



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:07 AM GMT

PDB ID : 2W1B  
Title : THE STRUCTURE OF THE EFFLUX PUMP ACRB IN COMPLEX WITH BILE ACID  
Authors : Drew, D.; Klepsch, M.M.; Newstead, S.; Flaig, R.; De Gier, J.W.; Iwata, S.; Beis, K.  
Deposited on : 2008-10-17  
Resolution : 3.85 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](mailto: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.

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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

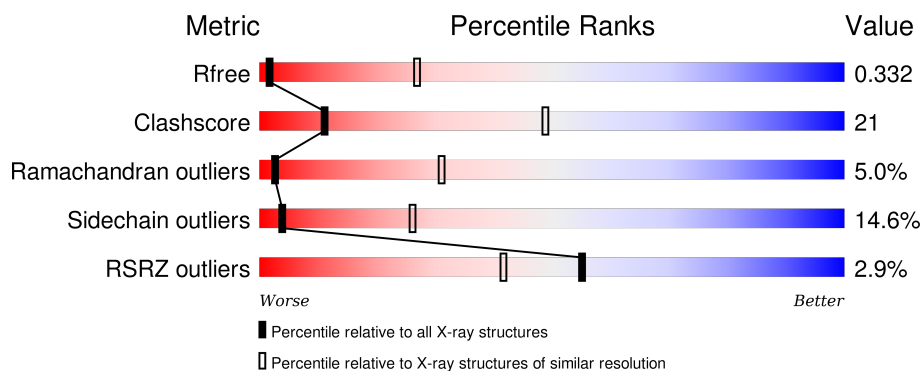
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*


The reported resolution of this entry is 3.85 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	1000 (4.20-3.52)
Clashscore	102246	1090 (4.20-3.52)
Ramachandran outliers	100387	1046 (4.20-3.52)
Sidechain outliers	100360	1038 (4.20-3.52)
RSRZ outliers	91569	1004 (4.20-3.52)

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  $\geq 3$ , 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  $\leq 5\%$ . 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	1049	

## 2 Entry composition [i](#)

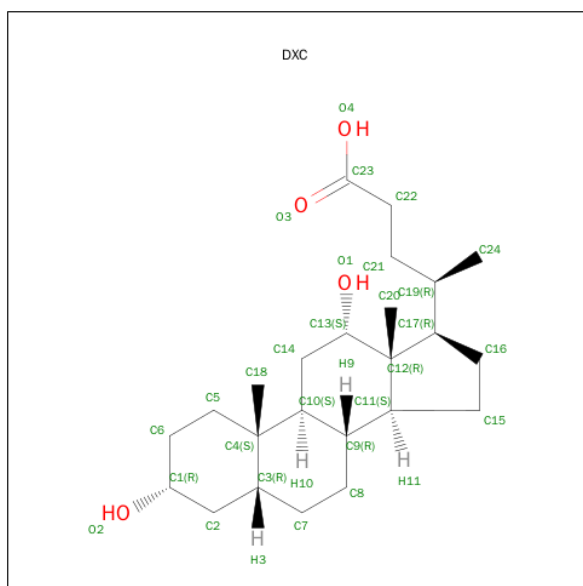
There are 2 unique types of molecules in this entry. The entry contains 7869 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 ACRIFLAVIN RESISTANCE PROTEIN B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1032	Total	C	N	O	S	0	0	0
			7841	5047	1294	1457	43			

- Molecule 2 is (3ALPHA,5BETA,12ALPHA)-3,12-DIHYDROXYCHOLAN-24-OIC ACID (three-letter code: DXC) (formula: C<sub>24</sub>H<sub>40</sub>O<sub>4</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			28	24	4		



L886	C887	L888	S894	H895	S896	I897	P898	F899	S900	V901	V904	V905	P906	L907	G908	V909	I910	L914	A915	G920	L921	D924	V925	Y926	V929	T934	I935	L944	E947	K950	E956	A963	T964	L965	V968	R971	L972	R973	P974	I975	L980	M987	P988		
L989	V990	I991	S992	T993	S997	G998	A999	Q1000	N1001	T1005	M1008	T1013	A1014	T1015	V1016	L1017	F1020	F1021	V1022	P1023	V1024	F1025	V1028	V1029	R1030	R1031	R1032	F1033	SER	ARG	LYS	ASN	GLU	ASP	ILE	GLU	HIS	SER	HIS	HIS	THR	VAL	ASP	HIS	HIS

## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	145.49Å 145.49Å 515.18Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	40.00 – 3.85 57.34 – 3.85	Depositor EDS
% Data completeness (in resolution range)	95.7 (40.00-3.85) 95.2 (57.34-3.85)	Depositor EDS
$R_{merge}$	0.23	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.45 (at 3.88Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.278 , 0.349 0.268 , 0.332	Depositor DCC
$R_{free}$ test set	977 reflections (5.30%)	DCC
Wilson B-factor (Å <sup>2</sup> )	96.0	Xtriage
Anisotropy	0.330	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 79.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.41$ , $\langle L^2 \rangle = 0.23$	Xtriage
Outliers	0 of 19432 reflections	Xtriage
$F_o, F_c$ correlation	0.84	EDS
Total number of atoms	7869	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	75.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.18% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  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: DXC

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	0.52	2/7991 (0.0%)	0.65	2/10852 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	386	PHE	CG-CD2	5.22	1.46	1.38
1	A	653	ARG	CB-CG	5.19	1.66	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	713	LEU	CA-CB-CG	5.32	127.54	115.30
1	A	972	LEU	CA-CB-CG	5.15	127.14	115.30

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	7841	0	7990	329	0
2	A	28	0	39	1	0
All	All	7869	0	8029	330	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 21.

