Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 08:19 PM GMT

PDB ID	:	1JRQ
Title	:	X-ray Structure Analysis of the Role of the Conserved Tyrosine-369 in Active Site of E. coli Amine Oxidase
Authors	:	Murray, J.M.; Kurtis, C.R.; Tambarajah, W.; Saysell, C.G.; Wilmot, C.M.; Parsons, M.R.; Phillips, S.E.V.; Knowles, P.F.; McPherson, M.J.
Deposited on	:	2001-08-14
Resolution	:	2.15 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

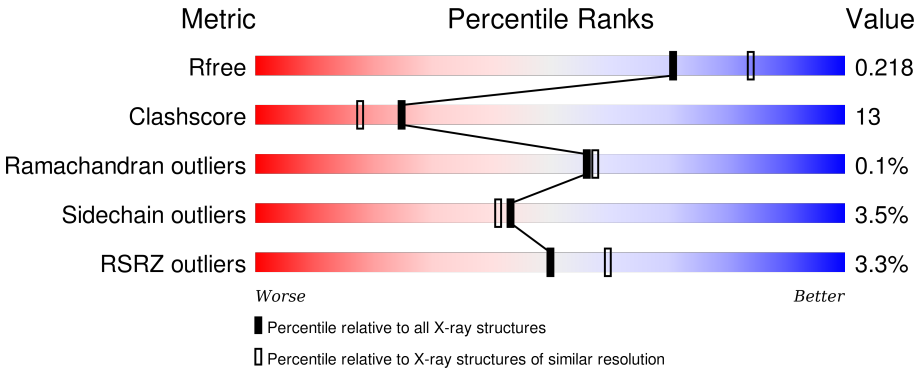
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 2.15 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	1045 (2.16-2.16)
Clashscore	102246	1152 (2.16-2.16)
Ramachandran outliers	100387	1131 (2.16-2.16)
Sidechain outliers	100360	1131 (2.16-2.16)
RSRZ outliers	91569	1050 (2.16-2.16)

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	727	2% 71% 25%
1	B	727	5% 74% 23%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 12750 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 Copper amine oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	718	Total	C	N	O	S	0	0	0
			5663	3602	964	1075	22			
1	B	721	Total	C	N	O	S	0	0	0
			5687	3617	970	1078	22			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	369	PHE	TYR	ENGINEERED	UNP P46883
A	466	TPQ	TYR	MODIFIED RESIDUE	UNP P46883
B	369	PHE	TYR	ENGINEERED	UNP P46883
B	466	TPQ	TYR	MODIFIED RESIDUE	UNP P46883

- Molecule 2 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

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

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total	Ca	0	0
			2	2		
3	A	2	Total	Ca	0	0
			2	2		

- Molecule 4 is water.

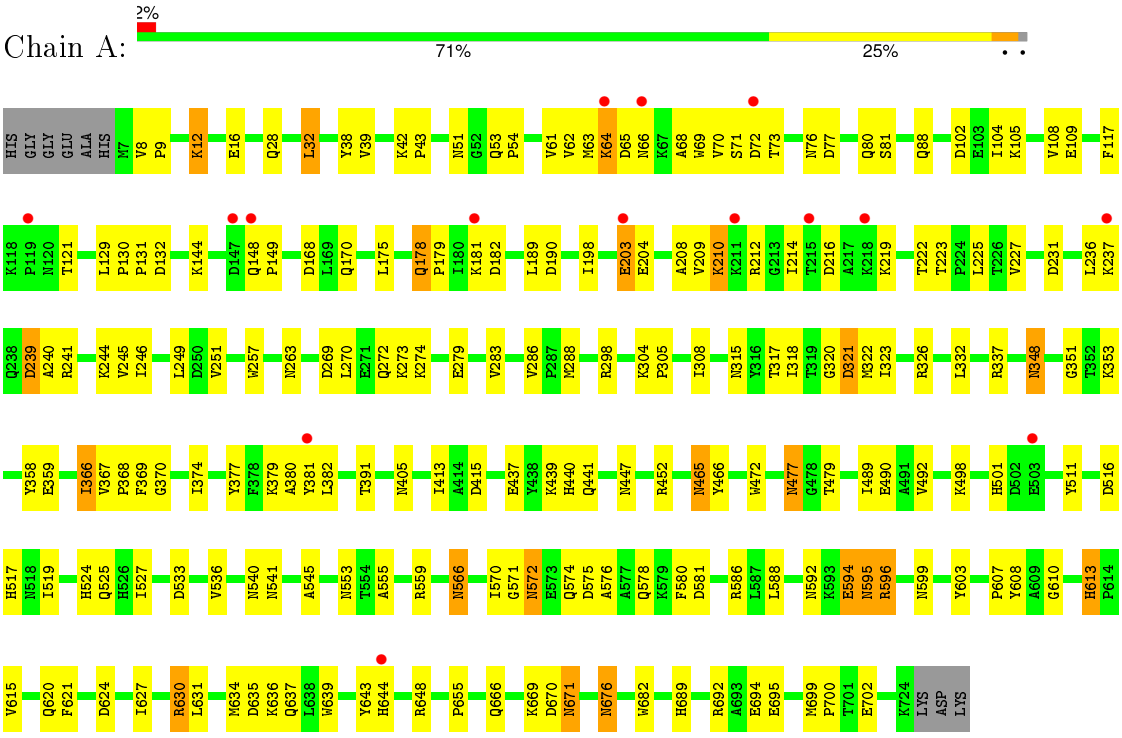
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	749	Total O 749 749	0	0
4	B	645	Total O 645 645	0	0

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.

