Full wwPDB X-ray Structure Validation Report ⓘ

Oct 4, 2023 – 04:18 AM EDT

PDB ID	:	6PEX
Title	:	An aldo keto reductase with 2-keto- L-gulonate reductase activity
Authors	:	Yong, J.; Crystal, S.; Robert, D.H.; John, B.B.
Deposited on	:	2019-06-21
Resolution	:	1.58 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

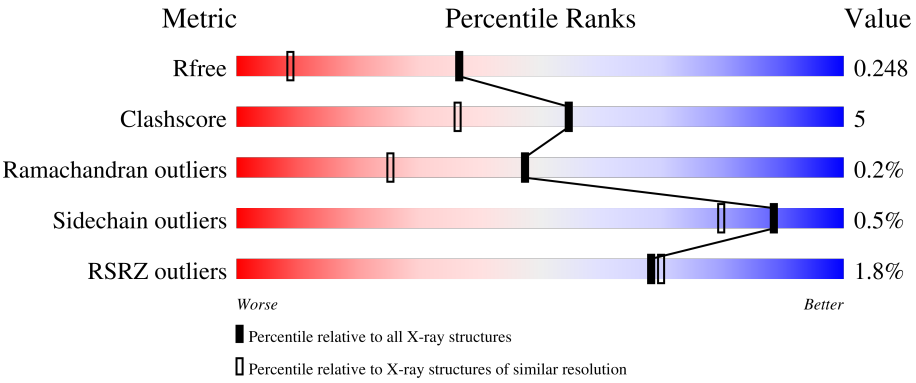
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

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 1.58 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	5534 (1.60-1.56)
Clashscore	141614	5861 (1.60-1.56)
Ramachandran outliers	138981	5708 (1.60-1.56)
Sidechain outliers	138945	5703 (1.60-1.56)
RSRZ outliers	127900	5431 (1.60-1.56)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	329	2% 82% 12% 5%
1	B	329	2% 85% 10% 5%
1	C	329	2% 84% 11% 5%
1	D	329	% 84% 11% 5%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 10588 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 aldo keto reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	312	Total	C	N	O	S	0	0	0
			2384	1517	411	445	11			
1	B	312	Total	C	N	O	S	0	0	0
			2382	1515	409	447	11			
1	C	312	Total	C	N	O	S	0	0	0
			2372	1509	409	443	11			
1	D	312	Total	C	N	O	S	0	0	0
			2373	1510	408	444	11			

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-15	HIS	-	expression tag	UNP A5CAL1
A	-14	HIS	-	expression tag	UNP A5CAL1
A	-13	HIS	-	expression tag	UNP A5CAL1
A	-12	HIS	-	expression tag	UNP A5CAL1
A	-11	HIS	-	expression tag	UNP A5CAL1
A	-10	SER	-	expression tag	UNP A5CAL1
A	-9	SER	-	expression tag	UNP A5CAL1
A	-8	GLY	-	expression tag	UNP A5CAL1
A	-7	LEU	-	expression tag	UNP A5CAL1
A	-6	VAL	-	expression tag	UNP A5CAL1
A	-5	PRO	-	expression tag	UNP A5CAL1
A	-4	ARG	-	expression tag	UNP A5CAL1
A	-3	GLY	-	expression tag	UNP A5CAL1
A	-2	SER	-	expression tag	UNP A5CAL1
A	-1	HIS	-	expression tag	UNP A5CAL1
A	0	MET	-	expression tag	UNP A5CAL1
B	-15	HIS	-	expression tag	UNP A5CAL1
B	-14	HIS	-	expression tag	UNP A5CAL1
B	-13	HIS	-	expression tag	UNP A5CAL1
B	-12	HIS	-	expression tag	UNP A5CAL1
B	-11	HIS	-	expression tag	UNP A5CAL1