All (330) 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:A:2034:DXC:C15	2:A:2034:DXC:C16	1.84	1.41
1:A:971:ARG:HG2	1:A:974:PRO:HG2	1.30	1.08
1:A:971:ARG:HB2	1:A:971:ARG:NH1	1.74	1.02
1:A:552:MET:HA	1:A:910:ILE:HD13	1.43	0.97
1:A:33:ALA:HA	1:A:300:LEU:HD13	1.44	0.95
1:A:904:VAL:HA	1:A:907:LEU:HD13	1.50	0.94
1:A:367:ILE:HB	1:A:368:PRO:HD3	1.52	0.91
1:A:897:ILE:H	1:A:898:PRO:HD2	1.38	0.88
1:A:676:THR:HA	1:A:862:MET:HB2	1.55	0.88
1:A:239:ARG:HH11	1:A:239:ARG:HG3	1.40	0.86
1:A:453:PHE:O	1:A:471:SER:OG	1.96	0.84
1:A:306:ILE:HG22	1:A:306:ILE:O	1.76	0.84
1:A:971:ARG:HB2	1:A:971:ARG:CZ	2.07	0.83
1:A:372:VAL:HB	1:A:373:PRO:HD3	1.63	0.81
1:A:393:LEU:HD23	1:A:470:PHE:HE1	1.45	0.81
1:A:684:LEU:O	1:A:824:SER:HA	1.80	0.81
1:A:393:LEU:HD23	1:A:470:PHE:CE1	2.15	0.81
1:A:451:ALA:HB1	1:A:883:VAL:HG13	1.64	0.80
1:A:330:THR:H	1:A:331:PRO:HD2	1.46	0.80
1:A:448:VAL:HG22	1:A:887:CYS:HB3	1.63	0.79
1:A:578:LEU:HB2	1:A:623:ASN:HB2	1.65	0.78
1:A:465:ALA:O	1:A:469:GLN:HG2	1.84	0.77
1:A:944:LEU:HB3	1:A:971:ARG:HH21	1.48	0.77
1:A:415:ASN:O	1:A:419:VAL:HG23	1.85	0.77
1:A:225:VAL:O	1:A:226:LYS:O	2.03	0.77
1:A:19:ILE:HG22	1:A:378:GLY:HA2	1.68	0.75
1:A:64:VAL:HG11	1:A:117:LEU:HB3	1.70	0.74
1:A:342:LYS:HG2	1:A:346:GLU:OE2	1.89	0.73
1:A:576:VAL:HA	1:A:663:VAL:HG22	1.71	0.73
1:A:626:ILE:HG12	1:A:628:PHE:CE1	2.24	0.73
1:A:743:ILE:HD12	1:A:743:ILE:H	1.53	0.72
1:A:144:ASN:HD22	1:A:149:MET:HG2	1.53	0.72
1:A:352:PHE:O	1:A:352:PHE:HD2	1.72	0.71
1:A:758:TYR:HE1	1:A:770:LYS:HB3	1.53	0.71
1:A:907:LEU:HD23	1:A:1017:LEU:HB3	1.73	0.71
1:A:402:ILE:O	1:A:406:VAL:HG23	1.91	0.69
1:A:915:ALA:HB1	1:A:1005:THR:HG22	1.73	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:574:THR:HG23	1:A:627:ALA:HB3	1.74	0.69
1:A:35:TYR:CE1	1:A:671:ILE:HG12	2.27	0.68
1:A:36:PRO:HG2	1:A:38:ILE:HG12	1.74	0.68
1:A:70:ASN:O	1:A:72:ILE:HG13	1.93	0.67
1:A:564:LEU:HG	1:A:926:TYR:CE1	2.30	0.67
1:A:411:VAL:O	1:A:415:ASN:HB2	1.95	0.67
1:A:356:TYR:O	1:A:360:GLN:N	2.24	0.67
1:A:989:LEU:O	1:A:1001:ASN:ND2	2.29	0.66
1:A:344:LEU:HD21	1:A:399:VAL:HA	1.77	0.66
1:A:186:ILE:HD13	1:A:262:LEU:HD21	1.77	0.66
1:A:682:PHE:HB3	1:A:827:ILE:HB	1.76	0.66
1:A:817:GLU:O	1:A:818:ARG:HG3	1.95	0.65
1:A:330:THR:N	1:A:331:PRO:HD2	2.12	0.65
1:A:897:ILE:N	1:A:898:PRO:HD2	2.12	0.65
1:A:697:GLN:O	1:A:700:ASN:HB2	1.95	0.64
1:A:58:GLN:HA	1:A:62:THR:OG1	1.98	0.64
1:A:1020:PHE:O	1:A:1023:PRO:HD2	1.98	0.64
1:A:449:LEU:O	1:A:453:PHE:HD1	1.81	0.63
1:A:195:LYS:HB3	1:A:196:PHE:CD1	2.34	0.63
1:A:149:MET:HG3	1:A:154:ILE:HG13	1.81	0.63
1:A:408:ASP:O	1:A:412:VAL:HG23	1.99	0.62
1:A:758:TYR:CE1	1:A:770:LYS:HB3	2.34	0.62
1:A:200:PRO:HD2	1:A:749:THR:HG22	1.81	0.62
1:A:368:PRO:HG3	1:A:413:VAL:HG21	1.80	0.62
1:A:68:ASN:O	1:A:110:LYS:HD2	2.00	0.62
1:A:239:ARG:HG3	1:A:239:ARG:NH1	2.14	0.61
1:A:758:TYR:OH	1:A:761:ASP:OD1	2.13	0.61
1:A:545:TYR:HA	1:A:548:ILE:HD12	1.82	0.61