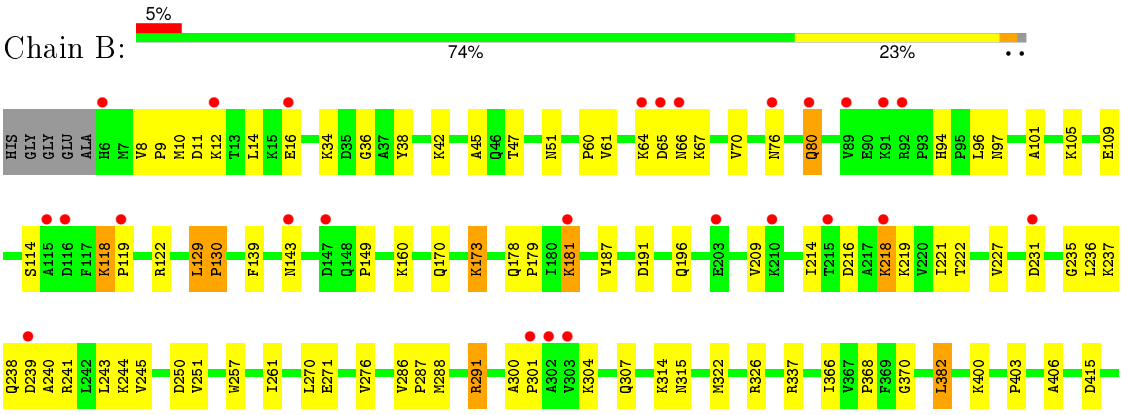
- Molecule 1: Copper amine oxidase

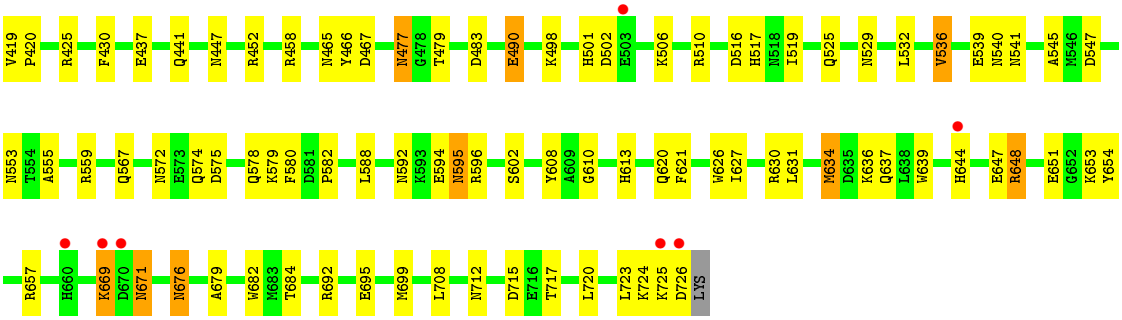
Chain A:



- Molecule 1: Copper amine oxidase

Chain B:




V419 P420 R425 F430 E437 Q441 N447 R452 R458 N465 Y466 D467 N477 G478 T479 D483 E490 K498 H501 D502 E503 K506 R510 D516 H517 N518 I519 Q525 N529 L532 V536 E539 N540 N541 A545 N546 D547
N553 T554 A555 R559 Q567 N572 E573 Q574 D575 Q578 K579 F580 D581 P582 L588 N592 K593 E594 N595 R596 S602 Y608 A609 G610 H613 Q620 F621 W626 I627 R630 L631 M634 D635 K636 Q637 L638 W639 H644 E647 R648 E651 G652 K653 Y654
R657 H660 K669 D670 N671 N676 A679 W682 M683 T684 R692 E695 M699 L708 N712 D715 E716 T717 L720 L723 K724 K725 D726 LYS

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	134.54Å 166.45Å 79.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.15 19.99 – 2.15	Depositor EDS
% Data completeness (in resolution range)	91.4 (20.00-2.15) 92.6 (19.99-2.15)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$< I/\sigma(I) >$ [1]	2.96 (at 2.15Å)	Xtriage
Refinement program	?	Depositor
R, R_{free}	0.195 , 0.235 0.178 , 0.218	Depositor DCC
R_{free} test set	3092 reflections (3.55%)	DCC
Wilson B-factor (Å^2)	24.2	Xtriage
Anisotropy	0.785	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.38 , 62.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Outliers	2 of 90148 reflections (0.002%)	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	12750	wwPDB-VP
Average B, all atoms (Å^2)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.94% 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 (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: CA, TPQ, CU

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	0.34	0/5792	1.24	32/7884 (0.4%)
1	B	0.33	0/5817	1.20	25/7917 (0.3%)
All	All	0.34	0/11609	1.22	57/15801 (0.4%)

There are no bond length outliers.

