



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 2, 2022 – 12:12 PM JST

PDB ID : 7F0E
Title : Crystal Structure of EnPKS2
Authors : Huang, S.X.; Yan, Y.J.
Deposited on : 2021-06-03
Resolution : 2.62 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	:	2.29
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)	:	2.29

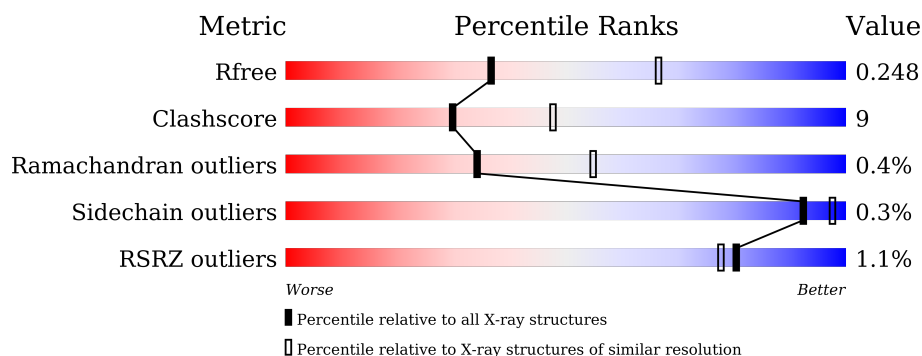
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.62 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	3797 (2.64-2.60)
Clashscore	141614	4168 (2.64-2.60)
Ramachandran outliers	138981	4093 (2.64-2.60)
Sidechain outliers	138945	4093 (2.64-2.60)
RSRZ outliers	127900	3731 (2.64-2.60)

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 $\leq 5\%$. 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	390	<div> <div style="width: 82%; background-color: green;"></div> <div style="width: 14%; background-color: yellow;"></div> <div style="width: 4%; background-color: orange;"></div> <div style="width: 2%; background-color: red;"></div> <div style="width: 2%; background-color: grey;"></div> </div> <div>82% 14% . .</div>
1	B	390	<div> <div style="width: 75%; background-color: green;"></div> <div style="width: 20%; background-color: yellow;"></div> <div style="width: 5%; background-color: orange;"></div> <div style="width: 2%; background-color: red;"></div> <div style="width: 2%; background-color: grey;"></div> </div> <div>% 75% 20% . .</div>
1	C	390	<div> <div style="width: 72%; background-color: green;"></div> <div style="width: 22%; background-color: yellow;"></div> <div style="width: 6%; background-color: orange;"></div> <div style="width: 2%; background-color: red;"></div> <div style="width: 2%; background-color: grey;"></div> </div> <div>% 72% 22% . .</div>
1	D	390	<div> <div style="width: 74%; background-color: green;"></div> <div style="width: 21%; background-color: yellow;"></div> <div style="width: 5%; background-color: orange;"></div> <div style="width: 2%; background-color: red;"></div> <div style="width: 2%; background-color: grey;"></div> </div> <div>2% 74% 21% . .</div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11598 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 EnPKS2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	374	Total	C	N	O	S	0	0	0
			2872	1827	493	535	17			
1	B	376	Total	C	N	O	S	0	0	0
			2889	1838	497	537	17			
1	C	374	Total	C	N	O	S	0	0	0
			2875	1830	493	535	17			
1	D	377	Total	C	N	O	S	0	0	0
			2900	1844	501	538	17			

