Full wwPDB X-ray Structure Validation Report (i)

Jan 1, 2017 – 04:42 AM EST

PDB ID	:	5TZ6
Title	:	Crystal Structure of CurJ Dehydratase H978F Inactive Mutant In Complex with Compound 21
Authors	:	Dodge, G.J.; Smith, J.L.
Deposited on	:	2016-11-21
Resolution	:	2.40 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:

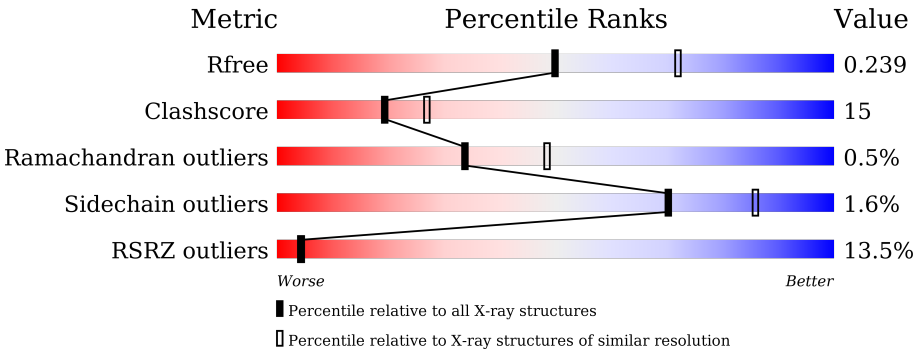
MolProbity	:	4.02b-467
Mogul	:	1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20028442
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20028442

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 2.40 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	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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	308	2% 76% 17% • 6%
1	B	308	23% 61% 30% 9%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4611 atoms, of which 13 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.

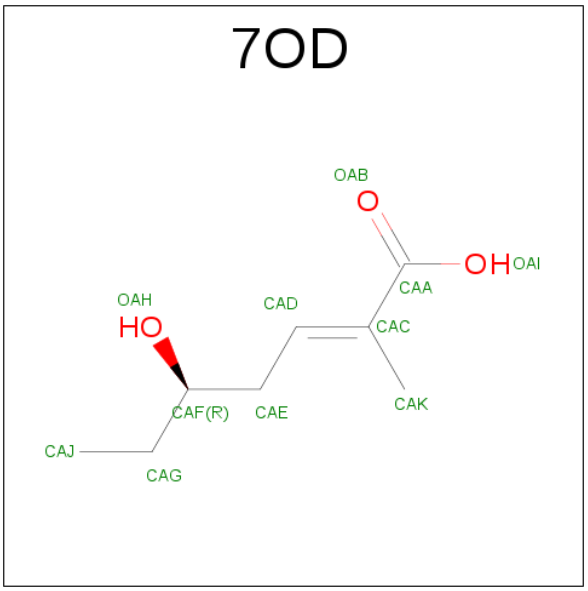
- Molecule 1 is a protein called CurJ.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	290	Total	C	N	O	S	0	0	0
			2311	1490	378	435	8			
1	B	281	Total	C	N	O	S	0	0	0
			2243	1448	365	422	8			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	938	SER	-	expression tag	UNP F4Y426
A	939	ASN	-	expression tag	UNP F4Y426
A	940	ALA	-	expression tag	UNP F4Y426
A	978	PHE	HIS	engineered mutation	UNP F4Y426
B	938	SER	-	expression tag	UNP F4Y426
B	939	ASN	-	expression tag	UNP F4Y426
B	940	ALA	-	expression tag	UNP F4Y426
B	978	PHE	HIS	engineered mutation	UNP F4Y426

- Molecule 2 is (2E,5R)-5-hydroxy-2-methylhept-2-enoic acid (three-letter code: 7OD) (formula: $C_8H_{14}O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C H O 24 8 13 3	0	0

- Molecule 3 is water.

