Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 12:46 PM GMT

PDB ID	:	3S4U
Title	:	Crystal structure of open, unliganded E. coli PhnD H157A
Authors	:	Alicea, I.; Schreiter, E.R.
Deposited on	:	2011-05-20
Resolution	:	3.30 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:

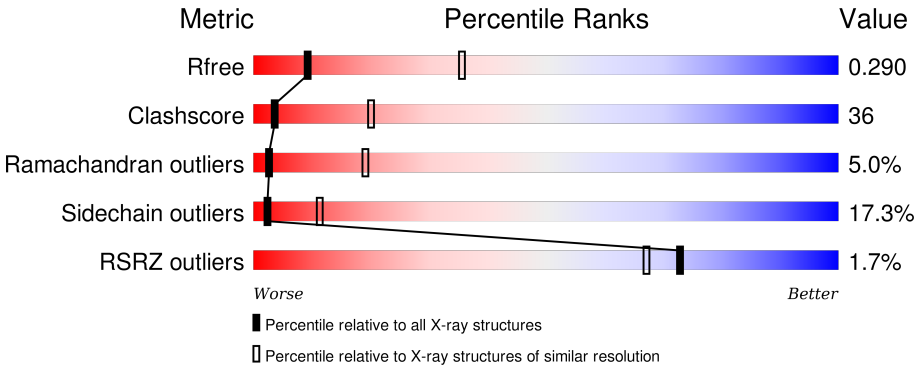
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 3.30 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	2060 (3.40-3.20)
Clashscore	102246	1058 (3.38-3.22)
Ramachandran outliers	100387	1038 (3.38-3.22)
Sidechain outliers	100360	1037 (3.38-3.22)
RSRZ outliers	91569	2070 (3.40-3.20)

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	321	2% 34% 49% 11% • 6%

2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 2360 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 PhnD, subunit of alkylphosphonate ABC transporter.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	302	2360	1500	403	451	6	0	0	0

There are 10 discrepancies between the modelled and reference sequences:

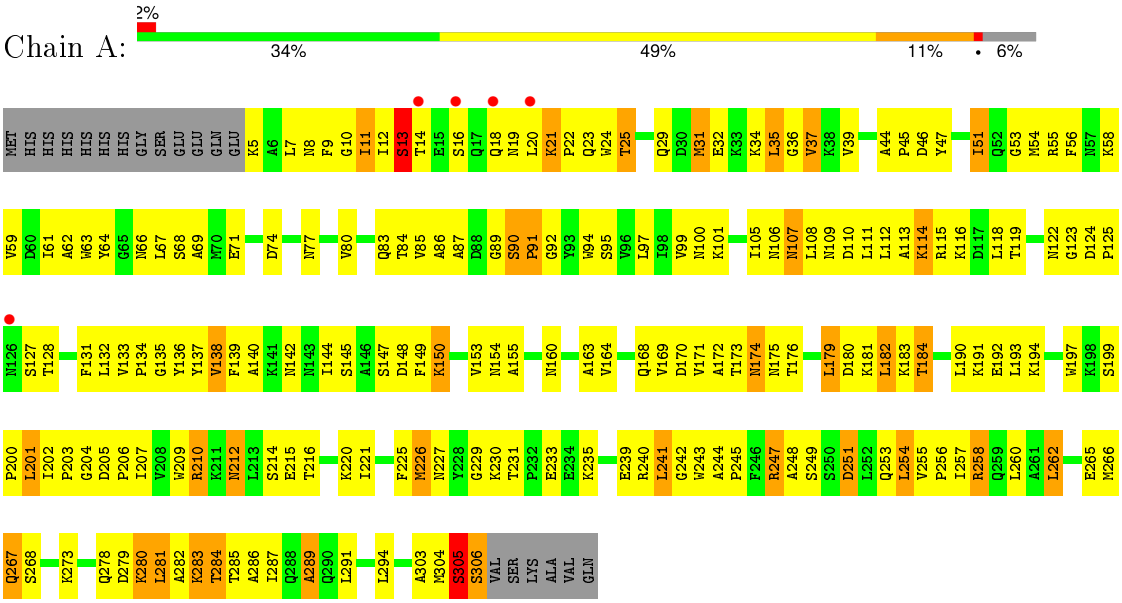
Chain	Residue	Modelled	Actual	Comment	Reference
A	-8	MET	-	EXPRESSION TAG	UNP Q1R3F7
A	-7	HIS	-	EXPRESSION TAG	UNP Q1R3F7
A	-6	HIS	-	EXPRESSION TAG	UNP Q1R3F7
A	-5	HIS	-	EXPRESSION TAG	UNP Q1R3F7
A	-4	HIS	-	EXPRESSION TAG	UNP Q1R3F7
A	-3	HIS	-	EXPRESSION TAG	UNP Q1R3F7
A	-2	HIS	-	EXPRESSION TAG	UNP Q1R3F7
A	-1	GLY	-	EXPRESSION TAG	UNP Q1R3F7
A	0	SER	-	EXPRESSION TAG	UNP Q1R3F7
A	157	ALA	HIS	ENGINEERED MUTATION	UNP Q1R3F7

