



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:26 AM GMT

PDB ID : 3IFR
Title : The crystal structure of xylulose kinase from Rhodospirillum rubrum
Authors : Zhang, Z.; Burley, S.K.; Swaminathan, S.; New York SGX Research Center
for Structural Genomics (NYSGXRC)
Deposited on : 2009-07-24
Resolution : 2.30 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 ⓘ symbol.

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

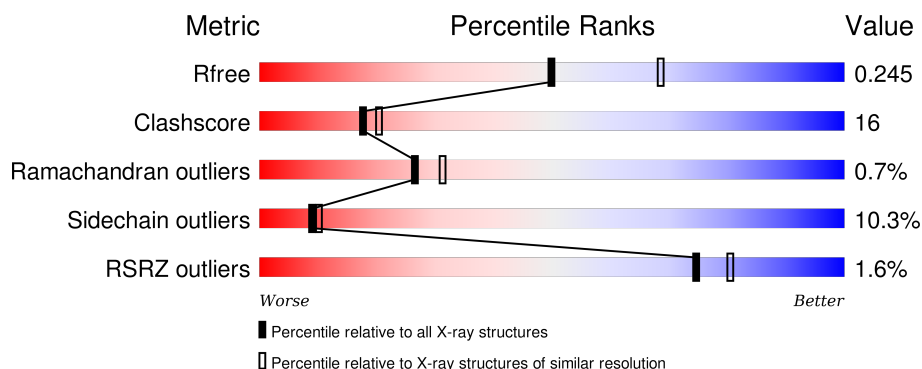
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

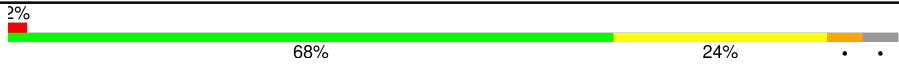

The reported resolution of this entry is 2.30 Å.

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	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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 $\leq 5\%$. 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	508	
1	B	508	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	PO4	A	509	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7532 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 Carbohydrate kinase, FGGY.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	488	Total	C	N	O	S	Se	0	0	0
			3624	2301	638	675	5	5			
1	B	480	Total	C	N	O	S	Se	0	0	0
			3564	2263	629	662	5	5			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	-	EXPRESSION TAG	UNP Q2RUX7
A	2	SER	-	EXPRESSION TAG	UNP Q2RUX7
A	3	LEU	-	EXPRESSION TAG	UNP Q2RUX7
A	501	GLU	-	EXPRESSION TAG	UNP Q2RUX7
A	502	GLY	-	EXPRESSION TAG	UNP Q2RUX7
A	503	HIS	-	EXPRESSION TAG	UNP Q2RUX7
A	504	HIS	-	EXPRESSION TAG	UNP Q2RUX7
A	505	HIS	-	EXPRESSION TAG	UNP Q2RUX7
A	506	HIS	-	EXPRESSION TAG	UNP Q2RUX7
A	507	HIS	-	EXPRESSION TAG	UNP Q2RUX7
A	508	HIS	-	EXPRESSION TAG	UNP Q2RUX7
B	1	MSE	-	EXPRESSION TAG	UNP Q2RUX7
B	2	SER	-	EXPRESSION TAG	UNP Q2RUX7
B	3	LEU	-	EXPRESSION TAG	UNP Q2RUX7
B	501	GLU	-	EXPRESSION TAG	UNP Q2RUX7
B	502	GLY	-	EXPRESSION TAG	UNP Q2RUX7
B	503	HIS	-	EXPRESSION TAG	UNP Q2RUX7
B	504	HIS	-	EXPRESSION TAG	UNP Q2RUX7
B	505	HIS	-	EXPRESSION TAG	UNP Q2RUX7
B	506	HIS	-	EXPRESSION TAG	UNP Q2RUX7
B	507	HIS	-	EXPRESSION TAG	UNP Q2RUX7
B	508	HIS	-	EXPRESSION TAG	UNP Q2RUX7