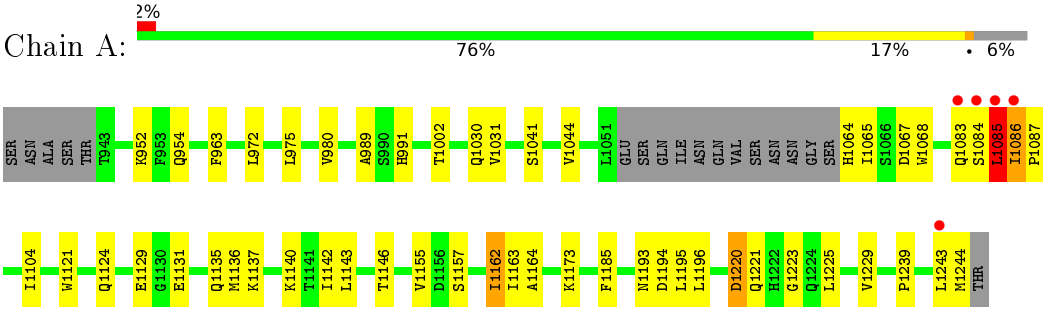
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	30	Total O 30 30	0	0
3	B	3	Total O 3 3	0	0

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.

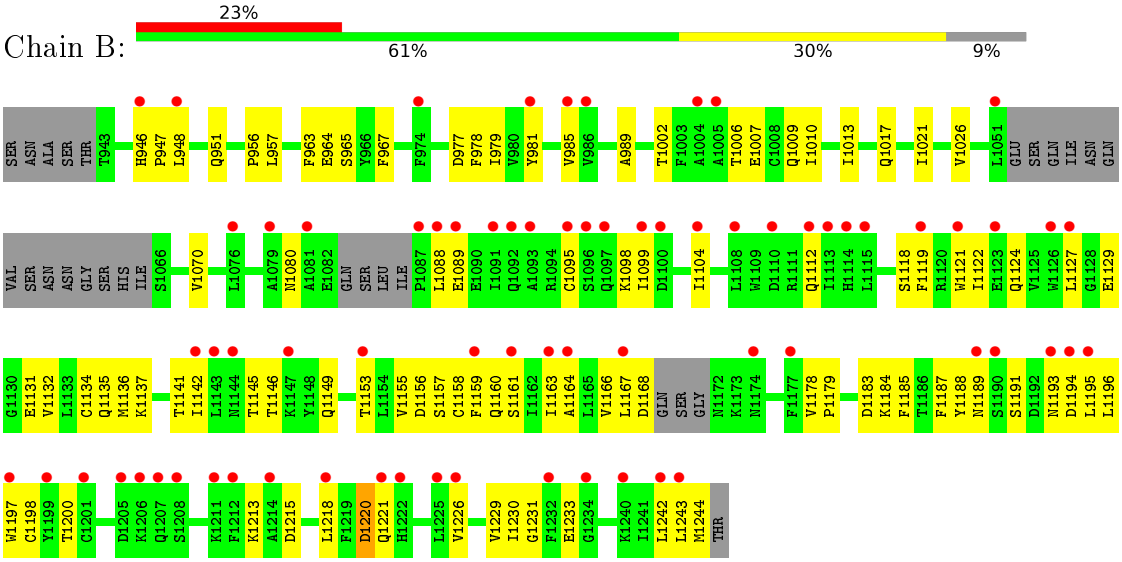
- Molecule 1: CurJ

Chain A:



- Molecule 1: CurJ

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	47.55Å 70.58Å 176.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.91 – 2.40 45.90 – 2.39	Depositor EDS
% Data completeness (in resolution range)	99.8 (45.91-2.40) 94.6 (45.90-2.39)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.29 (at 2.39Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.199 , 0.242 0.196 , 0.239	Depositor DCC
R_{free} test set	1899 reflections (8.33%)	DCC
Wilson B-factor (Å^2)	57.3	Xtriage
Anisotropy	0.414	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 77.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	4611	wwPDB-VP
Average B, all atoms (Å^2)	114.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.05% 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

Bond lengths and bond angles in the following residue types are not validated in this section: 7OD

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	0.30	0/2365	0.50	0/3211
1	B	0.26	0/2294	0.45	0/3111
All	All	0.28	0/4659	0.47	0/6322

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	2311	0	2279	60	1
1	B	2243	0	2209	80	1
2	A	11	13	0	0	0
3	A	30	0	0	4	0
3	B	3	0	0	3	0
All	All	4598	13	4488	137	1

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 15.

