1	B	122	TYR	3.0
1	A	719	GLY	2.9
1	B	130	GLY	2.7
1	B	138	GLN	2.6
1	A	135	ARG	2.5
1	A	178	SER	2.4
1	A	50	TYR	2.4
1	B	82	GLY	2.2
1	B	145	TRP	2.2
1	A	718	ALA	2.2
1	B	735	ARG	2.0
1	A	366	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 (i)

There are no carbohydrates 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. 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, 95^{th} 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	LLDF	B-factors(Å^2)	Q<0.9
2	ZN	A	801	1/1	0.99	0.14	-0.18	27,27,27,27	0
2	ZN	B	801	1/1	0.98	0.13	-0.77	34,34,34,34	0

6.5 Other polymers (i)

There are no such residues in this entry.