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	692	ARG	NE-CZ-NH2	-15.20	112.70	120.30
1	A	326	ARG	NE-CZ-NH2	-13.58	113.51	120.30
1	A	692	ARG	NE-CZ-NH2	-11.54	114.53	120.30
1	A	596	ARG	CD-NE-CZ	9.35	136.69	123.60
1	B	291	ARG	NE-CZ-NH2	8.30	124.45	120.30
1	A	630	ARG	NE-CZ-NH1	-7.83	116.39	120.30
1	B	122	ARG	NE-CZ-NH2	-7.81	116.39	120.30
1	B	415	ASP	CB-CG-OD1	7.81	125.33	118.30
1	B	657	ARG	NE-CZ-NH1	-7.65	116.48	120.30
1	A	692	ARG	NE-CZ-NH1	7.63	124.11	120.30
1	B	458	ARG	NE-CZ-NH1	6.76	123.68	120.30
1	A	72	ASP	CB-CG-OD1	6.72	124.35	118.30
1	A	586	ARG	NE-CZ-NH1	6.63	123.62	120.30
1	B	326	ARG	NE-CZ-NH2	-6.59	117.00	120.30
1	A	648	ARG	NE-CZ-NH1	6.41	123.50	120.30
1	A	415	ASP	CB-CG-OD1	6.31	123.98	118.30
1	A	648	ARG	NE-CZ-NH2	-6.31	117.15	120.30
1	B	483	ASP	CB-CG-OD1	6.28	123.95	118.30
1	B	425	ARG	NE-CZ-NH2	6.27	123.44	120.30
1	A	533	ASP	CB-CG-OD2	6.16	123.85	118.30
1	A	326	ARG	NE-CZ-NH1	6.15	123.38	120.30
1	A	588	LEU	N-CA-C	-6.14	94.42	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	689	HIS	CG-ND1-CE1	6.12	116.76	108.20
1	B	458	ARG	NE-CZ-NH2	-6.11	117.25	120.30
1	A	624	ASP	CB-CG-OD2	-6.08	112.83	118.30
1	A	269	ASP	CB-CG-OD1	6.04	123.73	118.30
1	A	102	ASP	CB-CG-OD1	6.01	123.71	118.30
1	A	231	ASP	CB-CG-OD1	5.98	123.68	118.30
1	B	467	ASP	CB-CG-OD2	5.96	123.66	118.30
1	B	490	GLU	OE1-CD-OE2	5.94	130.43	123.30
1	A	129	LEU	N-CA-C	-5.91	95.04	111.00
1	B	326	ARG	CD-NE-CZ	5.91	131.87	123.60
1	B	648	ARG	NE-CZ-NH1	5.90	123.25	120.30
1	A	337	ARG	NE-CZ-NH2	5.87	123.23	120.30
1	B	692	ARG	NH1-CZ-NH2	5.79	125.77	119.40
1	A	635	ASP	CB-CG-OD2	-5.79	113.09	118.30
1	B	588	LEU	N-CA-C	-5.74	95.52	111.00
1	A	581	ASP	CB-CG-OD2	-5.68	113.19	118.30
1	A	440	HIS	N-CA-CB	5.61	120.70	110.60
1	A	298	ARG	NE-CZ-NH2	-5.60	117.50	120.30
1	B	34	LYS	CA-CB-CG	-5.59	101.09	113.40
1	B	502	ASP	CB-CG-OD2	5.52	123.27	118.30
1	B	382	LEU	N-CA-C	-5.52	96.11	111.00
1	B	337	ARG	NE-CZ-NH1	5.48	123.04	120.30
1	A	571	GLY	N-CA-C	5.37	126.51	113.10
1	B	191	ASP	CB-CG-OD2	5.36	123.12	118.30
1	A	298	ARG	NE-CZ-NH1	5.33	122.97	120.30
1	A	472	TRP	N-CA-C	-5.30	96.69	111.00
1	B	261	ILE	N-CA-C	-5.28	96.74	111.00
1	A	635	ASP	CB-CG-OD1	5.28	123.05	118.30
1	B	241	ARG	NE-CZ-NH1	5.26	122.93	120.30
1	A	320	GLY	N-CA-C	-5.15	100.23	113.10
1	A	566	ASN	OD1-CG-ND2	5.14	133.73	121.90
1	B	467	ASP	CB-CG-OD1	-5.11	113.70	118.30
1	B	630	ARG	NE-CZ-NH2	-5.05	117.78	120.30
1	A	321	ASP	CB-CG-OD2	5.02	122.82	118.30
1	A	190	ASP	CB-CG-OD1	5.01	122.81	118.30

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	5663	0	5539	158	0
1	B	5687	0	5562	146	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	749	0	0	16	0
4	B	645	0	0	20	0
All	All	12750	0	11101	288	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 13.