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		

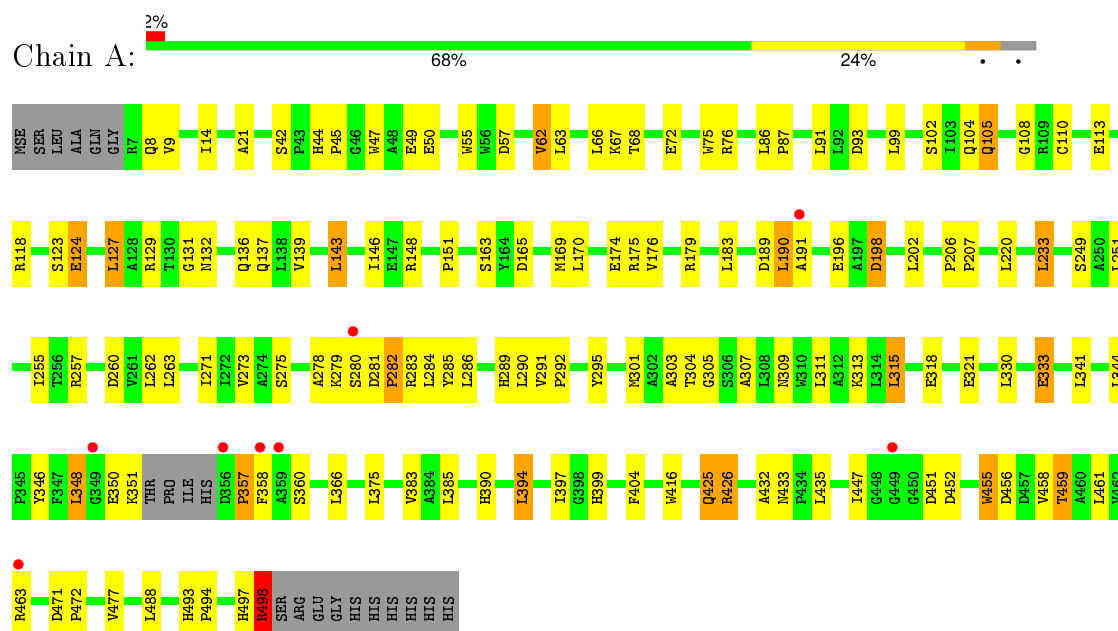
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	170	Total	O	0	0
			170	170		
3	B	169	Total	O	0	0
			169	169		

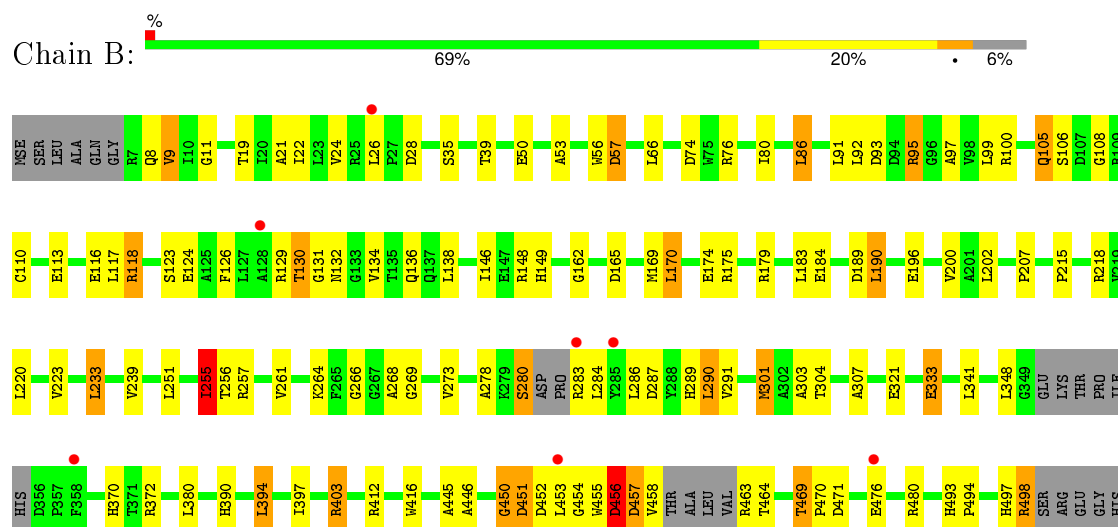
3 Residue-property plots

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: Carbohydrate kinase, FGGY



• Molecule 1: Carbohydrate kinase, FGGY



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4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	129.09 Å 149.81 Å 54.94 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	64.55 – 2.30 64.55 – 2.30	Depositor EDS
% Data completeness (in resolution range)	95.8 (64.55-2.30) 98.4 (64.55-2.30)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.05 (at 2.29 Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.206 , 0.260 0.202 , 0.245	Depositor DCC
R_{free} test set	4548 reflections (5.08%)	DCC
Wilson B-factor (Å ²)	27.7	Xtriage
Anisotropy	0.678	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 42.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 47562 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7532	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.59% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PO4