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-10	SER	-	expression tag	UNP A5CAL1
B	-9	SER	-	expression tag	UNP A5CAL1
B	-8	GLY	-	expression tag	UNP A5CAL1
B	-7	LEU	-	expression tag	UNP A5CAL1
B	-6	VAL	-	expression tag	UNP A5CAL1
B	-5	PRO	-	expression tag	UNP A5CAL1
B	-4	ARG	-	expression tag	UNP A5CAL1
B	-3	GLY	-	expression tag	UNP A5CAL1
B	-2	SER	-	expression tag	UNP A5CAL1
B	-1	HIS	-	expression tag	UNP A5CAL1
B	0	MET	-	expression tag	UNP A5CAL1
C	-15	HIS	-	expression tag	UNP A5CAL1
C	-14	HIS	-	expression tag	UNP A5CAL1
C	-13	HIS	-	expression tag	UNP A5CAL1
C	-12	HIS	-	expression tag	UNP A5CAL1
C	-11	HIS	-	expression tag	UNP A5CAL1
C	-10	SER	-	expression tag	UNP A5CAL1
C	-9	SER	-	expression tag	UNP A5CAL1
C	-8	GLY	-	expression tag	UNP A5CAL1
C	-7	LEU	-	expression tag	UNP A5CAL1
C	-6	VAL	-	expression tag	UNP A5CAL1
C	-5	PRO	-	expression tag	UNP A5CAL1
C	-4	ARG	-	expression tag	UNP A5CAL1
C	-3	GLY	-	expression tag	UNP A5CAL1
C	-2	SER	-	expression tag	UNP A5CAL1
C	-1	HIS	-	expression tag	UNP A5CAL1
C	0	MET	-	expression tag	UNP A5CAL1
D	-15	HIS	-	expression tag	UNP A5CAL1
D	-14	HIS	-	expression tag	UNP A5CAL1
D	-13	HIS	-	expression tag	UNP A5CAL1
D	-12	HIS	-	expression tag	UNP A5CAL1
D	-11	HIS	-	expression tag	UNP A5CAL1
D	-10	SER	-	expression tag	UNP A5CAL1
D	-9	SER	-	expression tag	UNP A5CAL1
D	-8	GLY	-	expression tag	UNP A5CAL1
D	-7	LEU	-	expression tag	UNP A5CAL1
D	-6	VAL	-	expression tag	UNP A5CAL1
D	-5	PRO	-	expression tag	UNP A5CAL1
D	-4	ARG	-	expression tag	UNP A5CAL1
D	-3	GLY	-	expression tag	UNP A5CAL1
D	-2	SER	-	expression tag	UNP A5CAL1
D	-1	HIS	-	expression tag	UNP A5CAL1

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Chain	Residue	Modelled	Actual	Comment	Reference
D	0	MET	-	expression tag	UNP A5CAL1

- Molecule 2 is water.

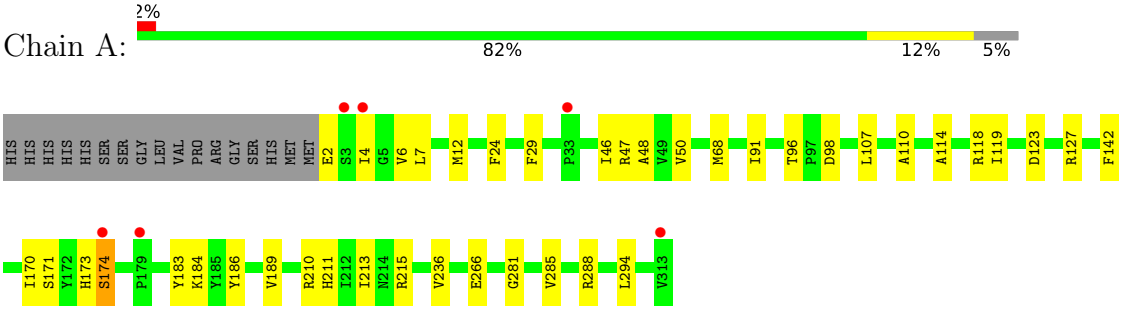
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	300	Total O 300 300	0	0
2	B	220	Total O 220 220	0	0
2	C	266	Total O 266 266	0	0
2	D	291	Total O 291 291	0	0

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.