All (288) 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:572:ASN:HD22	1:B:671:ASN:HD21	1.07	0.99
1:A:225:LEU:HD22	1:A:381:TYR:HE2	1.28	0.96
1:B:221:ILE:HD11	1:B:250:ASP:HB2	1.48	0.95
1:A:304:LYS:H	1:B:315:ASN:HD21	1.13	0.91
1:B:322:MET:HG3	4:B:1156:HOH:O	1.73	0.89
1:B:221:ILE:CD1	1:B:250:ASP:HB2	2.04	0.86
1:B:466:TPQ:O5	4:B:1201:HOH:O	1.93	0.85
1:A:249:LEU:HD23	1:A:288:MET:HE1	1.59	0.84
1:A:465:ASN:H	1:A:465:ASN:HD22	1.24	0.82
1:B:94:HIS:HD2	1:B:96:LEU:H	1.27	0.82
1:B:572:ASN:ND2	1:B:671:ASN:HD21	1.77	0.82
1:A:580:PHE:H	1:A:637:GLN:HE21	1.26	0.82
1:B:12:LYS:O	1:B:16:GLU:HG2	1.82	0.80
1:A:62:VAL:HG23	1:A:69:TRP:HB2	1.61	0.79
1:B:181:LYS:HE2	1:B:181:LYS:H	1.50	0.77
1:A:368:PRO:HG3	1:A:634:MET:CE	2.15	0.77
1:A:553:ASN:ND2	1:A:555:ALA:H	1.82	0.76
1:A:368:PRO:HG3	1:A:634:MET:HE3	1.67	0.75
1:B:271:GLU:HG2	4:B:811:HOH:O	1.87	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:490:GLU:HG2	4:B:818:HOH:O	1.87	0.74
1:A:527:ILE:HD12	1:A:634:MET:HE2	1.68	0.74
1:A:279:GLU:OE1	1:A:374:ILE:HD11	1.88	0.73
1:B:580:PHE:H	1:B:637:GLN:HE21	1.34	0.73
1:A:227:VAL:HG12	1:A:244:LYS:HG3	1.70	0.73
1:A:592:ASN:HD21	1:A:676:ASN:HD21	1.37	0.73
1:B:38:TYR:H	1:B:51:ASN:HD21	1.38	0.72
1:B:592:ASN:HD21	1:B:676:ASN:HD21	1.36	0.72
1:B:540:ASN:HB3	1:B:676:ASN:ND2	2.05	0.71
1:A:559:ARG:HH22	1:B:370:GLY:HA2	1.54	0.71
1:A:525:GLN:HE22	1:A:620:GLN:H	1.38	0.71
1:B:466:TPQ:C5	4:B:1201:HOH:O	2.39	0.70
1:A:596:ARG:NH1	4:A:1517:HOH:O	2.17	0.70
1:A:203:GLU:CD	1:A:203:GLU:H	1.95	0.70
1:B:209:VAL:HG13	1:B:214:ILE:HB	1.74	0.69
1:B:227:VAL:HG12	1:B:244:LYS:HG3	1.75	0.69
1:A:225:LEU:HD22	1:A:381:TYR:CE2	2.20	0.69
1:B:38:TYR:H	1:B:51:ASN:ND2	1.91	0.69
1:A:237:LYS:NZ	1:A:239:ASP:HB3	2.08	0.68
1:A:38:TYR:H	1:A:51:ASN:HD21	1.40	0.68
1:A:286:VAL:HG12	1:A:288:MET:HE3	1.75	0.68
1:B:572:ASN:HD22	1:B:671:ASN:ND2	1.85	0.68
1:A:527:ILE:HD12	1:A:634:MET:CE	2.24	0.68
1:A:322:MET:HE1	4:A:1140:HOH:O	1.94	0.67
1:B:648:ARG:NH1	4:B:1044:HOH:O	2.26	0.67
1:A:315:ASN:HD21	1:B:304:LYS:H	1.41	0.67
1:A:38:TYR:H	1:A:51:ASN:ND2	1.91	0.67
1:A:540:ASN:HB3	1:A:676:ASN:ND2	2.11	0.66
1:B:572:ASN:HB2	1:B:671:ASN:ND2	2.11	0.66
1:B:366:ILE:HD11	1:B:627:ILE:HD11	1.78	0.66
1:A:572:ASN:ND2	1:A:575:ASP:H	1.94	0.65
1:A:358:TYR:CD2	1:A:359:GLU:HG3	2.32	0.65
1:A:525:GLN:NE2	1:A:620:GLN:H	1.95	0.65
1:B:553:ASN:ND2	1:B:555:ALA:H	1.94	0.65
1:B:580:PHE:H	1:B:637:GLN:NE2	1.94	0.65
1:A:237:LYS:HZ3	1:A:239:ASP:HB3	1.61	0.65
1:B:237:LYS:HE2	1:B:239:ASP:HB3	1.77	0.65
1:A:225:LEU:CD2	1:A:381:TYR:HE2	2.08	0.64
1:A:272:GLN:HE21	1:A:274:LYS:HD3	1.62	0.64
1:B:216:ASP:OD1	1:B:218:LYS:HB2	1.98	0.64
1:A:131:PRO:HB3	1:A:148:GLN:NE2	2.12	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:465:ASN:H	1:A:465:ASN:ND2	1.95	0.64
1:A:43:PRO:HB3	1:A:63:MET:HG2	1.80	0.64
1:B:525:GLN:HE22	1:B:620:GLN:H	1.44	0.63
1:B:525:GLN:NE2	1:B:620:GLN:H	1.95	0.63
1:B:65:ASP:O	1:B:66:ASN:HB2	1.98	0.63
1:A:216:ASP:HB3	1:A:219:LYS:HD2	1.81	0.63
1:A:517:HIS:CE1	1:B:596:ARG:HH11	2.17	0.62
1:B:181:LYS:CE	1:B:181:LYS:H	2.12	0.62
1:B:61:VAL:HG22	1:B:70:VAL:HG12	1.81	0.62
1:B:8:VAL:HG23	1:B:9:PRO:HD2	1.81	0.62
1:A:382:LEU:HD13	1:A:655:PRO:HB2	1.81	0.62
1:B:498:LYS:O	1:B:517:HIS:HD2	1.83	0.62
1:A:572:ASN:CG	1:A:671:ASN:HD21	2.04	0.61
1:A:28:GLN:HG3	4:A:1170:HOH:O	2.00	0.61
1:B:196:GLN:HE22	1:B:222:THR:H	1.48	0.61
1:A:580:PHE:H	1:A:637:GLN:NE2	1.99	0.60
1:B:291:ARG:NH1	1:B:516:ASP:OD2	2.34	0.60
1:A:149:PRO:HB3	1:A:170:GLN:NE2	2.16	0.60
1:A:594:GLU:OE1	1:B:501:HIS:HE1	1.84	0.59
1:A:368:PRO:HB2	1:A:621:PHE:CZ	2.37	0.59
1:B:8:VAL:CG2	1:B:9:PRO:HD2	2.33	0.59
1:B:466:TPQ:O2	4:B:832:HOH:O	2.17	0.58
1:B:574:GLN:H	1:B:671:ASN:ND2	2.02	0.58
1:B:105:LYS:O	1:B:109:GLU:HG3	2.03	0.58
1:B:251:VAL:HG12	4:B:1373:HOH:O	2.04	0.58
1:B:516:ASP:HB3	1:B:519:ILE:HB	1.86	0.58
1:A:574:GLN:H	1:A:671:ASN:ND2	2.02	0.57
1:B:10:MET:HG2	1:B:14:LEU:HD12	1.87	0.57
1:A:644:HIS:HB3	4:A:1265:HOH:O	2.05	0.57
1:B:216:ASP:HB3	1:B:219:LYS:HD2	1.87	0.56
1:A:636:LYS:HE3	4:A:1465:HOH:O	2.05	0.56
1:A:382:LEU:CD1	1:A:655:PRO:HB2	2.35	0.56
1:B:366:ILE:HD12	1:B:631:LEU:CD1	2.35	0.56
1:A:366:ILE:HG13	1:A:367:VAL:N	2.21	0.56
1:B:67:LYS:HD3	4:B:1429:HOH:O	2.06	0.56
1:A:8:VAL:HG22	1:A:9:PRO:HD2	1.88	0.55