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	0.75	1/3707 (0.0%)	0.83	2/5065 (0.0%)
1	B	0.78	0/3644	0.82	0/4975
All	All	0.77	1/7351 (0.0%)	0.83	2/10040 (0.0%)

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	1
1	B	0	1
All	All	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	477	VAL	CB-CG2	-5.10	1.42	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	202	LEU	CA-CB-CG	5.23	127.33	115.30
1	A	498	ARG	NE-CZ-NH1	-5.13	117.73	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	455	TRP	Peptide
1	B	455	TRP	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	3624	0	3580	119	0
1	B	3564	0	3516	108	0
2	A	5	0	0	0	0
3	A	170	0	0	6	0
3	B	169	0	0	12	0
All	All	7532	0	7096	227	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 16.

All (227) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:301:MSE:HE2	1:B:390:HIS:CE1	1.80	1.16
1:B:95:ARG:O	1:B:95:ARG:HD2	1.61	1.01
1:B:261:VAL:HG22	1:B:273:VAL:HG22	1.40	1.00
1:B:458:VAL:O	1:B:458:VAL:HG12	1.59	0.98
1:A:426:ARG:HG3	1:A:426:ARG:HH11	1.30	0.97
1:B:190:LEU:HD21	1:B:286:LEU:HD23	1.44	0.96
1:B:26:LEU:HD21	1:B:446:ALA:HB1	1.51	0.93
1:B:130:THR:HG22	1:B:132:ASN:H	1.32	0.92
1:A:333:GLU:O	1:A:333:GLU:HG3	1.71	0.87
1:B:118:ARG:HD2	3:B:613:HOH:O	1.75	0.85
1:B:301:MSE:CE	1:B:390:HIS:CE1	2.59	0.84
1:A:131:GLY:HA2	1:A:280:SER:HB3	1.58	0.83
1:A:350:GLU:HG2	1:A:351:LYS:N	1.95	0.81
1:B:255:ILE:O	1:B:255:ILE:HG22	1.81	0.81
1:B:301:MSE:HE2	1:B:390:HIS:HE1	1.46	0.80
1:B:458:VAL:CG1	1:B:458:VAL:O	2.30	0.79
1:B:149:HIS:HD2	3:B:633:HOH:O	1.64	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:21:ALA:HB3	1:A:62:VAL:HG13	1.66	0.77
1:B:264:LYS:HE2	3:B:549:HOH:O	1.85	0.77
1:A:289:HIS:HD2	1:A:290:LEU:H	1.32	0.77
1:B:257:ARG:HD2	1:B:403:ARG:NH1	2.00	0.76
1:A:304:THR:HG22	1:A:305:GLY:N	1.99	0.76
1:B:189:ASP:HB2	1:B:196:GLU:HG3	1.69	0.75
1:A:341:LEU:HD21	1:A:366:LEU:HD23	1.73	0.70
1:B:130:THR:HG22	1:B:132:ASN:N	2.06	0.70
1:A:21:ALA:CB	1:A:62:VAL:HG13	2.22	0.69
1:B:179:ARG:HD2	1:B:215:PRO:HA	1.74	0.69
1:B:110:CYS:HB2	1:B:113:GLU:HG2	1.73	0.69
1:B:179:ARG:HD2	1:B:215:PRO:CA	2.24	0.68
1:B:257:ARG:HD2	1:B:403:ARG:HH12	1.59	0.67