1:A:230:LEU:HG	1:A:231:ASN:N	2.15	0.61
1:A:41:PRO:HB2	1:A:94:PHE:HB2	1.82	0.61
1:A:990:VAL:HG21	1:A:1008:MET:HE3	1.83	0.61
1:A:888:LEU:HB2	1:A:898:PRO:HB3	1.83	0.61
1:A:23:GLY:HA3	1:A:377:LEU:O	2.01	0.60
1:A:971:ARG:HB2	1:A:971:ARG:HH11	1.66	0.60
1:A:733:GLN:HE22	1:A:743:ILE:HG21	1.67	0.60
1:A:95:GLU:O	1:A:98:THR:HB	2.02	0.60
1:A:200:PRO:HD2	1:A:749:THR:CG2	2.33	0.59
1:A:5:PHE:HE2	1:A:11:PHE:HD1	1.51	0.59
1:A:73:ASP:HB2	1:A:106:GLN:HE22	1.66	0.59
1:A:190:PRO:HG3	1:A:789:TRP:CH2	2.38	0.59
1:A:35:TYR:HB3	1:A:36:PRO:CD	2.33	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:741:VAL:HG21	1:A:799:VAL:HG21	1.84	0.59
1:A:971:ARG:CB	1:A:971:ARG:CZ	2.81	0.59
1:A:344:LEU:CD2	1:A:402:ILE:HD11	2.33	0.59
1:A:5:PHE:CE1	1:A:487:ILE:HG12	2.38	0.58
1:A:679:GLY:HA3	1:A:830:GLN:HG2	1.85	0.58
1:A:527:TYR:O	1:A:531:VAL:HG23	2.04	0.58
1:A:411:VAL:HG22	1:A:974:PRO:HB3	1.85	0.58
1:A:997:SER:HA	1:A:1000:GLN:HB2	1.85	0.58
1:A:874:PRO:HA	1:A:877:TYR:HB2	1.85	0.58
1:A:58:GLN:HE22	1:A:63:GLN:HE21	1.52	0.58
1:A:753:ALA:O	1:A:775:SER:HB3	2.04	0.58
1:A:552:MET:CA	1:A:910:ILE:HD13	2.27	0.57
1:A:35:TYR:HB3	1:A:36:PRO:HD2	1.86	0.57
1:A:337:ILE:HG22	1:A:338:HIS:H	1.69	0.57
1:A:762:PHE:CE1	1:A:764:ASP:HB2	2.41	0.56
1:A:904:VAL:HA	1:A:907:LEU:CD1	2.29	0.56
1:A:527:TYR:OH	1:A:968:VAL:HG22	2.05	0.56
1:A:372:VAL:HB	1:A:373:PRO:CD	2.33	0.56
1:A:470:PHE:CD2	1:A:929:VAL:HG11	2.41	0.56
1:A:750:LEU:HD12	1:A:754:TRP:CD1	2.41	0.56
1:A:32:VAL:O	1:A:300:LEU:HB2	2.05	0.56
1:A:228:GLN:HG2	1:A:230:LEU:H	1.71	0.56
1:A:337:ILE:CG2	1:A:338:HIS:N	2.69	0.55
1:A:449:LEU:O	1:A:453:PHE:CD1	2.58	0.55
1:A:882:ILE:O	1:A:886:LEU:HG	2.06	0.55
1:A:395:MET:O	1:A:398:MET:HB2	2.07	0.55
1:A:714:THR:HG22	1:A:715:SER:H	1.72	0.55
1:A:393:LEU:HD13	1:A:469:GLN:HG3	1.88	0.55
1:A:513:PHE:O	1:A:516:PHE:HB2	2.06	0.55
1:A:144:ASN:ND2	1:A:149:MET:H	2.04	0.55
1:A:174:ASP:O	1:A:292:LYS:HB2	2.07	0.55
1:A:56:THR:HG22	1:A:56:THR:O	2.06	0.54
1:A:435:MET:C	1:A:437:GLN:H	2.11	0.54
1:A:5:PHE:HE2	1:A:11:PHE:CD1	2.26	0.54
1:A:414:GLU:HA	1:A:417:GLU:HG2	1.90	0.54
1:A:827:ILE:O	1:A:828:LEU:HD22	2.08	0.54
1:A:351:VAL:O	1:A:353:LEU:N	2.41	0.54
1:A:406:VAL:HG13	1:A:410:ILE:HD12	1.91	0.53
1:A:563:PHE:O	1:A:924:ASP:HB2	2.09	0.53
1:A:394:THR:O	1:A:398:MET:HG2	2.09	0.53
1:A:109:ASN:O	1:A:113:LEU:HG	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:53:ASP:O	1:A:57:VAL:HG23	2.09	0.53
1:A:35:TYR:CB	1:A:36:PRO:CD	2.86	0.53
1:A:598:TYR:HB3	1:A:606:VAL:HG11	1.89	0.53
1:A:989:LEU:HB3	1:A:1000:GLN:O	2.09	0.53
1:A:35:TYR:CB	1:A:36:PRO:HD2	2.39	0.53
1:A:408:ASP:N	1:A:408:ASP:OD1	2.41	0.53
1:A:671:ILE:HD12	1:A:674:LEU:HD22	1.91	0.53
1:A:69:MET:O	1:A:70:ASN:HB3	2.09	0.53
1:A:210:GLN:OE1	1:A:249:ILE:HA	2.08	0.52
1:A:415:ASN:OD1	1:A:418:ARG:NH2	2.40	0.52
1:A:5:PHE:CE2	1:A:11:PHE:HD1	2.27	0.52
1:A:456:MET:O	1:A:467:TYR:HB3	2.09	0.52
1:A:144:ASN:ND2	1:A:149:MET:HG2	2.22	0.52