All (137) 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:1084:SER:HB2	1:A:1085:LEU:HA	1.45	0.98
1:A:1086:ILE:HD11	1:A:1223:GLY:HA3	1.50	0.94
1:A:1084:SER:N	1:A:1085:LEU:HB2	1.87	0.90
1:A:1084:SER:HB2	1:A:1085:LEU:CA	2.05	0.86
1:A:1084:SER:H	1:A:1085:LEU:HB2	1.40	0.84
1:B:1007:GLU:HG3	1:B:1188:TYR:HA	1.62	0.82
1:B:989:ALA:HB1	1:B:1155:VAL:HG12	1.64	0.80
1:A:1086:ILE:H	1:A:1087:PRO:HA	1.47	0.79
1:B:1142:ILE:CG2	1:B:1145:THR:HG23	2.12	0.78
1:B:1142:ILE:HG21	1:B:1145:THR:HG23	1.64	0.78
1:B:1197:TRP:O	1:B:1218:LEU:HD12	1.83	0.77
1:A:1195:LEU:HD13	1:A:1221:GLN:HE22	1.53	0.73
1:B:1178:VAL:HG11	1:B:1242:LEU:HD21	1.71	0.71
1:B:1163:ILE:HA	1:B:1166:VAL:HG22	1.73	0.70
1:B:1189:ASN:O	1:B:1226:VAL:HG13	1.92	0.70
1:B:1137:LYS:HE2	1:B:1195:LEU:HD12	1.75	0.69
1:B:977:ASP:OD2	1:B:1118:SER:OG	2.08	0.68
1:B:985:VAL:HG21	1:B:1243:LEU:HD21	1.76	0.68
1:B:963:PHE:HZ	1:B:1002:THR:HG23	1.62	0.64
1:A:1084:SER:CA	1:A:1085:LEU:HB2	2.28	0.64
1:A:1086:ILE:HD11	1:A:1223:GLY:CA	2.26	0.63
1:B:1119:PHE:CZ	1:B:1153:THR:HG22	2.32	0.63
1:B:1215:ASP:OD1	1:B:1230:ILE:HA	1.98	0.63
1:A:1135:GLN:NE2	1:A:1195:LEU:HD11	2.14	0.63
1:A:1065:ILE:HG21	1:B:956:PRO:HG3	1.82	0.62
1:A:980:VAL:CG1	1:A:1243:LEU:HD11	2.31	0.61
1:A:980:VAL:HB	1:A:1243:LEU:HD11	1.82	0.61
1:B:1122:ILE:HG12	1:B:1136:MET:CE	2.31	0.61
1:A:1129:GLU:OE1	1:A:1129:GLU:HA	2.01	0.60
1:B:951:GLN:HG2	1:B:964:GLU:HB3	1.83	0.60
1:B:1009:GLN:OE1	1:B:1184:LYS:HD2	2.00	0.60
1:B:1099:ILE:HG21	3:B:1303:HOH:O	2.01	0.60
1:B:1193:ASN:ND2	1:B:1220:ASP:OD1	2.35	0.59
1:A:1243:LEU:O	1:A:1244:MET:HB2	2.02	0.59
1:B:1197:TRP:HE1	1:B:1221:GLN:CD	2.07	0.58
1:A:1085:LEU:HD13	1:A:1223:GLY:O	2.03	0.58
1:B:1185:PHE:CD2	1:B:1229:VAL:HG22	2.39	0.57
1:A:1085:LEU:HA	1:A:1086:ILE:HG13	1.87	0.57
1:B:1132:VAL:HG12	1:B:1200:THR:HB	1.86	0.57
1:B:989:ALA:HB1	1:B:1155:VAL:CG1	2.35	0.56
1:A:1140:LYS:HE3	3:A:1401:HOH:O	2.06	0.56
1:A:1142:ILE:O	1:A:1143:LEU:HD23	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1157:SER:HA	1:B:1160:GLN:CD	2.26	0.55
1:B:1121:TRP:O	1:B:1136:MET:HA	2.06	0.55
1:A:1088:LEU:HD12	1:A:1088:LEU:O	2.06	0.55