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: PhnD, subunit of alkylphosphonate ABC transporter

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, α, β, γ	103.73Å 103.73Å 58.42Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.30 19.55 – 3.30	Depositor EDS
% Data completeness (in resolution range)	99.2 (20.00-3.30) 99.8 (19.55-3.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.43 (at 3.29Å)	Xtriage
Refinement program	REFMAC 5.6.0111	Depositor
R, R_{free}	0.194 , 0.289 0.197 , 0.290	Depositor DCC
R_{free} test set	237 reflections (4.87%)	DCC
Wilson B-factor (Å^2)	58.0	Xtriage
Anisotropy	0.304	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 41.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.45$, $< L^2 > = 0.28$	Xtriage
Outliers	0 of 5104 reflections	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	2360	wwPDB-VP
Average B, all atoms (Å^2)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.45% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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	0.90	1/2406 (0.0%)	0.91	3/3259 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	192	GLU	CG-CD	5.79	1.60	1.51

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	210	ARG	NE-CZ-NH1	6.37	123.49	120.30
1	A	262	LEU	CA-CB-CG	5.41	127.73	115.30
1	A	304	MET	N-CA-C	-5.27	96.76	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	13	SER	Peptide
1	A	147	SER	Peptide

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	2360	0	2367	169	1
All	All	2360	0	2367	169	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 36.