1:B:289:HIS:HD2	1:B:290:LEU:H	1.42	0.67
1:A:275:SER:OG	1:A:399:HIS:HD2	1.78	0.67
1:A:289:HIS:CD2	1:A:290:LEU:H	2.12	0.67
1:B:26:LEU:HD21	1:B:446:ALA:CB	2.22	0.66
1:B:93:ASP:HB3	1:B:99:LEU:HD21	1.77	0.66
1:A:190:LEU:HD21	1:A:286:LEU:HD23	1.77	0.66
1:B:39:THR:O	1:B:50:GLU:HG2	1.96	0.66
1:A:91:LEU:HD11	1:A:146:ILE:HD13	1.75	0.66
1:A:284:LEU:HD13	1:A:394:LEU:HD13	1.78	0.65
1:B:450:GLY:O	1:B:452:ASP:N	2.29	0.65
1:B:471:ASP:C	1:B:471:ASP:OD2	2.32	0.65
1:A:426:ARG:HH11	1:A:426:ARG:CG	2.09	0.64
1:A:350:GLU:HG2	1:A:351:LYS:H	1.60	0.64
1:B:110:CYS:H	1:B:136:GLN:NE2	1.95	0.64
1:A:129:ARG:NH2	1:A:198:ASP:OD2	2.30	0.64
1:A:255:ILE:HD13	1:A:262:LEU:HB2	1.78	0.64
1:B:26:LEU:HD23	1:B:453:LEU:CD1	2.28	0.64
1:B:220:LEU:HD12	1:B:220:LEU:O	1.98	0.64
1:A:425:GLN:HE22	1:A:472:PRO:HA	1.63	0.63
1:A:110:CYS:HB2	1:A:113:GLU:HG2	1.79	0.63
1:A:289:HIS:HD2	1:A:290:LEU:N	1.97	0.61
1:B:333:GLU:O	1:B:333:GLU:HG3	1.99	0.61
1:A:357:PRO:HB2	1:A:497:HIS:CE1	2.37	0.60
1:B:498:ARG:O	1:B:498:ARG:HG2	1.99	0.60
1:A:426:ARG:HG3	1:A:426:ARG:NH1	2.08	0.60
1:B:289:HIS:CD2	1:B:290:LEU:H	2.17	0.59
1:A:350:GLU:CG	1:A:351:LYS:N	2.64	0.59
1:A:350:GLU:CG	1:A:351:LYS:H	2.16	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:148:ARG:HG2	1:B:149:HIS:ND1	2.18	0.59
1:A:255:ILE:CD1	1:A:262:LEU:HB2	2.33	0.59
1:B:289:HIS:HD2	1:B:290:LEU:N	2.02	0.58
1:B:200:VAL:HG11	1:B:207:PRO:HA	1.85	0.58
1:A:289:HIS:CD2	1:A:290:LEU:N	2.71	0.58
1:B:100:ARG:HG3	1:B:100:ARG:HH21	1.68	0.58
1:B:218:ARG:HB3	3:B:599:HOH:O	2.03	0.58
1:B:304:THR:HG22	1:B:304:THR:O	2.04	0.58
1:B:179:ARG:CD	1:B:215:PRO:HA	2.34	0.58
1:B:450:GLY:C	1:B:452:ASP:H	2.07	0.58
1:A:86:LEU:O	1:A:105:GLN:HG2	2.04	0.58
1:B:124:GLU:N	1:B:124:GLU:OE1	2.37	0.57
1:A:493:HIS:HB3	1:A:494:PRO:HD3	1.86	0.57
1:A:281:ASP:C	1:A:283:ARG:H	2.08	0.57
1:B:8:GLN:HB2	1:B:76:ARG:O	2.04	0.57
1:B:454:GLY:N	1:B:457:ASP:OD2	2.24	0.57
1:A:165:ASP:O	1:A:169:MSE:HG3	2.05	0.56
1:A:284:LEU:CD1	1:A:394:LEU:HD13	2.35	0.56
1:A:311:LEU:HG	1:A:315:LEU:HD22	1.86	0.56
1:A:21:ALA:HB2	1:A:62:VAL:CG1	2.36	0.56
1:A:278:ALA:HB2	1:A:295:TYR:CZ	2.41	0.55
1:A:251:LEU:O	1:A:458:VAL:HG11	2.06	0.55
1:B:301:MSE:CE	1:B:390:HIS:ND1	2.69	0.55