1:B:1095:CYS:SG	1:B:1131:GLU:HG3	2.46	0.55
1:B:1007:GLU:CG	1:B:1188:TYR:HA	2.36	0.54
1:B:1213:LYS:HE2	1:B:1233:GLU:OE2	2.08	0.54
1:B:1132:VAL:CG1	1:B:1200:THR:HB	2.37	0.54
1:B:1129:GLU:OE1	1:B:1129:GLU:HA	2.08	0.54
1:B:1088:LEU:O	1:B:1088:LEU:HD12	2.08	0.53
1:B:1124:GLN:HB2	1:B:1135:GLN:HB3	1.89	0.53
1:B:1104:ILE:HD11	1:B:1164:ALA:HB2	1.89	0.53
1:A:1243:LEU:O	1:A:1244:MET:CB	2.56	0.53
1:A:991:HIS:HE1	3:A:1403:HOH:O	1.91	0.53
1:A:1086:ILE:HD12	1:A:1086:ILE:N	2.24	0.53
1:A:1065:ILE:CG2	1:B:956:PRO:HG3	2.39	0.52
1:B:1243:LEU:O	1:B:1244:MET:HB2	2.09	0.52
1:B:1243:LEU:O	1:B:1244:MET:CB	2.58	0.52
1:B:1124:GLN:NE2	1:B:1135:GLN:OE1	2.31	0.52
1:B:1142:ILE:HG21	1:B:1145:THR:CG2	2.38	0.51
1:B:1134:CYS:HB2	1:B:1198:CYS:HB3	1.92	0.51
1:B:1127:LEU:HD12	3:B:1303:HOH:O	2.11	0.51
1:B:1183:ASP:OD2	1:B:1231:GLY:HA3	2.09	0.51
1:B:981:TYR:CE2	1:B:1112:GLN:HG3	2.45	0.51
1:A:1084:SER:HB2	1:A:1085:LEU:CB	2.41	0.50
1:A:1104:ILE:HD11	1:A:1164:ALA:HB2	1.92	0.50
1:B:1119:PHE:CE1	1:B:1153:THR:HG22	2.46	0.50
1:A:980:VAL:CB	1:A:1243:LEU:HD11	2.42	0.50
1:B:1193:ASN:HB2	1:B:1196:LEU:CD2	2.42	0.49
1:B:1010:ILE:HG22	1:B:1013:ILE:HG12	1.94	0.49
1:B:1118:SER:HB3	1:B:1141:THR:OG1	2.13	0.49
1:B:1193:ASN:HB2	1:B:1196:LEU:HD21	1.93	0.49
1:A:989:ALA:HB1	1:A:1155:VAL:HG12	1.95	0.49
1:A:1195:LEU:HD22	1:A:1221:GLN:NE2	2.28	0.49
1:B:1122:ILE:HG12	1:B:1136:MET:HE2	1.93	0.49
1:A:1239:PRO:O	1:A:1243:LEU:HG	2.13	0.48
1:A:1086:ILE:H	1:A:1087:PRO:CA	2.18	0.48
1:B:1166:VAL:C	1:B:1168:ASP:H	2.16	0.48
1:B:989:ALA:CB	1:B:1155:VAL:HG12	2.40	0.48
1:B:963:PHE:HZ	1:B:1002:THR:CG2	2.26	0.48
1:B:1007:GLU:HA	1:B:1187:PHE:O	2.14	0.48
1:B:1122:ILE:HG12	1:B:1136:MET:HE3	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:946:HIS:CG	1:B:947:PRO:HD2	2.49	0.48
1:B:1099:ILE:HD13	1:B:1127:LEU:CD1	2.44	0.47
1:A:1041:SER:HB2	3:A:1422:HOH:O	2.14	0.47
1:A:1195:LEU:CD1	1:A:1221:GLN:HE22	2.24	0.47
1:B:948:LEU:O	1:B:965:SER:OG	2.12	0.47
1:A:1068:TRP:CZ3	1:B:957:LEU:HD11	2.50	0.47
1:A:1195:LEU:HD22	1:A:1221:GLN:HE22	1.80	0.46
1:B:1157:SER:O	1:B:1160:GLN:HB2	2.14	0.46
1:A:1162:ILE:HG13	1:A:1163:ILE:N	2.27	0.46