All (169) 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:226:MET:HE1	1:A:248:ALA:HA	1.23	1.09
1:A:226:MET:HE1	1:A:248:ALA:CA	1.97	0.94
1:A:85:VAL:HG12	1:A:244:ALA:HB3	1.51	0.93
1:A:118:LEU:HD22	1:A:170:ASP:CG	1.93	0.89
1:A:122:ASN:O	1:A:154:ASN:HA	1.73	0.87
1:A:134:PRO:O	1:A:138:VAL:HG23	1.75	0.85
1:A:23:GLN:O	1:A:241:LEU:HD11	1.81	0.80
1:A:128:THR:O	1:A:133:VAL:HG23	1.81	0.80
1:A:47:TYR:O	1:A:51:ILE:HG23	1.83	0.78
1:A:255:VAL:HG13	1:A:294:LEU:HD22	1.64	0.77
1:A:206:PRO:C	1:A:207:ILE:HD13	2.06	0.76
1:A:5:LYS:O	1:A:37:VAL:CG2	2.34	0.76
1:A:97:LEU:HD21	1:A:173:THR:HG22	1.68	0.76
1:A:262:LEU:HD12	1:A:291:LEU:HD21	1.68	0.75
1:A:101:LYS:NZ	1:A:191:LYS:O	2.19	0.75
1:A:12:ILE:HA	1:A:44:ALA:HB3	1.69	0.73
1:A:66:ASN:HA	1:A:69:ALA:HB3	1.70	0.73
1:A:241:LEU:N	1:A:241:LEU:HD13	2.03	0.73
1:A:25:THR:HG22	1:A:29:GLN:HE21	1.54	0.73
1:A:262:LEU:HD12	1:A:291:LEU:CD2	2.19	0.73
1:A:201:LEU:HD12	1:A:202:ILE:N	2.05	0.70
1:A:97:LEU:CD2	1:A:173:THR:HG22	2.21	0.70
1:A:5:LYS:O	1:A:37:VAL:HG23	1.92	0.70
1:A:251:ASP:O	1:A:254:LEU:HD12	1.93	0.68
1:A:118:LEU:HD13	1:A:170:ASP:HB3	1.75	0.68
1:A:199:SER:OG	1:A:200:PRO:O	2.11	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:201:LEU:HD12	1:A:202:ILE:C	2.16	0.66
1:A:11:ILE:HD12	1:A:11:ILE:N	2.10	0.65
1:A:19:ASN:O	1:A:23:GLN:NE2	2.29	0.65
1:A:254:LEU:HD12	1:A:254:LEU:H	1.61	0.65
1:A:107:ASN:HB3	1:A:109:ASN:H	1.62	0.65
1:A:10:GLY:C	1:A:11:ILE:HD12	2.17	0.64
1:A:226:MET:HE3	1:A:248:ALA:HB2	1.80	0.64
1:A:97:LEU:HD23	1:A:173:THR:HA	1.80	0.63
1:A:113:ALA:O	1:A:115:ARG:N	2.29	0.62
1:A:226:MET:HE1	1:A:248:ALA:CB	2.29	0.62
1:A:226:MET:CE	1:A:248:ALA:CB	2.78	0.62
1:A:283:LYS:O	1:A:287:ILE:HD12	2.01	0.61
1:A:281:LEU:O	1:A:285:THR:HG23	2.01	0.61
1:A:286:ALA:O	1:A:289:ALA:HB3	1.99	0.61
1:A:207:ILE:N	1:A:207:ILE:HD13	2.14	0.61
1:A:133:VAL:HG11	1:A:202:ILE:HD11	1.83	0.60
1:A:99:VAL:HG23	1:A:100:ASN:O	2.02	0.60
1:A:149:PHE:O	1:A:150:LYS:C	2.40	0.59
1:A:35:LEU:HB3	1:A:37:VAL:HG13	1.83	0.59
1:A:125:PRO:HA	1:A:131:PHE:CD1	2.37	0.59
1:A:77:ASN:O	1:A:210:ARG:HA	2.02	0.59
1:A:254:LEU:HD12	1:A:254:LEU:N	2.17	0.59
1:A:5:LYS:O	1:A:37:VAL:HG21	2.03	0.58
1:A:111:LEU:HD13	1:A:111:LEU:C	2.24	0.58
1:A:66:ASN:CA	1:A:69:ALA:HB3	2.34	0.58
1:A:31:MET:HG3	1:A:35:LEU:CD2	2.34	0.58
1:A:266:MET:HE1	1:A:287:ILE:HB	1.86	0.58
1:A:239:GLU:O	1:A:242:GLY:N	2.37	0.58
1:A:171:VAL:HG22	1:A:172:ALA:N	2.19	0.58
1:A:32:GLU:O	1:A:36:GLY:HA2	2.04	0.57