1:A:3:ASN:HA	1:A:6:ILE:HG12	1.91	0.52
1:A:699:ARG:HE	1:A:718:PRO:HB3	1.75	0.52
1:A:356:TYR:O	1:A:357:LEU:C	2.48	0.52
1:A:262:LEU:HG	1:A:268:ILE:HD11	1.90	0.52
1:A:57:VAL:HG11	1:A:86:GLY:O	2.10	0.52
1:A:872:GLN:HG2	1:A:872:GLN:O	2.10	0.52
1:A:57:VAL:HG21	1:A:86:GLY:HA2	1.92	0.52
1:A:79:SER:OG	1:A:80:SER:N	2.43	0.52
1:A:546:LEU:O	1:A:550:VAL:HG23	2.10	0.52
1:A:534:ILE:HG22	1:A:541:TYR:CZ	2.45	0.51
1:A:1013:THR:C	1:A:1015:THR:H	2.14	0.51
1:A:34:GLN:HA	1:A:333:VAL:HG11	1.92	0.51
1:A:563:PHE:CE2	1:A:674:LEU:HD21	2.44	0.51
1:A:684:LEU:O	1:A:824:SER:CA	2.57	0.51
1:A:489:THR:HB	1:A:490:PRO:CD	2.40	0.51
1:A:582:ALA:HB1	1:A:586:ARG:NE	2.26	0.51
1:A:367:ILE:HB	1:A:368:PRO:CD	2.35	0.51
1:A:563:PHE:HE2	1:A:674:LEU:HD21	1.75	0.51
1:A:195:LYS:HB3	1:A:196:PHE:HD1	1.73	0.51
1:A:184:MET:HB2	1:A:762:PHE:CE2	2.46	0.51
1:A:351:VAL:C	1:A:353:LEU:H	2.14	0.51
1:A:88:VAL:HG13	1:A:88:VAL:O	2.10	0.51
1:A:897:ILE:H	1:A:898:PRO:CD	2.18	0.51
1:A:1017:LEU:O	1:A:1021:PHE:HD1	1.93	0.50
1:A:699:ARG:O	1:A:703:LEU:HG	2.11	0.50
1:A:187:TRP:HA	1:A:187:TRP:CE3	2.46	0.50
1:A:896:SER:O	1:A:897:ILE:HG13	2.11	0.50
1:A:187:TRP:HZ3	1:A:774:MET:HB3	1.76	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:310:LEU:O	1:A:314:GLU:HG3	2.12	0.50
1:A:183:ALA:HB2	1:A:273:GLU:CG	2.42	0.50
1:A:470:PHE:O	1:A:472:ILE:N	2.44	0.50
1:A:817:GLU:C	1:A:818:ARG:HG3	2.32	0.50
1:A:485:ALA:HA	1:A:489:THR:HB	1.93	0.50
1:A:971:ARG:CG	1:A:974:PRO:HG2	2.22	0.50
1:A:68:ASN:O	1:A:110:LYS:CD	2.60	0.50
1:A:595:THR:HG22	1:A:599:LEU:HD12	1.94	0.50
1:A:337:ILE:CG2	1:A:338:HIS:H	2.25	0.49
1:A:905:VAL:HG22	1:A:935:ILE:HG23	1.92	0.49
1:A:545:TYR:O	1:A:548:ILE:HB	2.13	0.49
1:A:352:PHE:O	1:A:352:PHE:CD2	2.59	0.49
1:A:351:VAL:C	1:A:353:LEU:N	2.66	0.49
1:A:306:ILE:CG2	1:A:306:ILE:O	2.47	0.49
1:A:187:TRP:HA	1:A:187:TRP:HE3	1.77	0.48
1:A:947:GLU:O	1:A:950:LYS:HB3	2.12	0.48
1:A:450:SER:HA	1:A:453:PHE:HB2	1.94	0.48
1:A:61:VAL:HG13	1:A:118:LEU:HD22	1.95	0.48
1:A:999:ALA:O	1:A:1001:ASN:N	2.46	0.48
1:A:188:MET:HB2	1:A:775:SER:HA	1.95	0.48
1:A:564:LEU:HG	1:A:926:TYR:HE1	1.78	0.48
1:A:1030:ARG:C	1:A:1032:ARG:H	2.16	0.48
1:A:293:LEU:HD11	1:A:299:ALA:HB2	1.96	0.48
1:A:213:GLN:HG2	1:A:239:ARG:HG2	1.94	0.48
1:A:314:GLU:HA	1:A:317:PHE:CE2	2.49	0.48
1:A:344:LEU:HD21	1:A:402:ILE:HD11	1.96	0.47
1:A:791:VAL:O	1:A:798:MET:HA	2.14	0.47
1:A:404:LEU:HD23	1:A:478:MET:HG3	1.97	0.47
1:A:166:ILE:HG22	1:A:175:VAL:HG21	1.95	0.47
1:A:591:LEU:HD13	1:A:611:ALA:HB1	1.94	0.47
1:A:576:VAL:HB	1:A:625:GLY:HA3	1.95	0.47
1:A:817:GLU:O	1:A:818:ARG:CG	2.62	0.47
1:A:421:ALA:O	1:A:503:GLY:HA2	2.14	0.47
1:A:218:GLN:HG2	1:A:233:SER:HA	1.95	0.47
1:A:719:ASN:HB2	1:A:828:LEU:HD23	1.96	0.47
1:A:534:ILE:HG22	1:A:541:TYR:CE1	2.49	0.47
1:A:383:LEU:O	1:A:384:ALA:C	2.52	0.47
1:A:690:LEU:HD11	1:A:854:GLY:HA3	1.96	0.47
1:A:470:PHE:C	1:A:472:ILE:H	2.18	0.47