1:B:26:LEU:HD23	1:B:453:LEU:HD12	1.87	0.55
1:A:91:LEU:HD11	1:A:146:ILE:CD1	2.37	0.55
1:A:309:ASN:O	1:A:313:LYS:HG3	2.06	0.55
1:A:189:ASP:HB2	1:A:196:GLU:HG3	1.88	0.55
1:A:127:LEU:HD23	1:A:132:ASN:C	2.27	0.55
1:B:289:HIS:CD2	1:B:290:LEU:N	2.75	0.55
1:A:304:THR:HG22	1:A:305:GLY:H	1.69	0.55
1:A:108:GLY:C	1:A:136:GLN:HE22	2.10	0.54
1:B:126:PHE:O	1:B:130:THR:HB	2.08	0.54
1:A:278:ALA:HB2	1:A:295:TYR:CE1	2.43	0.54
1:A:190:LEU:HD21	1:A:286:LEU:CD2	2.38	0.54
1:B:53:ALA:O	1:B:57:ASP:HB2	2.07	0.54
1:A:447:ILE:HD13	1:A:455:TRP:CD2	2.43	0.54
1:B:91:LEU:HD11	1:B:146:ILE:HD13	1.90	0.54
1:B:450:GLY:C	1:B:452:ASP:N	2.61	0.53
1:B:321:GLU:HB2	3:B:616:HOH:O	2.07	0.53
1:B:19:THR:O	1:B:35:SER:HA	2.09	0.53
1:B:450:GLY:O	1:B:451:ASP:C	2.47	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:22:ILE:HD12	1:B:24:VAL:CG2	2.40	0.52
1:A:284:LEU:HD12	1:A:397:ILE:HD11	1.91	0.52
1:B:86:LEU:O	1:B:105:GLN:HG2	2.10	0.52
1:B:266:GLY:O	1:B:304:THR:O	2.28	0.52
1:A:113:GLU:H	1:A:113:GLU:CD	2.12	0.52
1:A:281:ASP:OD2	1:A:283:ARG:NE	2.42	0.51
1:A:471:ASP:OD1	1:A:471:ASP:C	2.48	0.51
1:B:190:LEU:HD21	1:B:286:LEU:CD2	2.30	0.51
1:A:21:ALA:CB	1:A:62:VAL:CG1	2.87	0.51
1:A:432:ALA:HB3	1:A:463:ARG:HB3	1.92	0.51
1:A:303:ALA:HB1	1:A:307:ALA:HB2	1.93	0.51
1:B:169:MSE:HG2	1:B:174:GLU:O	2.11	0.51
1:B:493:HIS:HB3	1:B:494:PRO:HD3	1.93	0.51
1:B:56:TRP:CD2	1:B:170:LEU:HD23	2.46	0.51
1:A:189:ASP:OD2	1:A:191:ALA:HB3	2.11	0.50
1:A:110:CYS:H	1:A:136:GLN:NE2	2.10	0.50
1:B:108:GLY:C	1:B:136:GLN:HE22	2.15	0.50
1:A:281:ASP:O	1:A:283:ARG:N	2.43	0.50
1:A:357:PRO:HB2	1:A:497:HIS:HE1	1.75	0.50
1:B:341:LEU:HD21	1:B:370:HIS:CE1	2.47	0.50
1:B:284:LEU:HD13	1:B:394:LEU:HD13	1.94	0.49
1:A:67:LYS:HG3	1:A:75:TRP:CE2	2.47	0.49
1:B:284:LEU:CD1	1:B:394:LEU:HD13	2.42	0.49
1:A:175:ARG:HD3	3:A:597:HOH:O	2.12	0.49
1:A:447:ILE:HD11	1:A:455:TRP:HA	1.94	0.49
1:A:220:LEU:O	1:A:220:LEU:HD12	2.13	0.49
1:A:383:VAL:HG11	1:A:416:TRP:HZ2	1.78	0.49
1:A:14:ILE:HG22	1:A:14:ILE:O	2.13	0.49
1:A:124:GLU:CA	1:A:124:GLU:OE2	2.60	0.48
1:A:55:TRP:CZ3	1:A:163:SER:HA	2.48	0.48
1:A:42:SER:CB	1:A:47:TRP:O	2.61	0.48
1:A:139:VAL:HG12	1:A:143:LEU:HD22	1.93	0.48
1:B:456:ASP:N	1:B:456:ASP:OD2	2.42	0.48