1:A:135:GLY:HA2	1:A:139:PHE:HD2	1.69	0.57
1:A:105:ILE:HG23	1:A:110:ASP:HB2	1.86	0.57
1:A:90:SER:HB2	1:A:91:PRO:HB3	1.85	0.57
1:A:305:SER:N	1:A:306:SER:HB2	2.19	0.57
1:A:280:LYS:O	1:A:282:ALA:N	2.38	0.57
1:A:180:ASP:O	1:A:181:LYS:C	2.43	0.57
1:A:226:MET:CE	1:A:248:ALA:HB2	2.35	0.56
1:A:280:LYS:O	1:A:281:LEU:C	2.42	0.56
1:A:20:LEU:HD22	1:A:24:TRP:CD1	2.40	0.56
1:A:115:ARG:HG3	1:A:116:LYS:N	2.21	0.56
1:A:227:ASN:OD1	1:A:230:LYS:NZ	2.36	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:229:GLY:O	1:A:235:LYS:HG3	2.06	0.55
1:A:108:LEU:HD22	1:A:197:TRP:CD1	2.42	0.55
1:A:266:MET:SD	1:A:284:THR:HG22	2.46	0.55
1:A:32:GLU:HG3	1:A:39:VAL:HB	1.87	0.55
1:A:176:THR:HA	1:A:179:LEU:HB2	1.89	0.55
1:A:37:VAL:O	1:A:39:VAL:HG23	2.06	0.55
1:A:74:ASP:OD1	1:A:258:ARG:NH2	2.38	0.55
1:A:171:VAL:CG2	1:A:172:ALA:N	2.70	0.54
1:A:111:LEU:HD21	1:A:171:VAL:HG11	1.89	0.53
1:A:97:LEU:HD11	1:A:138:VAL:HG11	1.90	0.53
1:A:127:SER:O	1:A:131:PHE:HB3	2.09	0.53
1:A:119:THR:O	1:A:169:VAL:HG22	2.09	0.53
1:A:134:PRO:C	1:A:138:VAL:HG23	2.28	0.52
1:A:201:LEU:C	1:A:201:LEU:HD12	2.31	0.51
1:A:23:GLN:HB3	1:A:241:LEU:HD12	1.93	0.51
1:A:221:ILE:HG23	1:A:225:PHE:CE1	2.46	0.51
1:A:32:GLU:O	1:A:36:GLY:CA	2.59	0.50
1:A:35:LEU:HD12	1:A:220:LYS:HB3	1.93	0.50
1:A:262:LEU:CD1	1:A:291:LEU:CD2	2.88	0.50
1:A:118:LEU:HD22	1:A:170:ASP:OD1	2.12	0.50
1:A:39:VAL:O	1:A:39:VAL:HG12	2.12	0.50
1:A:95:SER:HB2	1:A:199:SER:HB3	1.94	0.50
1:A:231:THR:HG22	1:A:233:GLU:H	1.77	0.49
1:A:9:PHE:CD2	1:A:61:ILE:HG23	2.48	0.49
1:A:206:PRO:O	1:A:207:ILE:HD13	2.13	0.49
1:A:105:ILE:HG23	1:A:110:ASP:CB	2.42	0.49
1:A:133:VAL:HG21	1:A:202:ILE:CD1	2.43	0.48
1:A:266:MET:HB2	1:A:287:ILE:HG21	1.94	0.48
1:A:241:LEU:CD1	1:A:241:LEU:N	2.73	0.48
1:A:182:LEU:HB2	1:A:190:LEU:HD22	1.96	0.48
1:A:8:ASN:O	1:A:61:ILE:HG22	2.14	0.47
1:A:163:ALA:HA	1:A:168:GLN:OE1	2.15	0.47
1:A:251:ASP:O	1:A:254:LEU:CD1	2.62	0.47
1:A:226:MET:HE3	1:A:248:ALA:CB	2.43	0.47
1:A:107:ASN:HB2	1:A:110:ASP:OD1	2.14	0.47
1:A:83:GLN:NE2	1:A:91:PRO:O	2.48	0.47
1:A:280:LYS:O	1:A:283:LYS:N	2.48	0.47
1:A:21:LYS:HB2	1:A:22:PRO:HD3	1.96	0.47
1:A:46:ASP:N	1:A:46:ASP:OD1	2.48	0.46
1:A:182:LEU:O	1:A:184:THR:N	2.48	0.46
1:A:214:SER:O	1:A:215:GLU:C	2.54	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:175:ASN:O	1:A:179:LEU:HD23	2.16	0.46
1:A:133:VAL:N	1:A:134:PRO:HD2	2.30	0.46
1:A:63:TRP:C	1:A:63:TRP:CD1	2.87	0.46
1:A:99:VAL:HG22	1:A:194:LYS:HG2	1.98	0.46
1:A:134:PRO:CA	1:A:138:VAL:HG23	2.45	0.46