1:A:448:VAL:HG22	1:A:887:CYS:CB	2.38	0.47
1:A:346:GLU:O	1:A:350:LEU:HD12	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:699:ARG:NE	1:A:718:PRO:HB3	2.30	0.47
1:A:192:GLU:O	1:A:265:VAL:HG12	2.15	0.47
1:A:682:PHE:CZ	1:A:857:TYR:HB2	2.49	0.47
1:A:351:VAL:O	1:A:355:MET:HB2	2.15	0.47
1:A:677:ALA:HA	1:A:867:ARG:HH21	1.80	0.47
1:A:888:LEU:CB	1:A:898:PRO:HB3	2.45	0.47
1:A:832:ALA:HA	1:A:833:PRO:HD3	1.79	0.46
1:A:158:VAL:HG13	1:A:162:MET:HE2	1.97	0.46
1:A:578:LEU:O	1:A:580:ALA:N	2.49	0.46
1:A:611:ALA:HA	1:A:627:ALA:HA	1.96	0.46
1:A:98:THR:HG23	1:A:99:ASP:N	2.31	0.46
1:A:49:TYR:HA	1:A:50:PRO:HD2	1.64	0.46
1:A:196:PHE:N	1:A:196:PHE:CD1	2.84	0.46
1:A:17:ILE:O	1:A:20:MET:HB2	2.16	0.46
1:A:33:ALA:HA	1:A:300:LEU:CD1	2.32	0.46
1:A:30:LEU:HD12	1:A:31:PRO:HD2	1.97	0.46
1:A:467:TYR:HE1	1:A:868:LEU:HD21	1.81	0.46
1:A:521:GLU:O	1:A:524:THR:HG22	2.16	0.46
1:A:448:VAL:O	1:A:884:VAL:HG22	2.14	0.45
1:A:240:LEU:HB3	1:A:245:GLU:HG3	1.98	0.45
1:A:576:VAL:HG21	1:A:591:LEU:HD23	1.98	0.45
1:A:156:ASP:O	1:A:157:TYR:C	2.52	0.45
1:A:149:MET:HB2	1:A:153:ASP:HB3	1.97	0.45
1:A:905:VAL:HG13	1:A:935:ILE:HD12	1.98	0.45
1:A:338:HIS:CD2	1:A:339:GLU:HG2	2.51	0.45
1:A:559:LEU:HA	1:A:560:PRO:HD3	1.81	0.45
1:A:157:TYR:O	1:A:161:ASN:HB2	2.17	0.45
1:A:463:THR:HG23	1:A:563:PHE:HE1	1.82	0.45
1:A:183:ALA:HB2	1:A:273:GLU:HG2	1.98	0.45
1:A:111:LEU:HD11	1:A:127:VAL:HG11	1.99	0.45
1:A:199:THR:OG1	1:A:201:VAL:HB	2.17	0.45
1:A:239:ARG:CG	1:A:239:ARG:NH1	2.80	0.44
1:A:39:ALA:HA	1:A:40:PRO:HD2	1.81	0.44
1:A:228:GLN:HG2	1:A:230:LEU:N	2.32	0.44
1:A:782:LEU:O	1:A:785:ASP:HB2	2.17	0.44
1:A:19:ILE:HA	1:A:22:ALA:HB3	1.99	0.44
1:A:795:ASP:OD2	1:A:796:GLY:N	2.42	0.44
1:A:987:MET:N	1:A:988:PRO:CD	2.81	0.44
1:A:633:ASP:O	1:A:635:ALA:N	2.50	0.44
1:A:971:ARG:HG2	1:A:974:PRO:CG	2.23	0.44
1:A:182:TYR:HD1	1:A:271:GLY:O	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:420:MET:SD	1:A:499:PRO:HA	2.58	0.44
1:A:281:PHE:CZ	1:A:324:VAL:HG21	2.52	0.44
1:A:154:ILE:HG22	1:A:287:SER:HB3	2.00	0.44
1:A:40:PRO:HA	1:A:41:PRO:HD2	1.67	0.44
1:A:300:LEU:HD12	1:A:300:LEU:H	1.82	0.44
1:A:897:ILE:HD11	1:A:1030:ARG:HH11	1.83	0.44
1:A:626:ILE:HG12	1:A:628:PHE:HE1	1.76	0.44
1:A:377:LEU:O	1:A:380:PHE:HB2	2.19	0.43
1:A:28:LEU:O	1:A:30:LEU:N	2.50	0.43
1:A:491:ALA:O	1:A:494:ALA:HB3	2.18	0.43
1:A:43:VAL:HG12	1:A:44:THR:N	2.33	0.43
1:A:901:VAL:O	1:A:904:VAL:HG23	2.18	0.43
1:A:282:ASN:HD21	1:A:608:SER:HB3	1.83	0.43
1:A:758:TYR:CE1	1:A:770:LYS:HD3	2.52	0.43
1:A:183:ALA:N	1:A:271:GLY:O	2.41	0.43
1:A:914:LEU:O	1:A:915:ALA:C	2.57	0.43
1:A:783:PRO:O	1:A:786:ILE:HG12	2.18	0.43
1:A:672:VAL:O	1:A:673:GLU:HB3	2.18	0.43
1:A:410:ILE:HG22	1:A:411:VAL:N	2.34	0.43
1:A:158:VAL:HG22	1:A:162:MET:CE	2.49	0.43
1:A:431:THR:O	1:A:435:MET:N	2.48	0.43
1:A:813:SER:HB3	1:A:816:LEU:HD23	2.01	0.43
1:A:1021:PHE:O	1:A:1025:PHE:N	2.46	0.43
1:A:610:PHE:N	1:A:628:PHE:O	2.44	0.43