1:B:506:LYS:HE2	1:B:510:ARG:HH12	1.71	0.55
1:B:181:LYS:HE2	1:B:181:LYS:N	2.19	0.55
1:A:368:PRO:HB2	1:A:621:PHE:HZ	1.71	0.55
1:A:288:MET:HA	1:A:288:MET:HE2	1.89	0.55
1:A:608:TYR:HE2	1:B:608:TYR:HE2	1.53	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:288:MET:CE	1:A:288:MET:HA	2.37	0.55
1:B:679:ALA:HB2	4:B:1167:HOH:O	2.07	0.55
1:A:198:ILE:HD11	1:A:273:LYS:HA	1.90	0.54
1:A:73:THR:HG23	1:A:77:ASP:OD2	2.07	0.54
1:A:62:VAL:CG2	1:A:69:TRP:HB2	2.34	0.54
1:A:149:PRO:HB3	1:A:170:GLN:HE21	1.72	0.53
1:B:45:ALA:O	1:B:60:PRO:HB3	2.07	0.53
1:B:717:THR:HB	1:B:720:LEU:HG	1.91	0.53
1:A:237:LYS:HD3	1:A:240:ALA:HB2	1.90	0.53
1:B:243:LEU:CD1	1:B:270:LEU:HD11	2.38	0.53
1:B:307:GLN:HB3	4:B:1093:HOH:O	2.09	0.53
1:A:219:LYS:HD3	4:A:1177:HOH:O	2.08	0.53
1:B:437:GLU:HA	1:B:452:ARG:HB2	1.89	0.52
1:A:366:ILE:HD13	1:A:631:LEU:CD1	2.38	0.52
1:A:466:TPQ:O5	1:A:490:GLU:HA	2.10	0.52
1:A:639:TRP:HB2	1:A:682:TRP:HB2	1.91	0.52
1:B:231:ASP:HB2	1:B:626:TRP:CZ2	2.44	0.52
1:B:368:PRO:HB2	1:B:621:PHE:CZ	2.45	0.52
1:A:615:VAL:CG2	1:B:582:PRO:HB2	2.39	0.52
1:A:615:VAL:HG23	1:B:582:PRO:HB2	1.92	0.52
1:A:214:ILE:HD12	1:A:214:ILE:N	2.25	0.52
1:A:12:LYS:O	1:A:16:GLU:HG3	2.08	0.51
1:B:580:PHE:N	1:B:637:GLN:HE21	2.05	0.51
1:A:498:LYS:O	1:A:517:HIS:HD2	1.93	0.51
1:B:639:TRP:HB2	1:B:682:TRP:HB2	1.92	0.51
1:B:139:PHE:O	1:B:143:ASN:HA	2.10	0.51
1:A:203:GLU:N	1:A:203:GLU:CD	2.62	0.51
1:A:318:ILE:HG12	1:A:323:ILE:HG12	1.92	0.51
1:A:536:VAL:H	1:A:541:ASN:HD21	1.59	0.51
1:B:94:HIS:HB3	1:B:97:ASN:ND2	2.26	0.51
1:A:251:VAL:HG11	1:A:288:MET:HE2	1.91	0.51
1:B:94:HIS:CD2	1:B:96:LEU:H	2.16	0.50
1:A:131:PRO:HA	1:A:148:GLN:HE22	1.75	0.50
1:A:88:GLN:HG2	1:A:317:THR:CG2	2.42	0.50
1:A:203:GLU:HA	4:A:1374:HOH:O	2.12	0.50
1:B:214:ILE:H	1:B:214:ILE:HD12	1.76	0.50
1:B:36:GLY:HA2	1:B:314:LYS:HE2	1.94	0.50
1:A:492:VAL:HB	1:A:519:ILE:HG23	1.94	0.50
1:B:699:MET:CE	4:B:1017:HOH:O	2.60	0.50
1:B:572:ASN:HB2	1:B:671:ASN:HD21	1.74	0.50
1:A:545:ALA:HB2	1:A:570:ILE:HD11	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210:LYS:HE2	1:A:214:ILE:O	2.12	0.50
1:A:477:ASN:HD22	1:A:477:ASN:C	2.15	0.50
1:B:574:GLN:HB2	1:B:671:ASN:CG	2.33	0.49
1:A:572:ASN:C	1:A:572:ASN:HD22	2.16	0.49
1:A:61:VAL:HG22	1:A:70:VAL:HG12	1.94	0.49
1:A:366:ILE:HD11	1:A:380:ALA:HB1	1.94	0.49
1:A:214:ILE:HD12	1:A:214:ILE:H	1.77	0.49
1:A:477:ASN:HD22	1:A:479:THR:H	1.61	0.49
1:A:501:HIS:NE2	1:B:594:GLU:OE1	2.41	0.49
4:A:1238:HOH:O	1:B:400:LYS:HE2	2.12	0.49
1:A:498:LYS:HB3	1:A:498:LYS:NZ	2.27	0.49
1:A:379:LYS:HD3	1:A:381:TYR:OH	2.12	0.49
1:B:441:GLN:OE1	1:B:447:ASN:HB2	2.13	0.49
1:A:348:ASN:HD21	1:A:351:GLY:CA	2.26	0.49
1:A:670:ASP:HB2	4:A:1523:HOH:O	2.13	0.49
1:A:241:ARG:HG2	1:A:270:LEU:HD12	1.94	0.48
1:A:580:PHE:N	1:A:637:GLN:HE21	2.03	0.48
1:A:553:ASN:HD21	1:A:555:ALA:H	1.55	0.48
1:B:76:ASN:O	1:B:80:GLN:HB2	2.13	0.48
1:A:257:TRP:HB2	1:A:489:ILE:HG21	1.96	0.48
1:A:369:PHE:CD2	1:A:524:HIS:HB3	2.49	0.48
1:A:286:VAL:O	1:A:288:MET:HG2	2.14	0.48
1:A:437:GLU:HA	1:A:452:ARG:HB2	1.96	0.48
1:B:595:ASN:HB2	1:B:715:ASP:OD1	2.14	0.48
1:A:272:GLN:NE2	1:A:274:LYS:HD3	2.27	0.48
1:B:406:ALA:HA	1:B:430:PHE:HB3	1.95	0.48
1:B:101:ALA:O	1:B:105:LYS:HG3	2.14	0.47
1:B:726:ASP:O	4:B:1239:HOH:O	2.20	0.47
1:A:64:LYS:HD3	1:A:65:ASP:N	2.29	0.47
1:B:631:LEU:O	1:B:634:MET:HG3	2.14	0.47
1:A:439:LYS:HZ2	1:A:447:ASN:HD21	1.62	0.47
1:A:516:ASP:HB3	1:A:519:ILE:HB	1.96	0.47
1:B:51:ASN:N	1:B:51:ASN:HD22	2.13	0.47
1:B:237:LYS:HD3	1:B:240:ALA:HB2	1.96	0.47
1:A:595:ASN:ND2	1:A:599:ASN:H	2.13	0.47
1:A:368:PRO:HG3	1:A:634:MET:HE1	1.95	0.47
1:A:263:ASN:HD22	1:A:374:ILE:HG22	1.81	0.46
1:A:32:LEU:HB2	1:A:39:VAL:HB	1.96	0.46
1:A:370:GLY:HA2	1:B:559:ARG:HH22	1.81	0.46
1:B:65:ASP:O	1:B:66:ASN:CB	2.64	0.46
1:B:695:GLU:HB3	1:B:699:MET:HG3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:203:GLU:OE2	1:A:204:GLU:HG3	2.16	0.46
1:A:144:LYS:HE3	4:A:1495:HOH:O	2.15	0.46
1:B:129:LEU:HA	1:B:130:PRO:HD2	1.79	0.46
1:A:53:GLN:HA	1:A:54:PRO:HD3	1.72	0.46
1:A:38:TYR:N	1:A:51:ASN:HD21	2.10	0.45
1:B:725:LYS:O	1:B:726:ASP:CB	2.64	0.45
1:B:644:HIS:HB2	1:B:647:GLU:HG3	1.97	0.45
1:A:553:ASN:HD21	1:A:555:ALA:HB3	1.82	0.45
1:A:594:GLU:OE1	1:B:501:HIS:CE1	2.68	0.45
1:B:276:VAL:HG11	4:B:1275:HOH:O	2.17	0.45