1:A:97:LEU:HD21	1:A:173:THR:CG2	2.43	0.46
1:A:244:ALA:HB1	1:A:245:PRO:CD	2.46	0.45
1:A:83:GLN:HB3	1:A:247:ARG:HB3	1.98	0.45
1:A:112:LEU:CD2	1:A:144:ILE:HD13	2.46	0.45
1:A:7:LEU:HD12	1:A:37:VAL:HG22	1.98	0.45
1:A:31:MET:HE3	1:A:39:VAL:HG11	1.98	0.45
1:A:12:ILE:O	1:A:13:SER:CB	2.63	0.45
1:A:61:ILE:O	1:A:62:ALA:HB2	2.17	0.45
1:A:190:LEU:O	1:A:190:LEU:HD12	2.17	0.45
1:A:25:THR:O	1:A:29:GLN:HG2	2.16	0.45
1:A:155:ALA:HB3	1:A:160:ASN:OD1	2.17	0.45
1:A:54:MET:O	1:A:56:PHE:N	2.49	0.44
1:A:71:GLU:HA	1:A:74:ASP:HB2	1.99	0.44
1:A:87:ALA:HB2	1:A:243:TRP:N	2.32	0.44
1:A:90:SER:CB	1:A:91:PRO:HB3	2.48	0.44
1:A:115:ARG:HE	1:A:148:ASP:HB3	1.81	0.44
1:A:164:VAL:HG22	1:A:169:VAL:CG1	2.48	0.44
1:A:267:GLN:HA	1:A:267:GLN:HE21	1.82	0.44
1:A:281:LEU:HD12	1:A:285:THR:HG21	2.00	0.44
1:A:107:ASN:HB3	1:A:109:ASN:N	2.31	0.43
1:A:94:TRP:O	1:A:176:THR:HG23	2.19	0.43
1:A:255:VAL:N	1:A:256:PRO:CD	2.82	0.43
1:A:10:GLY:CA	1:A:11:ILE:HD12	2.48	0.43
1:A:240:ARG:C	1:A:241:LEU:HD13	2.37	0.43
1:A:92:GLY:HA3	1:A:203:PRO:HA	2.00	0.43
1:A:94:TRP:NE1	1:A:201:LEU:HB2	2.34	0.43
1:A:97:LEU:CD2	1:A:173:THR:CG2	2.95	0.43
1:A:133:VAL:HG21	1:A:202:ILE:HD11	2.00	0.43
1:A:266:MET:HE1	1:A:284:THR:O	2.18	0.43
1:A:305:SER:HB2	1:A:306:SER:HA	2.01	0.43
1:A:84:THR:HG23	1:A:205:ASP:O	2.18	0.42
1:A:204:GLY:O	1:A:253:GLN:NE2	2.52	0.42
1:A:86:ALA:O	1:A:89:GLY:N	2.48	0.42
1:A:182:LEU:C	1:A:184:THR:N	2.71	0.42
1:A:32:GLU:O	1:A:36:GLY:N	2.53	0.42
1:A:63:TRP:HE1	1:A:205:ASP:CB	2.33	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:281:LEU:O	1:A:285:THR:CG2	2.66	0.42
1:A:212:ASN:OD1	1:A:212:ASN:C	2.58	0.42
1:A:138:VAL:O	1:A:142:ASN:ND2	2.47	0.42
1:A:305:SER:CB	1:A:306:SER:HA	2.49	0.42
1:A:101:LYS:CE	1:A:191:LYS:O	2.68	0.42
1:A:63:TRP:CG	1:A:63:TRP:O	2.73	0.42
1:A:56:PHE:C	1:A:58:LYS:H	2.23	0.42
1:A:97:LEU:HD23	1:A:173:THR:HG22	1.99	0.41
1:A:80:VAL:HG13	1:A:206:PRO:HB2	2.01	0.41
1:A:109:ASN:O	1:A:110:ASP:C	2.59	0.41
1:A:137:TYR:HB2	1:A:197:TRP:HZ2	1.84	0.41
1:A:53:GLY:C	1:A:59:VAL:HG12	2.40	0.41
1:A:173:THR:O	1:A:174:ASN:ND2	2.36	0.41
1:A:94:TRP:CZ2	1:A:201:LEU:HD22	2.56	0.41
1:A:11:ILE:CD1	1:A:11:ILE:N	2.79	0.41
1:A:193:LEU:HD23	1:A:193:LEU:HA	1.92	0.41
1:A:64:TYR:O	1:A:205:ASP:HB3	2.21	0.40
1:A:136:TYR:HA	1:A:140:ALA:HB3	2.03	0.40
1:A:123:GLY:O	1:A:124:ASP:C	2.59	0.40
1:A:122:ASN:O	1:A:154:ASN:CA	2.58	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:247:ARG:NH1	1:A:303:ALA:O[7_555]	2.11	0.09