1:A:1083:GLN:HB3	1:A:1084:SER:OG	2.15	0.46
1:B:1006:THR:HG23	1:B:1007:GLU:CD	2.36	0.46
1:B:978:PHE:C	1:B:979:ILE:HD13	2.37	0.46
1:A:1195:LEU:HB3	1:A:1221:GLN:NE2	2.30	0.46
1:B:1179:PRO:HG2	3:B:1301:HOH:O	2.15	0.46
1:A:963:PHE:HZ	1:A:1002:THR:HG23	1.81	0.46
1:B:1009:GLN:HA	1:B:1185:PHE:O	2.16	0.45
1:A:952:LYS:HD2	1:A:963:PHE:CE2	2.52	0.45
1:A:1064:HIS:CG	1:A:1065:ILE:N	2.83	0.44
1:B:1098:LYS:NZ	1:B:1124:GLN:OE1	2.50	0.44
1:B:1166:VAL:O	1:B:1168:ASP:N	2.47	0.44
1:A:1031:VAL:HG22	1:A:1044:VAL:HG22	1.99	0.44
1:A:1185:PHE:CD2	1:A:1229:VAL:HG22	2.53	0.44
1:A:1084:SER:CB	1:A:1085:LEU:HB2	2.48	0.43
1:B:1137:LYS:HA	1:B:1194:ASP:O	2.18	0.43
1:B:989:ALA:HB2	1:B:1159:PHE:CE2	2.53	0.43
1:A:972:LEU:HB2	1:A:975:LEU:HD12	2.00	0.43
1:B:1010:ILE:HB	1:B:1013:ILE:HD11	2.00	0.43
1:A:1193:ASN:ND2	1:A:1220:ASP:OD1	2.52	0.43
1:A:963:PHE:O	1:A:1030:GLN:HA	2.19	0.43
1:B:946:HIS:NE2	1:B:1149:GLN:OE1	2.51	0.43
1:A:1195:LEU:HD23	1:A:1196:LEU:N	2.34	0.42
1:B:1156:ASP:O	1:B:1160:GLN:HG3	2.20	0.42
1:A:980:VAL:HG12	1:A:1243:LEU:HD11	1.99	0.42
1:A:1083:GLN:HA	1:A:1084:SER:HA	1.69	0.42
1:A:954:GLN:HB2	3:A:1410:HOH:O	2.17	0.42
1:B:946:HIS:ND1	1:B:947:PRO:HD2	2.35	0.41
1:A:1137:LYS:HA	1:A:1194:ASP:O	2.20	0.41
1:B:1007:GLU:HG3	1:B:1188:TYR:CA	2.41	0.41
1:A:1084:SER:HB3	1:A:1225:LEU:HB2	2.02	0.41
1:A:1136:MET:HE2	1:A:1157:SER:HB3	2.01	0.41
1:B:1191:SER:HB2	1:B:1226:VAL:HG21	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1026:VAL:HG13	1:B:1026:VAL:O	2.21	0.41
1:B:1158:CYS:O	1:B:1161:SER:OG	2.21	0.41
1:A:1195:LEU:O	1:A:1196:LEU:HD23	2.20	0.41
1:B:1089:GLU:CD	1:B:1089:GLU:H	2.24	0.41
1:B:1137:LYS:CE	1:B:1195:LEU:HD12	2.46	0.41
1:B:1195:LEU:O	1:B:1196:LEU:HD23	2.20	0.41
1:A:1173:LYS:N	1:A:1173:LYS:HD2	2.36	0.41
1:A:1121:TRP:HB2	1:A:1157:SER:OG	2.20	0.41
1:B:967:PHE:HB3	1:B:1021:ILE:HD11	2.03	0.41
1:A:1084:SER:OG	1:A:1085:LEU:HD22	2.21	0.41
1:A:1095:CYS:SG	1:A:1131:GLU:HG3	2.61	0.41
1:A:1124:GLN:HB2	1:A:1135:GLN:HB3	2.02	0.40
1:B:1017:GLN:HB3	1:B:1070:VAL:O	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1096:SER:OG	1:B:1221:GLN:O[2_454]	2.03	0.17