1:B:268:ALA:O	1:B:304:THR:OG1	2.29	0.48
1:A:273:VAL:HG21	1:A:394:LEU:HG	1.97	0.47
1:A:257:ARG:CD	3:A:601:HOH:O	2.63	0.47
1:B:175:ARG:HD3	3:B:660:HOH:O	2.14	0.47
1:A:425:GLN:NE2	1:A:472:PRO:HA	2.28	0.47
1:B:124:GLU:N	1:B:124:GLU:CD	2.68	0.47
1:B:100:ARG:HG3	1:B:100:ARG:NH2	2.28	0.47
1:B:303:ALA:HB1	1:B:307:ALA:HB2	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:93:ASP:HB3	1:A:99:LEU:HD21	1.96	0.47
1:A:344:LEU:HD11	1:A:488:LEU:HG	1.96	0.47
1:B:268:ALA:O	1:B:301:MSE:O	2.33	0.47
1:A:42:SER:HB3	1:A:47:TRP:O	2.15	0.46
1:B:278:ALA:HB1	1:B:286:LEU:HD21	1.97	0.46
1:A:44:HIS:HB3	1:A:45:PRO:HD2	1.97	0.46
1:B:80:ILE:CG2	1:B:239:VAL:HG22	2.46	0.46
1:A:275:SER:OG	1:A:399:HIS:CD2	2.65	0.46
1:B:268:ALA:O	1:B:304:THR:HB	2.16	0.46
1:A:383:VAL:HG11	1:A:416:TRP:CZ2	2.51	0.46
1:A:124:GLU:OE2	1:A:124:GLU:HA	2.16	0.46
1:A:113:GLU:CD	1:A:113:GLU:N	2.69	0.46
1:A:390:HIS:O	1:A:394:LEU:HD22	2.15	0.45
1:A:281:ASP:C	1:A:283:ARG:N	2.69	0.45
1:B:9:VAL:HG22	1:B:445:ALA:HB3	1.98	0.45
1:A:426:ARG:NH1	1:A:426:ARG:CG	2.72	0.45
1:A:207:PRO:HD2	3:A:582:HOH:O	2.16	0.45
1:A:86:LEU:HB2	1:A:87:PRO:HA	1.97	0.45
1:A:459:THR:HG23	1:A:459:THR:O	2.16	0.45
1:B:189:ASP:CB	1:B:196:GLU:HG3	2.44	0.45
1:A:255:ILE:HG13	1:A:260:ASP:HB3	1.99	0.45
1:B:134:VAL:HA	1:B:138:LEU:HD11	1.98	0.45
1:A:151:PRO:HD2	3:A:642:HOH:O	2.16	0.45
1:A:278:ALA:HB2	1:A:295:TYR:CE2	2.52	0.45
1:B:289:HIS:CD2	1:B:291:VAL:H	2.35	0.45
1:B:165:ASP:HA	3:B:562:HOH:O	2.17	0.45
1:A:279:LYS:HD3	1:A:397:ILE:HG22	1.99	0.44
1:A:284:LEU:HD12	1:A:397:ILE:CD1	2.47	0.44
1:B:220:LEU:HD12	1:B:220:LEU:C	2.38	0.44
1:A:206:PRO:HA	1:A:207:PRO:HD3	1.84	0.44
1:B:268:ALA:O	1:B:304:THR:CB	2.66	0.44
1:A:67:LYS:HE3	1:A:75:TRP:NE1	2.32	0.44
1:A:291:VAL:HA	1:A:292:PRO:HD3	1.89	0.44
1:B:301:MSE:HE1	1:B:390:HIS:ND1	2.32	0.43
1:A:105:GLN:NE2	1:A:105:GLN:H	2.16	0.43
1:A:44:HIS:O	1:A:47:TRP:HB2	2.18	0.43
1:B:497:HIS:O	1:B:498:ARG:C	2.56	0.43
1:B:56:TRP:CG	1:B:170:LEU:HD23	2.54	0.43
1:A:91:LEU:CD1	1:A:146:ILE:HD13	2.45	0.43
1:B:321:GLU:CD	3:B:616:HOH:O	2.57	0.43
1:A:174:GLU:HA	1:A:174:GLU:OE1	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:118:ARG:HG3	1:A:123:SER:OG	2.19	0.43