1:A:208:ALA:O	1:A:212:ARG:HG3	2.16	0.45
1:B:466:TPQ:O4	4:B:818:HOH:O	1.99	0.45
1:A:576:ALA:O	1:A:578:GLN:HG2	2.17	0.45
1:B:368:PRO:HB2	1:B:621:PHE:HZ	1.82	0.45
1:B:536:VAL:H	1:B:541:ASN:HD21	1.64	0.45
1:A:168:ASP:HB2	1:A:175:LEU:HD21	1.99	0.45
1:B:579:LYS:HA	1:B:637:GLN:NE2	2.30	0.45
1:A:608:TYR:HE2	1:B:608:TYR:CE2	2.32	0.45
1:A:104:ILE:O	1:A:108:VAL:HG23	2.17	0.45
1:B:118:LYS:HA	1:B:119:PRO:HD3	1.70	0.45
1:B:477:ASN:HD22	1:B:477:ASN:C	2.21	0.44
1:A:572:ASN:HD22	1:A:575:ASP:H	1.65	0.44
1:A:694:GLU:O	1:B:717:THR:HA	2.17	0.44
1:B:61:VAL:HG22	1:B:70:VAL:CG1	2.46	0.44
1:B:160:LYS:HD3	1:B:271:GLU:OE1	2.18	0.44
1:A:222:THR:HB	1:A:245:VAL:HG13	1.98	0.44
1:A:441:GLN:OE1	1:A:447:ASN:HB2	2.18	0.44
1:A:613:HIS:CE1	1:A:702:GLU:HG3	2.52	0.44
1:A:308:ILE:O	1:A:405:ASN:HB3	2.18	0.44
1:A:643:TYR:O	1:A:644:HIS:CD2	2.70	0.44
1:B:723:LEU:HD12	1:B:723:LEU:HA	1.86	0.44
1:B:235:GLY:O	1:B:236:LEU:HD23	2.16	0.44
1:B:669:LYS:CE	1:B:669:LYS:HA	2.47	0.44
1:A:236:LEU:HD11	1:A:244:LYS:HE3	2.00	0.43
1:A:210:LYS:HA	1:A:210:LYS:HE2	1.99	0.43
1:B:238:GLN:HB2	4:B:1089:HOH:O	2.17	0.43
1:B:196:GLN:NE2	1:B:222:THR:H	2.15	0.43
1:A:77:ASP:O	1:A:81:SER:HB3	2.17	0.43
1:B:149:PRO:HB3	1:B:170:GLN:HB3	1.99	0.43
1:A:391:THR:HA	1:A:413:ILE:HD11	2.00	0.43
1:A:699:MET:HA	1:A:700:PRO:HD3	1.88	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:610:GLY:HA3	1:B:610:GLY:HA3	2.01	0.43
1:B:300:ALA:HA	1:B:301:PRO:HD2	1.90	0.43
1:B:251:VAL:O	1:B:251:VAL:HG12	2.19	0.43
1:B:572:ASN:OD1	1:B:575:ASP:OD2	2.37	0.43
1:B:257:TRP:CE2	1:B:465:ASN:HB3	2.53	0.43
1:B:720:LEU:HA	1:B:720:LEU:HD23	1.87	0.43
1:A:88:GLN:HG2	1:A:317:THR:HG23	2.01	0.43
1:A:148:GLN:NE2	4:A:1166:HOH:O	2.42	0.42
1:A:178:GLN:HA	1:A:179:PRO:HD3	1.88	0.42
1:B:129:LEU:HD12	1:B:130:PRO:N	2.35	0.42
1:A:181:LYS:O	1:A:182:ASP:HB2	2.19	0.42
1:B:173:LYS:HB3	1:B:173:LYS:NZ	2.34	0.42
1:B:547:ASP:HB3	4:B:1127:HOH:O	2.19	0.42
1:A:63:MET:SD	1:A:68:ALA:HB2	2.59	0.42
1:A:71:SER:OG	1:A:73:THR:HG22	2.20	0.42
1:A:304:LYS:H	1:B:315:ASN:ND2	1.96	0.42
1:B:506:LYS:HD3	1:B:506:LYS:C	2.39	0.42
1:B:699:MET:HE1	4:B:1017:HOH:O	2.19	0.42
1:A:42:LYS:HE2	4:B:822:HOH:O	2.19	0.42
1:B:286:VAL:O	1:B:288:MET:HG2	2.19	0.42
1:A:131:PRO:HA	1:A:148:GLN:NE2	2.34	0.42
1:A:580:PHE:CE2	1:A:607:PRO:HD2	2.55	0.42
1:B:553:ASN:HD21	1:B:555:ALA:H	1.68	0.42
1:A:105:LYS:O	1:A:109:GLU:HG3	2.20	0.42
1:B:653:LYS:HG2	1:B:654:TYR:CE2	2.55	0.42
1:A:353:LYS:NZ	4:A:836:HOH:O	2.48	0.42
1:A:63:MET:HG3	4:A:1492:HOH:O	2.20	0.42
1:B:477:ASN:HD22	1:B:479:THR:H	1.68	0.42
1:B:286:VAL:HA	1:B:287:PRO:HD3	1.90	0.41
1:B:221:ILE:N	1:B:221:ILE:HD12	2.35	0.41
1:B:129:LEU:HD12	1:B:129:LEU:C	2.40	0.41
1:A:511:TYR:HB3	4:A:906:HOH:O	2.20	0.41
1:B:403:PRO:HG3	1:B:430:PHE:CD2	2.55	0.41
1:A:246:ILE:HD13	1:A:246:ILE:HA	1.95	0.41
1:B:237:LYS:HG2	1:B:240:ALA:H	1.84	0.41
1:A:209:VAL:HG13	1:A:214:ILE:HB	2.01	0.41
1:A:130:PRO:HA	1:A:131:PRO:HD3	1.87	0.41
1:A:348:ASN:HD22	1:A:348:ASN:C	2.23	0.41
1:B:382:LEU:HD23	1:B:651:GLU:HG3	2.01	0.41
1:B:578:GLN:HG3	1:B:579:LYS:O	2.21	0.41
1:A:666:GLN:O	1:A:669:LYS:HB2	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:559:ARG:NH2	1:B:370:GLY:HA2	2.28	0.41
1:B:545:ALA:O	1:B:567:GLN:HA	2.21	0.41
1:A:572:ASN:HB2	1:A:671:ASN:ND2	2.35	0.41
1:B:532:LEU:HD13	1:B:708:LEU:HD11	2.03	0.41
1:B:637:GLN:HA	1:B:684:THR:HB	2.03	0.41
1:B:529:ASN:HD21	1:B:682:TRP:HE3	1.69	0.41
1:B:595:ASN:HB3	1:B:712:ASN:O	2.20	0.41
1:A:132:ASP:HB3	4:A:1511:HOH:O	2.21	0.41
1:B:419:VAL:HA	1:B:420:PRO:HD3	1.96	0.41
1:B:222:THR:HB	1:B:245:VAL:CG1	2.51	0.40
1:A:117:PHE:CZ	1:A:121:THR:HB	2.56	0.40
1:B:178:GLN:HA	1:B:179:PRO:HD3	1.79	0.40
1:A:76:ASN:O	1:A:80:GLN:HB2	2.20	0.40
1:A:321:ASP:HB3	1:A:332:LEU:O	2.22	0.40
1:A:490:GLU:OE2	1:A:695:GLU:HB2	2.22	0.40
1:B:536:VAL:HG11	1:B:602:SER:HA	2.03	0.40
1:B:42:LYS:HB3	4:B:1419:HOH:O	2.21	0.40
1:A:305:PRO:HB2	4:A:1391:HOH:O	2.20	0.40
1:B:366:ILE:HD12	1:B:631:LEU:HD13	2.03	0.40
1:B:578:GLN:HA	1:B:636:LYS:HD2	2.03	0.40
1:B:187:VAL:HG22	1:B:243:LEU:HD21	2.03	0.40
1:A:212:ARG:NH1	1:A:283:VAL:HA	2.37	0.40
1:A:630:ARG:HD3	1:A:630:ARG:HH11	1.70	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	715/727 (98%)	688 (96%)	27 (4%)	0	100 100
1	B	718/727 (99%)	693 (96%)	23 (3%)	2 (0%)	46 42