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	286/308 (93%)	275 (96%)	9 (3%)	2 (1%)	26 38
1	B	273/308 (89%)	265 (97%)	7 (3%)	1 (0%)	39 56
All	All	559/616 (91%)	540 (97%)	16 (3%)	3 (0%)	34 48

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1085	LEU
1	B	1167	LEU
1	A	1086	ILE

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	259/275 (94%)	254 (98%)	5 (2%)	65 83
1	B	251/275 (91%)	248 (99%)	3 (1%)	78 90
All	All	510/550 (93%)	502 (98%)	8 (2%)	70 86

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1067	ASP
1	A	1085	LEU
1	A	1146	THR
1	A	1162	ILE
1	A	1220	ASP
1	B	1080	ASN
1	B	1146	THR
1	B	1220	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	991	HIS
1	A	1221	GLN
1	B	1193	ASN

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 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	7OD	A	1301	-	7,10,10	3.59	1 (14%)	8,12,12	2.52	2 (25%)

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	7OD	A	1301	-	-	0/7/11/11	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1301	7OD	CAD-CAC	9.05	1.54	1.33

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	1301	7OD	CAE-CAD-CAC	-6.39	118.41	126.59
2	A	1301	7OD	CAF-CAE-CAD	2.01	116.23	112.06

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	290/308 (94%)	-0.04	5 (1%) 73 72	42, 68, 132, 205	0
1	B	281/308 (91%)	1.36	72 (25%) 1 1	65, 155, 218, 265	0
All	All	571/616 (92%)	0.65	77 (13%) 4 4	42, 105, 204, 265	0

All (77) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1206	LYS	12.5
1	B	1242	LEU	9.2
1	B	1163	ILE	8.2
1	B	1207	GLN	8.2
1	A	1084	SER	7.1
1	B	1164	ALA	7.0
1	B	1089	GLU	6.6
1	B	1004	ALA	6.2
1	B	1113	ILE	6.2
1	B	1232	PHE	6.1
1	B	1100	ASP	6.0
1	B	1193	ASN	5.9
1	B	1095	CYS	5.2
1	B	1177	PHE	5.2
1	B	1226	VAL	5.1
1	B	1212	PHE	5.1
1	B	1221	GLN	5.0
1	B	1104	ILE	5.0
1	B	1093	ALA	4.8
1	B	1161	SER	4.5
1	B	1222	HIS	4.2
1	B	1110	ASP	4.2
1	B	1225	LEU	4.2
1	B	1199	TYR	4.2

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Mol	Chain	Res	Type	RSRZ
1	A	1083	GLN	4.0
1	B	1167	LEU	4.0
1	B	1119	PHE	3.9
1	B	1218	LEU	3.9
1	B	1091	ILE	3.9
1	A	1086	ILE	3.9
1	B	1143	LEU	3.8
1	B	1194	ASP	3.7
1	B	1195	LEU	3.6
1	B	1190	SER	3.6
1	B	948	LEU	3.6
1	B	1081	ALA	3.5
1	B	1243	LEU	3.5
1	B	1234	GLY	3.5
1	B	1099	ILE	3.4
1	B	981	TYR	3.4
1	A	1085	LEU	3.3
1	B	985	VAL	3.3
1	B	1211	LYS	3.2
1	B	1126	TRP	3.1
1	B	1115	LEU	3.1
1	B	1108	LEU	3.0
1	B	1005	ALA	3.0
1	B	1214	ALA	3.0
1	B	1205	ASP	2.9
1	B	1121	TRP	2.8
1	B	1112	GLN	2.8
1	B	1097	GLN	2.8
1	B	1189	ASN	2.7
1	B	1088	LEU	2.7
1	B	1153	THR	2.6
1	B	1051	LEU	2.6
1	B	1114	HIS	2.6
1	B	1240	LYS	2.5
1	B	1092	GLN	2.5
1	B	1201	CYS	2.4
1	B	1076	LEU	2.4
1	B	1123	GLU	2.4
1	B	986	VAL	2.3
1	B	1144	ASN	2.3
1	B	946	HIS	2.3
1	B	1096	SER	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	1127	LEU	2.2
1	B	1208	SER	2.2
1	B	1147	LYS	2.2
1	B	974	PHE	2.2
1	B	1079	ALA	2.2
1	B	1197	TRP	2.1
1	B	1174	ASN	2.1
1	A	1243	LEU	2.1
1	B	1087	PRO	2.1
1	B	1142	ILE	2.1
1	B	1159	PHE	2.1

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^{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	7OD	A	1301	11/11	0.77	0.20	1.21	97,116,125,125	0

6.5 Other polymers (i)

There are no such residues in this entry.