1:B:179:ARG:HD3	3:B:521:HOH:O	2.19	0.43
1:A:397:ILE:HG22	1:A:397:ILE:O	2.19	0.43
1:B:92:LEU:HA	1:B:97:ALA:O	2.19	0.43
1:A:257:ARG:HD3	3:A:601:HOH:O	2.17	0.42
1:B:9:VAL:HG22	1:B:445:ALA:CB	2.49	0.42
1:B:162:GLY:HA3	3:B:556:HOH:O	2.18	0.42
1:A:63:LEU:HD13	1:A:233:LEU:HD22	2.01	0.42
1:A:110:CYS:HB2	1:A:113:GLU:CG	2.46	0.42
1:B:458:VAL:CG2	3:B:675:HOH:O	2.66	0.42
1:A:341:LEU:HA	1:A:341:LEU:HD23	1.76	0.42
1:B:380:LEU:HD22	1:B:416:TRP:CG	2.54	0.42
1:A:301:MSE:HB2	1:A:301:MSE:HE2	1.76	0.42
1:A:289:HIS:CD2	1:A:291:VAL:H	2.37	0.42
1:A:271:ILE:HD13	1:A:390:HIS:HB3	2.01	0.42
1:A:346:TYR:O	1:A:360:SER:HA	2.20	0.42
1:B:110:CYS:HB2	1:B:113:GLU:CG	2.48	0.42
1:A:284:LEU:HD13	1:A:394:LEU:CD1	2.49	0.42
1:B:131:GLY:HA2	1:B:280:SER:HB3	2.01	0.42
1:A:281:ASP:CG	1:A:282:PRO:HD2	2.41	0.42
1:B:397:ILE:CG2	1:B:397:ILE:O	2.66	0.42
1:A:348:LEU:O	1:A:358:PHE:CZ	2.72	0.42
1:A:176:VAL:HG21	1:A:220:LEU:HD13	2.02	0.41
1:A:49:GLU:HA	1:A:102:SER:O	2.19	0.41
1:B:183:LEU:HD23	1:B:287:ASP:HA	2.02	0.41
1:A:50:GLU:HG3	1:A:104:GLN:CD	2.40	0.41
1:A:143:LEU:HD12	1:A:143:LEU:HA	1.91	0.41
1:B:117:LEU:HA	1:B:117:LEU:HD23	1.93	0.41
1:B:223:VAL:HG21	1:B:233:LEU:HD23	2.02	0.41
1:B:22:ILE:HD12	1:B:24:VAL:HG21	2.03	0.41
1:B:469:THR:HA	1:B:470:PRO:HD3	1.94	0.41
1:B:264:LYS:O	1:B:269:GLY:HA2	2.20	0.41
1:A:179:ARG:NH1	3:A:653:HOH:O	2.31	0.41
1:A:404:PHE:CE2	1:A:426:ARG:HD3	2.57	0.41
1:B:290:LEU:HA	1:B:290:LEU:HD12	1.88	0.40
1:B:458:VAL:HG22	3:B:675:HOH:O	2.21	0.40
1:B:493:HIS:CE1	1:B:497:HIS:CE1	3.10	0.40
1:A:435:LEU:N	1:A:435:LEU:CD1	2.83	0.40
1:B:251:LEU:HD12	1:B:256:THR:HG22	2.03	0.40
1:A:8:GLN:HG3	1:A:76:ARG:NH2	2.36	0.40
1:B:190:LEU:HD12	1:B:190:LEU:HA	1.74	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:498:ARG:HD2	1:A:498:ARG:HH11	1.68	0.40
1:A:278:ALA:HB2	1:A:295:TYR:CD1	2.57	0.40
1:A:67:LYS:HG3	1:A:75:TRP:CD2	2.56	0.40
1:B:184:GLU:OE2	1:B:287:ASP:OD2	2.39	0.40
1:B:11:GLY:O	1:B:21:ALA:HA	2.22	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	484/508 (95%)	463 (96%)	18 (4%)	3 (1%)	30	36
1	B	472/508 (93%)	450 (95%)	18 (4%)	4 (1%)	24	27
All	All	956/1016 (94%)	913 (96%)	36 (4%)	7 (1%)	26	31