1:A:925:VAL:O	1:A:926:TYR:C	2.57	0.43
1:A:944:LEU:HB3	1:A:971:ARG:NH2	2.27	0.42
1:A:162:MET:HA	1:A:313:MET:SD	2.60	0.42
1:A:813:SER:HB3	1:A:816:LEU:CD2	2.48	0.42
1:A:445:ILE:HB	1:A:449:LEU:HD11	2.01	0.42
1:A:223:PRO:HA	1:A:224:PRO:HD3	1.89	0.42
1:A:180:SER:OG	1:A:273:GLU:HB2	2.19	0.42
1:A:325:TYR:HA	1:A:326:PRO:HD3	1.84	0.42
1:A:304:ALA:HA	1:A:307:ARG:HH21	1.84	0.42
1:A:897:ILE:O	1:A:900:SER:HB3	2.20	0.42
1:A:915:ALA:HB1	1:A:1005:THR:CG2	2.45	0.42
1:A:488:LEU:O	1:A:492:LEU:HG	2.20	0.42
1:A:965:LEU:O	1:A:968:VAL:HG12	2.19	0.42
1:A:310:LEU:C	1:A:312:LYS:H	2.23	0.42
1:A:365:THR:O	1:A:368:PRO:HD2	2.19	0.42
1:A:776:GLU:HG2	1:A:777:ALA:H	1.84	0.42
1:A:564:LEU:HD12	1:A:925:VAL:HB	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:467:TYR:CE1	1:A:868:LEU:HD21	2.54	0.42
1:A:330:THR:N	1:A:331:PRO:CD	2.80	0.42
1:A:23:GLY:O	1:A:27:ILE:HG13	2.20	0.42
1:A:963:ALA:O	1:A:964:THR:C	2.58	0.42
1:A:454:VAL:HB	1:A:455:PRO:HD3	2.00	0.42
1:A:988:PRO:O	1:A:991:ILE:HG12	2.19	0.41
1:A:45:ILE:HG23	1:A:129:VAL:HG22	2.00	0.41
1:A:519:MET:H	1:A:519:MET:HG2	1.58	0.41
1:A:278:ILE:HG13	1:A:613:ASN:HB3	2.01	0.41
1:A:300:LEU:HD12	1:A:300:LEU:N	2.34	0.41
1:A:114:ALA:HA	1:A:117:LEU:HD12	2.02	0.41
1:A:75:LEU:CD1	1:A:92:LEU:HD23	2.50	0.41
1:A:721:LEU:HD13	1:A:815:ARG:HB2	2.02	0.41
1:A:1013:THR:O	1:A:1015:THR:N	2.53	0.41
1:A:474:ILE:O	1:A:478:MET:N	2.49	0.41
1:A:151:GLN:HB3	1:A:285:PRO:HB3	2.03	0.41
1:A:216:ALA:HB1	1:A:234:ILE:O	2.21	0.41
1:A:605:ASN:HB3	1:A:637:ARG:HD3	2.02	0.41
1:A:1015:THR:C	1:A:1017:LEU:H	2.24	0.41
1:A:184:MET:CE	1:A:184:MET:HA	2.50	0.41
1:A:47:ALA:HB3	1:A:88:VAL:CG1	2.51	0.41
1:A:603:LYS:O	1:A:604:ASN:C	2.59	0.41
1:A:367:ILE:CB	1:A:368:PRO:HD3	2.37	0.41
1:A:183:ALA:HB2	1:A:273:GLU:HG3	2.02	0.41
1:A:475:VAL:O	1:A:478:MET:HB3	2.21	0.41
1:A:326:PRO:O	1:A:630:SER:HB2	2.20	0.41
1:A:605:ASN:ND2	1:A:637:ARG:HG2	2.35	0.41
1:A:975:ILE:H	1:A:975:ILE:HG12	1.65	0.41
1:A:616:GLY:C	1:A:618:ALA:H	2.24	0.41
1:A:883:VAL:HA	1:A:886:LEU:HD12	2.03	0.41
1:A:818:ARG:NH1	1:A:823:PRO:HG3	2.36	0.41
1:A:349:ILE:O	1:A:353:LEU:HG	2.21	0.41
1:A:65:ILE:C	1:A:67:GLN:H	2.24	0.41
1:A:100:ALA:O	1:A:103:ALA:HB3	2.21	0.41
1:A:908:GLY:HA2	1:A:1014:ALA:HB2	2.03	0.41
1:A:445:ILE:O	1:A:449:LEU:HG	2.21	0.41
1:A:685:ILE:HG22	1:A:687:GLN:HG3	2.02	0.41
1:A:574:THR:HB	1:A:665:ALA:CB	2.51	0.40
1:A:845:GLU:HG3	1:A:859:TRP:HE1	1.86	0.40
1:A:610:PHE:HB3	1:A:628:PHE:HB2	2.03	0.40
1:A:1001:ASN:O	1:A:1005:THR:N	2.50	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:144:ASN:ND2	1:A:149:MET:N	2.69	0.40
1:A:211:ASN:HB2	1:A:240:LEU:HD12	2.03	0.40
1:A:157:TYR:CD2	1:A:157:TYR:C	2.95	0.40
1:A:184:MET:HB2	1:A:762:PHE:CD2	2.57	0.40
1:A:782:LEU:HB3	1:A:783:PRO:HD2	2.03	0.40
1:A:164:ASP:O	1:A:168:ARG:NH2	2.55	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	1030/1049 (98%)	782 (76%)	196 (19%)	52 (5%)	<b>3</b>	31