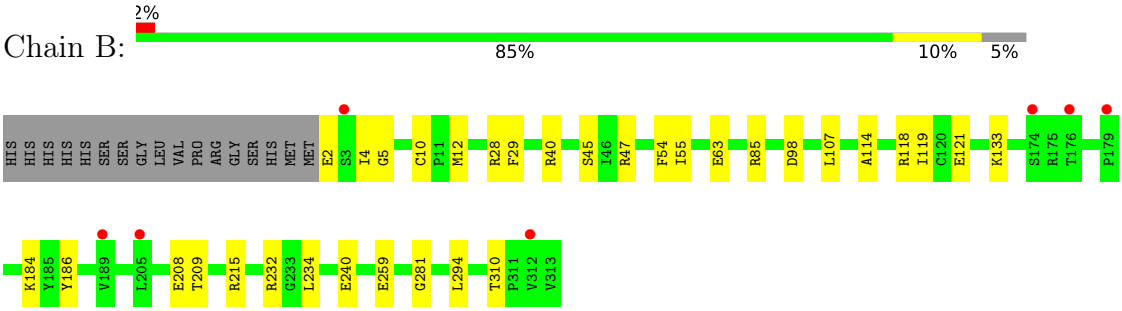
• Molecule 1: aldo keto reductase

Chain A:



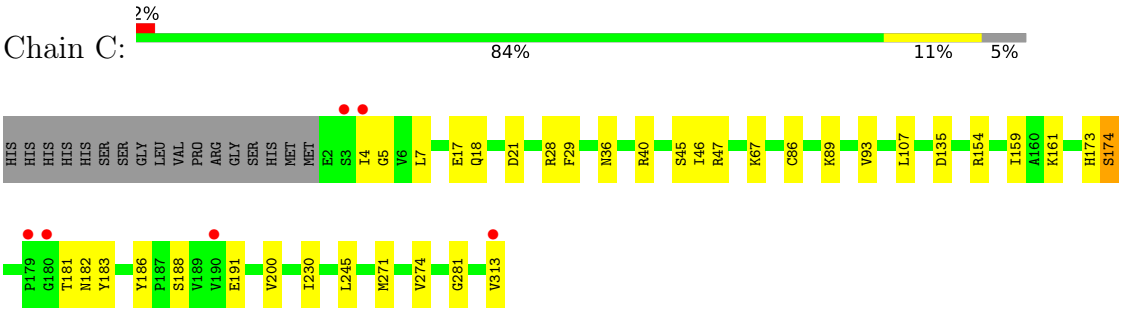
• Molecule 1: aldo keto reductase

Chain B:



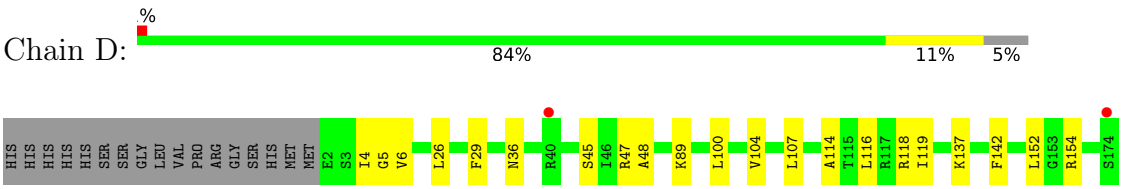
• Molecule 1: aldo keto reductase

Chain C:



• Molecule 1: aldo keto reductase

Chain D:




Q197
N214
R215
L221
I230
L245
V246
E247
P265
L268
L269
A270
M271
D272
N273
L277
G281
V313

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	73.05Å 85.72Å 112.90Å 90.00° 89.91° 90.00°	Depositor
Resolution (Å)	27.70 – 1.58 27.70 – 1.58	Depositor EDS
% Data completeness (in resolution range)	99.5 (27.70-1.58) 99.1 (27.70-1.58)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.84 (at 1.58Å)	Xtriage
Refinement program	PHENIX 1.15.2_3472	Depositor
R, R_{free}	0.230 , 0.251 0.242 , 0.248	Depositor DCC
R_{free} test set	9168 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å^2)	16.1	Xtriage
Anisotropy	0.385	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 35.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.447 for h,-k,-l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	10588	wwPDB-VP
Average B, all atoms (Å^2)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 18.04% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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	0.24	0/2427	0.42	0/3293
1	B	0.24	0/2424	0.43	0/3289
1	C	0.24	0/2415	0.42	0/3280
1	D	0.24	0/2415	0.44	0/3279
All	All	0.24	0/9681	0.43	0/13141

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

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	2384	0	2424	29	0
1	B	2382	0	2421	20	0
1	C	2372	0	2398	28	0
1	D	2373	0	2403	29	0
2	A	300	0	0	8	0
2	B	220	0	0	3	0
2	C	266	0	0	11	0
2	D	291	0	0	13	0
All	All	10588	0	9646	100	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 5.