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1433/1454 (99%)	1381 (96%)	50 (4%)	2 (0%)	56 57

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	536	VAL
1	B	130	PRO

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	609/615 (99%)	585 (96%)	24 (4%)	39 36
1	B	611/615 (99%)	592 (97%)	19 (3%)	47 47
All	All	1220/1230 (99%)	1177 (96%)	43 (4%)	43 41

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	LYS
1	A	32	LEU
1	A	64	LYS
1	A	66	ASN
1	A	178	GLN
1	A	189	LEU
1	A	203	GLU
1	A	210	LYS
1	A	223	THR
1	A	239	ASP
1	A	348	ASN
1	A	366	ILE
1	A	377	TYR
1	A	465	ASN
1	A	477	ASN
1	A	566	ASN

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Mol	Chain	Res	Type
1	A	572	ASN
1	A	594	GLU
1	A	595	ASN
1	A	603	TYR
1	A	613	HIS
1	A	627	ILE
1	A	671	ASN
1	A	676	ASN
1	B	11	ASP
1	B	47	THR
1	B	64	LYS
1	B	80	GLN
1	B	114	SER
1	B	118	LYS
1	B	129	LEU
1	B	173	LYS
1	B	181	LYS
1	B	218	LYS
1	B	477	ASN
1	B	539	GLU
1	B	595	ASN
1	B	613	HIS
1	B	634	MET
1	B	669	LYS
1	B	671	ASN
1	B	676	ASN
1	B	724	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (60) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	51	ASN
1	A	97	ASN
1	A	148	GLN
1	A	170	GLN
1	A	196	GLN
1	A	200	ASN
1	A	201	ASN
1	A	263	ASN
1	A	272	GLN
1	A	307	GLN
1	A	315	ASN