All (52) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	28	LEU
1	A	133	SER
1	A	226	LYS
1	A	227	GLY
1	A	255	GLN
1	A	352	PHE
1	A	471	SER
1	A	634	TRP
1	A	868	LEU
1	A	992	SER
1	A	29	LYS
1	A	41	PRO
1	A	80	SER
1	A	254	ASN
1	A	308	ALA

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Mol	Chain	Res	Type
1	A	383	LEU
1	A	406	VAL
1	A	674	LEU
1	A	862	MET
1	A	897	ILE
1	A	1000	GLN
1	A	1014	ALA
1	A	35	TYR
1	A	50	PRO
1	A	384	ALA
1	A	1017	LEU
1	A	66	GLU
1	A	387	GLY
1	A	439	GLN
1	A	579	PRO
1	A	671	ILE
1	A	991	ILE
1	A	1025	PHE
1	A	62	THR
1	A	306	ILE
1	A	716	VAL
1	A	796	GLY
1	A	833	PRO
1	A	834	GLY
1	A	1016	VAL
1	A	428	LYS
1	A	464	GLY
1	A	759	VAL
1	A	852	PRO
1	A	51	GLY
1	A	320	GLY
1	A	427	PRO
1	A	173	GLY
1	A	920	GLY
1	A	249	ILE
1	A	326	PRO
1	A	337	ILE