All (100) 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:D:271:MET:SD	2:D:674:HOH:O	2.32	0.86
1:C:86:CYS:SG	2:C:629:HOH:O	2.41	0.79
1:B:114:ALA:HA	1:B:119:ILE:HD12	1.69	0.74
1:C:4:ILE:HG23	1:C:47:ARG:HD3	1.69	0.72
1:C:28:ARG:NH2	2:C:401:HOH:O	2.21	0.72
1:C:181:THR:HG22	1:C:183:TYR:H	1.56	0.71
1:D:118:ARG:NH1	2:D:404:HOH:O	2.24	0.71
1:D:6:VAL:HG22	1:D:48:ALA:HB3	1.73	0.69
1:B:4:ILE:HG23	1:B:47:ARG:HD3	1.77	0.66
1:C:67:LYS:NZ	1:D:221:LEU:O	2.29	0.65
1:D:230:ILE:C	1:D:230:ILE:HD12	2.18	0.64
1:D:4:ILE:HG23	1:D:47:ARG:HD3	1.80	0.63
1:C:67:LYS:HD3	1:D:197:GLN:HG2	1.81	0.62
1:B:234:LEU:HD23	1:B:259:GLU:HG2	1.81	0.61
1:C:154:ARG:NH2	2:C:419:HOH:O	2.32	0.61
1:A:118:ARG:NH1	2:A:413:HOH:O	2.33	0.61
1:C:182:ASN:ND2	2:C:422:HOH:O	2.34	0.60
1:A:210:ARG:NH2	2:A:419:HOH:O	2.35	0.59
1:A:114:ALA:HA	1:A:119:ILE:HD12	1.85	0.58
1:D:6:VAL:HB	2:D:516:HOH:O	2.05	0.57
1:B:208:GLU:HG2	1:B:209:THR:HG23	1.86	0.57
1:C:7:LEU:HB2	1:C:46:ILE:HG12	1.86	0.56
1:B:5:GLY:HA3	1:B:45:SER:O	2.06	0.55
1:B:98:ASP:N	2:B:414:HOH:O	2.39	0.54
1:C:18:GLN:NE2	2:C:430:HOH:O	2.41	0.53
1:B:215:ARG:NH1	1:B:240:GLU:OE1	2.41	0.52
1:A:4:ILE:HG23	1:A:47:ARG:HD3	1.91	0.52
1:A:68:MET:HE3	1:A:91:ILE:HG21	1.91	0.52
1:D:114:ALA:HA	1:D:119:ILE:HD12	1.91	0.52
1:A:171:SER:HB3	1:A:184:LYS:HB3	1.92	0.51
1:D:247:GLU:HB3	2:D:500:HOH:O	2.11	0.51
1:A:142:PHE:N	2:A:405:HOH:O	2.43	0.51
1:D:214:ASN:ND2	2:D:424:HOH:O	2.43	0.51
1:D:107:LEU:HD22	1:D:281:GLY:HA2	1.93	0.50
1:C:173:HIS:O	1:C:174:SER:HB2	2.10	0.50
1:B:232:ARG:NH1	2:B:419:HOH:O	2.43	0.50
1:B:118:ARG:NH1	1:B:121:GLU:OE1	2.40	0.49
1:A:98:ASP:N	2:A:434:HOH:O	2.45	0.49
1:A:6:VAL:HG22	1:A:48:ALA:HB3	1.95	0.48
1:A:173:HIS:ND1	1:A:186:TYR:O	2.41	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:93:VAL:HB	1:C:313:VAL:HG22	1.95	0.48
1:B:294:LEU:HD21	1:B:310:THR:HG21	1.95	0.48
1:D:26:LEU:HA	2:D:516:HOH:O	2.13	0.48
1:B:63:GLU:OE2	1:B:85:ARG:HD3	2.14	0.47
1:D:5:GLY:HA3	1:D:45:SER:O	2.14	0.47
1:A:123:ASP:OD1	1:A:127:ARG:NE	2.48	0.47
1:B:40:ARG:HH22	1:D:270:ALA:HB2	1.80	0.47
1:A:173:HIS:CD2	1:A:189:VAL:HG22	2.50	0.46
1:A:173:HIS:CG	1:A:174:SER:N	2.83	0.46
1:B:107:LEU:HD22	1:B:281:GLY:HA2	1.96	0.46
1:A:288:ARG:NH2	2:A:440:HOH:O	2.49	0.46
1:D:104:VAL:CG1	1:D:230:ILE:HD13	2.46	0.45
1:B:2:GLU:O	1:B:4:ILE:N	2.48	0.45
1:D:230:ILE:HD12	1:D:230:ILE:O	2.17	0.45
1:C:17:GLU:HG2	2:C:401:HOH:O	2.16	0.45
1:C:154:ARG:NH1	2:C:447:HOH:O	2.50	0.45
1:A:215:ARG:NH2	2:A:436:HOH:O	2.45	0.45
1:A:266:GLU:H	1:A:266:GLU:CD	2.20	0.45
1:C:40:ARG:HA	1:C:40:ARG:HD3	1.80	0.45
1:C:159:ILE:HD13	1:C:200:VAL:HG11	1.99	0.45
1:C:161:LYS:NZ	2:C:409:HOH:O	2.26	0.45
1:C:5:GLY:HA3	1:C:45:SER:O	2.17	0.44
1:C:173:HIS:ND1	1:C:186:TYR:O	2.45	0.44
1:A:211:HIS:NE2	2:A:408:HOH:O	2.36	0.44
1:C:188:SER:OG	1:C:191:GLU:HB2	2.17	0.44
1:D:89:LYS:NZ	2:D:440:HOH:O	2.50	0.44
1:A:119:ILE:HG23	1:D:277:LEU:HD21	2.00	0.44
1:C:135:ASP:O	2:C:402:HOH:O	2.21	0.44
1:C:107:LEU:HD22	1:C:281:GLY:HA2	2.00	0.44
1:A:107:LEU:HD22	1:A:281:GLY:HA2	2.00	0.44
1:B:28:ARG:NH1	2:B:428:HOH:O	2.48	0.43
1:B:133:LYS:HA	1:B:133:LYS:HD3	1.86	0.43
1:D:152:LEU:HD12	2:D:448:HOH:O	2.18	0.43
1:A:96:THR:HA	1:A:294:LEU:HD23	2.00	0.43
1:A:4:ILE:O	1:A:24:PHE:HB3	2.19	0.43
1:C:245:LEU:HB3	2:C:451:HOH:O	2.18	0.43
1:A:170:ILE:HB	1:A:183:TYR:CD1	2.54	0.43
1:D:245:LEU:HD13	2:D:674:HOH:O	2.19	0.43
1:D:273:ASN:ND2	2:D:431:HOH:O	2.47	0.42
1:B:10:CYS:O	1:B:12:MET:N	2.52	0.42
1:B:208:GLU:HG2	1:B:209:THR:N	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2:GLU:N	2:A:443:HOH:O	2.52	0.42
1:B:54:PHE:CD2	1:B:55:ILE:HG13	2.54	0.42
1:B:184:LYS:HG2	1:B:186:TYR:CE2	2.55	0.42
1:C:271:MET:HB2	1:C:274:VAL:HG22	2.02	0.42
1:A:285:VAL:HG11	1:C:89:LYS:HG2	2.02	0.42
1:D:137:LYS:NZ	2:D:416:HOH:O	2.39	0.42
1:D:265:PRO:HD2	1:D:268:LEU:HD12	2.02	0.41
1:A:213:ILE:HB	1:A:236:VAL:HG22	2.03	0.41
1:A:2:GLU:O	1:A:4:ILE:N	2.53	0.41
1:D:116:LEU:HD23	1:D:142:PHE:HA	2.03	0.41
1:A:12:MET:HE1	1:A:50:VAL:O	2.21	0.41
1:A:110:ALA:HB1	1:D:119:ILE:HD13	2.03	0.41
1:C:36:ASN:O	1:C:40:ARG:HG2	2.21	0.41
1:D:215:ARG:NH1	2:D:445:HOH:O	2.53	0.41
1:A:7:LEU:HB2	1:A:46:ILE:HG12	2.03	0.40
1:C:230:ILE:H	1:C:230:ILE:HG13	1.75	0.40
1:C:21:ASP:HB2	2:C:401:HOH:O	2.20	0.40
1:D:36:ASN:H	1:D:36:ASN:HD22	1.69	0.40
1:D:154:ARG:NH1	2:D:420:HOH:O	2.41	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	310/329 (94%)	294 (95%)	15 (5%)	1 (0%)	41 21
1	B	310/329 (94%)	294 (95%)	16 (5%)	0	100 100
1	C	310/329 (94%)	294 (95%)	15 (5%)	1 (0%)	41 21
1	D	310/329 (94%)	294 (95%)	16 (5%)	0	100 100
All	All	1240/1316 (94%)	1176 (95%)	62 (5%)	2 (0%)	47 25