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	198	ASP
1	A	456	ASP
1	B	456	ASP
1	B	451	ASP
1	B	450	GLY
1	A	282	PRO
1	B	255	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	363/374 (97%)	326 (90%)	37 (10%)	9	10
1	B	356/374 (95%)	319 (90%)	37 (10%)	9	10
All	All	719/748 (96%)	645 (90%)	74 (10%)	9	10

All (74) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	VAL
1	A	57	ASP
1	A	62	VAL
1	A	66	LEU
1	A	68	THR
1	A	72	GLU
1	A	105	GLN
1	A	124	GLU
1	A	127	LEU
1	A	137	GLN
1	A	143	LEU
1	A	148	ARG
1	A	170	LEU
1	A	183	LEU
1	A	190	LEU
1	A	233	LEU
1	A	249	SER
1	A	263	LEU
1	A	285	TYR
1	A	315	LEU
1	A	318	GLU
1	A	321	GLU
1	A	330	LEU
1	A	333	GLU
1	A	348	LEU
1	A	357	PRO
1	A	375	LEU
1	A	385	LEU
1	A	394	LEU
1	A	425	GLN
1	A	426	ARG
1	A	433	ASN
1	A	451	ASP

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Mol	Chain	Res	Type
1	A	452	ASP
1	A	459	THR
1	A	461	LEU
1	A	498	ARG
1	B	9	VAL
1	B	28	ASP
1	B	57	ASP
1	B	66	LEU
1	B	74	ASP
1	B	86	LEU
1	B	95	ARG
1	B	105	GLN
1	B	106	SER
1	B	116	GLU
1	B	118	ARG
1	B	123	SER
1	B	129	ARG
1	B	130	THR
1	B	170	LEU
1	B	190	LEU
1	B	202	LEU
1	B	233	LEU
1	B	255	ILE
1	B	280	SER
1	B	283	ARG
1	B	290	LEU
1	B	301	MSE
1	B	333	GLU
1	B	348	LEU
1	B	372	ARG
1	B	394	LEU
1	B	403	ARG
1	B	412	ARG
1	B	456	ASP
1	B	457	ASP
1	B	463	ARG
1	B	464	THR
1	B	469	THR
1	B	476	GLU
1	B	480	ARG
1	B	498	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	58	ASN
1	A	136	GLN
1	A	289	HIS
1	A	390	HIS
1	A	399	HIS
1	A	425	GLN
1	A	493	HIS
1	A	497	HIS
1	B	44	HIS
1	B	58	ASN
1	B	105	GLN
1	B	136	GLN
1	B	149	HIS
1	B	214	HIS
1	B	289	HIS
1	B	298	ASN
1	B	425	GLN
1	B	433	ASN
1	B	493	HIS
1	B	497	HIS

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	PO4	A	509	-	4,4,4	0.25	0	6,6,6	0.27	0

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	PO4	A	509	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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, 95th 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	483/508 (95%)	0.09	8 (1%) 73 79	15, 25, 40, 54	0
1	B	475/508 (93%)	0.10	7 (1%) 76 81	16, 26, 40, 51	0
All	All	958/1016 (94%)	0.09	15 (1%) 74 80	15, 25, 40, 54	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	26	LEU	4.5
1	A	359	ALA	3.7
1	A	349	GLY	3.1
1	A	449	GLY	3.1
1	A	358	PHE	3.0
1	A	280	SER	2.9
1	A	356	ASP	2.7
1	B	453	LEU	2.7
1	A	463	ARG	2.4
1	B	358	PHE	2.3
1	B	283	ARG	2.3
1	B	128	ALA	2.2
1	A	191	ALA	2.2
1	B	476	GLU	2.1
1	B	285	TYR	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, 95th 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(\AA^2)	Q<0.9
2	PO4	A	509	5/5	0.70	0.34	7.72	30,30,30,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.