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	300/321 (94%)	234 (78%)	51 (17%)	15 (5%)	3 19

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	13	SER
1	A	16	SER
1	A	21	LYS
1	A	150	LYS
1	A	281	LEU
1	A	114	LYS
1	A	305	SER
1	A	55	ARG
1	A	91	PRO
1	A	183	LYS
1	A	212	ASN
1	A	280	LYS
1	A	289	ALA
1	A	90	SER
1	A	45	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	255/272 (94%)	211 (83%)	44 (17%)	2 11

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	ILE
1	A	14	THR
1	A	18	GLN
1	A	25	THR
1	A	31	MET
1	A	34	LYS
1	A	35	LEU
1	A	37	VAL
1	A	51	ILE
1	A	67	LEU

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Mol	Chain	Res	Type
1	A	68	SER
1	A	106	ASN
1	A	107	ASN
1	A	114	LYS
1	A	132	LEU
1	A	138	VAL
1	A	145	SER
1	A	153	VAL
1	A	174	ASN
1	A	179	LEU
1	A	182	LEU
1	A	184	THR
1	A	201	LEU
1	A	209	TRP
1	A	216	THR
1	A	226	MET
1	A	241	LEU
1	A	247	ARG
1	A	249	SER
1	A	251	ASP
1	A	254	LEU
1	A	257	ILE
1	A	258	ARG
1	A	260	LEU
1	A	265	GLU
1	A	267	GLN
1	A	268	SER
1	A	273	LYS
1	A	278	GLN
1	A	279	ASP
1	A	283	LYS
1	A	284	THR
1	A	305	SER
1	A	306	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	19	ASN
1	A	29	GLN
1	A	106	ASN
1	A	107	ASN

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Mol	Chain	Res	Type
1	A	122	ASN
1	A	178	ASN
1	A	267	GLN
1	A	299	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

There are no carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

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	302/321 (94%)	-0.37	5 (1%)	73 67	25, 49, 78, 128	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	14	THR	3.3
1	A	16	SER	3.3
1	A	20	LEU	2.8
1	A	126	ASN	2.2
1	A	18	GLN	2.1

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

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

6.3 Carbohydrates

There are no carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.