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	174	SER
1	C	174	SER

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	262/281 (93%)	261 (100%)	1 (0%)	91	84
1	B	262/281 (93%)	261 (100%)	1 (0%)	91	84
1	C	259/281 (92%)	258 (100%)	1 (0%)	91	84
1	D	259/281 (92%)	257 (99%)	2 (1%)	81	68
All	All	1042/1124 (93%)	1037 (100%)	5 (0%)	88	80

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	PHE
1	B	29	PHE
1	C	29	PHE
1	D	29	PHE
1	D	100	LEU

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

Mol	Chain	Res	Type
1	C	18	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

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^{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	312/329 (94%)	-0.13	6 (1%)	66	68	5, 17, 31, 46	0
1	B	312/329 (94%)	0.12	7 (2%)	62	63	13, 23, 35, 53	0
1	C	312/329 (94%)	0.02	6 (1%)	66	68	12, 21, 34, 44	0
1	D	312/329 (94%)	-0.14	3 (0%)	82	83	5, 17, 30, 43	0
All	All	1248/1316 (94%)	-0.03	22 (1%)	68	70	5, 20, 33, 53	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	174	SER	3.4
1	D	174	SER	3.4
1	B	179	PRO	3.2
1	B	3	SER	3.2
1	B	176	THR	3.1
1	C	179	PRO	3.0
1	A	4	ILE	2.8
1	C	3	SER	2.8
1	C	180	GLY	2.5
1	C	313	VAL	2.5
1	A	313	VAL	2.4
1	A	174	SER	2.4
1	A	33	PRO	2.3
1	B	205	LEU	2.3
1	B	312	VAL	2.2
1	C	190	VAL	2.2
1	C	4	ILE	2.2
1	B	189	VAL	2.2
1	D	313	VAL	2.1
1	D	40	ARG	2.0
1	A	3	SER	2.0
1	A	179	PRO	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 monosaccharides in this entry.

6.4 Ligands (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.