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Mol	Chain	Res	Type
1	A	324	HIS
1	A	327	ASN
1	A	348	ASN
1	A	447	ASN
1	A	465	ASN
1	A	477	ASN
1	A	517	HIS
1	A	525	GLN
1	A	529	ASN
1	A	541	ASN
1	A	553	ASN
1	A	567	GLN
1	A	572	ASN
1	A	595	ASN
1	A	599	ASN
1	A	604	GLN
1	A	637	GLN
1	A	660	HIS
1	A	671	ASN
1	A	676	ASN
1	B	51	ASN
1	B	94	HIS
1	B	97	ASN
1	B	143	ASN
1	B	196	GLN
1	B	197	ASN
1	B	200	ASN
1	B	238	GLN
1	B	263	ASN
1	B	315	ASN
1	B	327	ASN
1	B	350	ASN
1	B	447	ASN
1	B	477	ASN
1	B	501	HIS
1	B	517	HIS
1	B	525	GLN
1	B	529	ASN
1	B	541	ASN
1	B	553	ASN
1	B	567	GLN
1	B	572	ASN

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Mol	Chain	Res	Type
1	B	595	ASN
1	B	599	ASN
1	B	613	HIS
1	B	637	GLN
1	B	644	HIS
1	B	671	ASN
1	B	676	ASN

5.3.3 RNA

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains

2 non-standard protein/DNA/RNA residues 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	$\#\|Z\| > 2$
1	TPQ	A	466	1	13,14,15	2.26	4 (30%)	15,19,21	2.26	4 (26%)
1	TPQ	B	466	1	13,14,15	2.42	5 (38%)	15,19,21	2.57	7 (46%)

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
1	TPQ	A	466	1	-	0/4/22/24	0/1/1/1
1	TPQ	B	466	1	-	0/4/22/24	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	466	TPQ	C1-C2	-5.75	1.41	1.49
1	A	466	TPQ	C1-C2	-5.65	1.41	1.49
1	B	466	TPQ	C6-C5	-3.91	1.33	1.44
1	A	466	TPQ	C6-C5	-2.72	1.36	1.44
1	A	466	TPQ	C3-C4	2.06	1.39	1.35
1	B	466	TPQ	C3-C4	2.19	1.39	1.35
1	B	466	TPQ	O5-C5	2.36	1.31	1.24
1	B	466	TPQ	CB-C1	2.93	1.57	1.50
1	A	466	TPQ	CB-C1	3.24	1.57	1.50

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	466	TPQ	CA-CB-C1	-4.92	103.76	113.63
1	B	466	TPQ	CA-CB-C1	-4.91	103.79	113.63
1	B	466	TPQ	C3-C4-C5	-2.88	118.17	121.18
1	B	466	TPQ	C1-C6-C5	-2.70	121.33	122.97
1	B	466	TPQ	O-C-CA	-2.17	119.84	125.49
1	A	466	TPQ	O2-C2-C3	-2.06	117.26	121.89
1	A	466	TPQ	CB-C1-C2	2.25	121.85	118.33
1	B	466	TPQ	C6-C5-C4	3.02	122.51	117.34
1	B	466	TPQ	CB-C1-C2	3.02	123.04	118.33
1	B	466	TPQ	C3-C2-C1	4.66	121.79	118.30
1	A	466	TPQ	C3-C2-C1	4.79	121.89	118.30

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	466	TPQ	1	0
1	B	466	TPQ	4	0

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 6 ligands modelled in this entry, 6 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	717/727 (98%)	-0.31	15 (2%)	67	74	15, 26, 46, 62	0
1	B	720/727 (99%)	-0.18	33 (4%)	36	47	17, 29, 50, 65	0
All	All	1437/1454 (98%)	-0.24	48 (3%)	50	60	15, 28, 48, 65	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	6	HIS	7.7
1	B	726	ASP	5.6
1	B	115	ALA	5.4
1	B	91	LYS	4.5
1	B	302	ALA	4.5
1	A	64	LYS	4.4
1	B	65	ASP	4.2
1	B	215	THR	4.1
1	B	301	PRO	4.0
1	A	119	PRO	3.7
1	A	148	GLN	3.7
1	A	147	ASP	3.4
1	B	64	LYS	3.3
1	A	218	LYS	3.2
1	B	16	GLU	3.2
1	B	239	ASP	3.0
1	B	644	HIS	3.0
1	B	116	ASP	2.9
1	A	381	TYR	2.9
1	B	80	GLN	2.6
1	A	237	LYS	2.6
1	B	503	GLU	2.5
1	B	218	LYS	2.5
1	B	66	ASN	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	92	ARG	2.5
1	A	215	THR	2.4
1	B	143	ASN	2.4
1	B	89	VAL	2.3
1	A	181	LYS	2.3
1	A	211	LYS	2.3
1	B	76	ASN	2.3
1	A	644	HIS	2.3
1	A	503	GLU	2.3
1	B	12	LYS	2.2
1	B	725	LYS	2.2
1	B	670	ASP	2.2
1	B	669	LYS	2.2
1	B	147	ASP	2.2
1	B	303	VAL	2.2
1	B	231	ASP	2.1
1	B	181	LYS	2.1
1	B	210	LYS	2.1
1	B	660	HIS	2.1
1	B	119	PRO	2.1
1	A	203	GLU	2.0
1	A	72	ASP	2.0
1	B	203	GLU	2.0
1	A	66	ASN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (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
1	TPQ	A	466	14/15	0.78	0.21	-	20,43,47,48	0
1	TPQ	B	466	14/15	0.87	0.18	-	22,45,49,50	0

6.3 Carbohydrates (i)

There are no carbohydrates in this entry.

6.4 Ligands

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 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
3	CA	A	802	1/1	0.99	0.12	0.18	26,26,26,26	0
3	CA	A	803	1/1	0.94	0.12	-0.04	54,54,54,54	0
3	CA	B	802	1/1	0.99	0.13	-0.07	29,29,29,29	0
3	CA	B	803	1/1	0.92	0.09	-0.92	54,54,54,54	0
2	CU	B	801	1/1	1.00	0.05	-	28,28,28,28	0
2	CU	A	801	1/1	0.99	0.07	-	28,28,28,28	0

6.5 Other polymers

There are no such residues in this entry.