### 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	838/855 (98%)	716 (85%)	122 (15%)	4 26

All (122) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	PHE
1	A	15	ILE
1	A	17	ILE
1	A	19	ILE
1	A	21	LEU
1	A	38	ILE
1	A	48	SER
1	A	58	GLN
1	A	60	THR
1	A	64	VAL
1	A	73	ASP
1	A	89	GLN
1	A	91	THR
1	A	98	THR
1	A	102	ILE
1	A	109	ASN
1	A	118	LEU
1	A	122	VAL
1	A	138	MET
1	A	149	MET
1	A	151	GLN
1	A	152	GLU
1	A	164	ASP
1	A	182	TYR
1	A	187	TRP
1	A	195	LYS
1	A	196	PHE
1	A	197	GLN
1	A	205	THR
1	A	222	THR
1	A	225	VAL
1	A	241	THR
1	A	243	THR
1	A	246	PHE

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Mol	Chain	Res	Type
1	A	259	ARG
1	A	267	LYS
1	A	276	ASP
1	A	291	ILE
1	A	301	ASP
1	A	302	THR
1	A	307	ARG
1	A	321	LEU
1	A	323	ILE
1	A	338	HIS
1	A	343	THR
1	A	352	PHE
1	A	355	MET
1	A	362	PHE
1	A	365	THR
1	A	386	PHE
1	A	389	SER
1	A	400	LEU
1	A	408	ASP
1	A	410	ILE
1	A	439	GLN
1	A	452	VAL
1	A	459	PHE
1	A	463	THR
1	A	476	SER
1	A	486	LEU
1	A	502	LYS
1	A	519	MET
1	A	538	THR
1	A	544	LEU
1	A	546	LEU
1	A	564	LEU
1	A	566	ASP
1	A	568	ASP
1	A	573	MET
1	A	583	THR
1	A	589	LYS
1	A	595	THR
1	A	603	LYS
1	A	608	SER
1	A	613	ASN
1	A	620	ARG

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Mol	Chain	Res	Type
1	A	624	THR
1	A	641	GLU
1	A	658	ILE
1	A	659	LYS
1	A	660	ASP
1	A	666	PHE
1	A	668	LEU
1	A	673	GLU
1	A	674	LEU
1	A	695	LEU
1	A	696	THR
1	A	702	LEU
1	A	713	LEU
1	A	715	SER
1	A	723	ASP
1	A	724	THR
1	A	743	ILE
1	A	759	VAL
1	A	760	ASN
1	A	768	VAL
1	A	773	VAL
1	A	778	LYS
1	A	795	ASP
1	A	797	GLN
1	A	799	VAL
1	A	804	PHE
1	A	815	ARG
1	A	843	LEU
1	A	853	THR
1	A	872	GLN
1	A	880	SER
1	A	894	SER
1	A	901	VAL
1	A	921	LEU
1	A	934	THR
1	A	968	VAL
1	A	971	ARG
1	A	972	LEU
1	A	975	ILE
1	A	980	LEU
1	A	990	VAL
1	A	991	ILE

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Mol	Chain	Res	Type
1	A	993	THR
1	A	1016	VAL
1	A	1017	LEU
1	A	1028	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	63	GLN
1	A	67	GLN
1	A	68	ASN
1	A	106	GLN
1	A	124	GLN
1	A	125	GLN
1	A	176	GLN
1	A	194	ASN
1	A	218	GLN
1	A	229	GLN
1	A	237	GLN
1	A	517	ASN
1	A	604	ASN
1	A	605	ASN
1	A	622	GLN
1	A	642	ASN
1	A	692	HIS
1	A	733	GLN
1	A	871	ASN
1	A	928	GLN
1	A	1001	ASN

### 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 [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	DXC	A	2034	-	28,31,31	2.21	3 (10%)	46,49,49	1.73	7 (15%)

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
2	DXC	A	2034	-	-	0/7/71/71	0/4/4/4

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	2034	DXC	C16-C17	2.37	1.59	1.54
2	A	2034	DXC	C15-C11	2.63	1.60	1.54
2	A	2034	DXC	C16-C15	10.83	1.84	1.54

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	2034	DXC	C15-C16-C17	-3.63	97.81	105.12
2	A	2034	DXC	C16-C15-C11	-2.81	99.46	105.12
2	A	2034	DXC	C22-C21-C19	-2.17	112.19	114.75
2	A	2034	DXC	C21-C19-C17	2.59	115.68	110.24
2	A	2034	DXC	C16-C17-C19	2.90	117.23	112.05
2	A	2034	DXC	C4-C10-C9	2.95	115.68	112.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	2034	DXC	C15-C11-C12	6.53	110.09	103.60

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
2	A	2034	DXC	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	1032/1049 (98%)	-0.12	30 (2%) 55 42	48, 74, 98, 106	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	866	GLU	3.9
1	A	714	THR	3.9
1	A	833	PRO	3.6
1	A	423	GLU	3.3
1	A	425	LEU	3.2
1	A	709	HIS	3.1
1	A	708	LYS	3.1
1	A	835	LYS	3.1
1	A	852	PRO	3.0
1	A	712	MET	2.9
1	A	701	GLN	2.9
1	A	843	LEU	2.9
1	A	865	GLN	2.7
1	A	834	GLY	2.6
1	A	500	ILE	2.6
1	A	420	MET	2.5
1	A	460	GLY	2.4
1	A	840	ALA	2.4
1	A	832	ALA	2.3
1	A	842	GLU	2.3
1	A	841	MET	2.3
1	A	408	ASP	2.3
1	A	448	VAL	2.2
1	A	956	GLU	2.2
1	A	851	LEU	2.1
1	A	711	ASP	2.1
1	A	870	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	702	LEU	2.1
1	A	303	ALA	2.1
1	A	831	ALA	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 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<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
2	DXC	A	2034	28/28	0.87	0.26	-1.09	99,104,105,105	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.