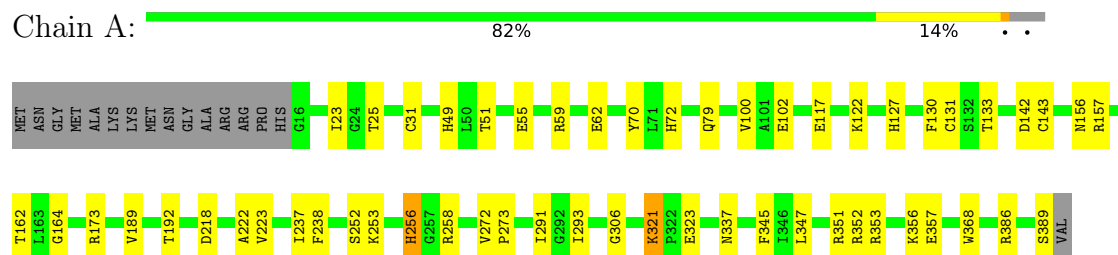
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	25	Total	O	0	0
			25	25		
2	B	21	Total	O	0	0
			21	21		
2	C	13	Total	O	0	0
			13	13		
2	D	3	Total	O	0	0
			3	3		

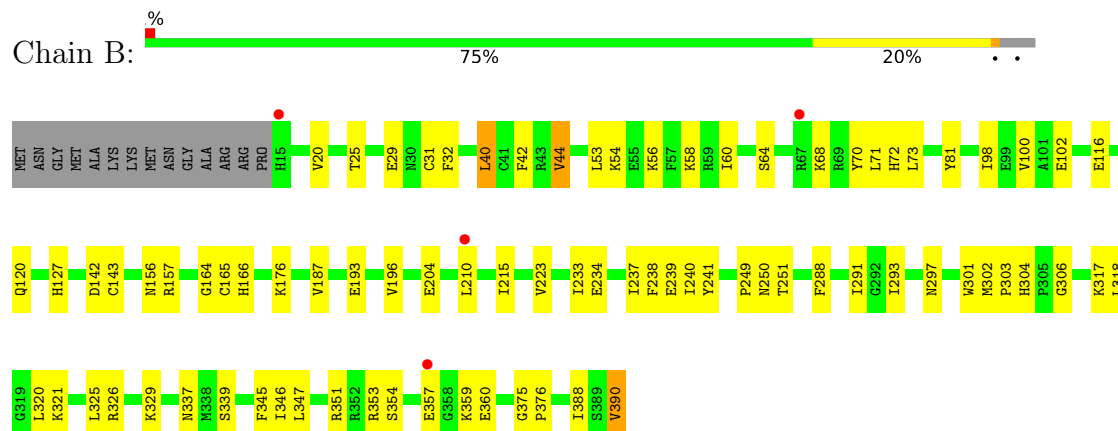
3 Residue-property plots [i](#)

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.

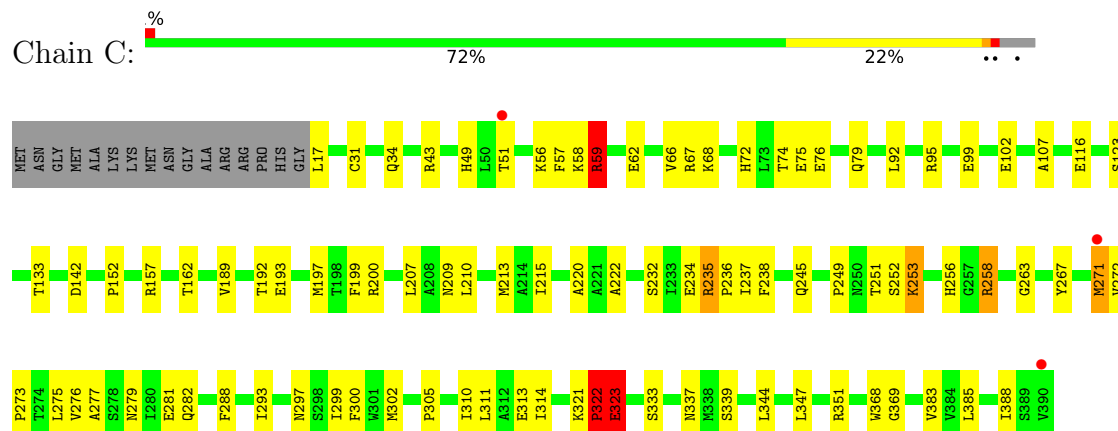
• Molecule 1: EnPKS2



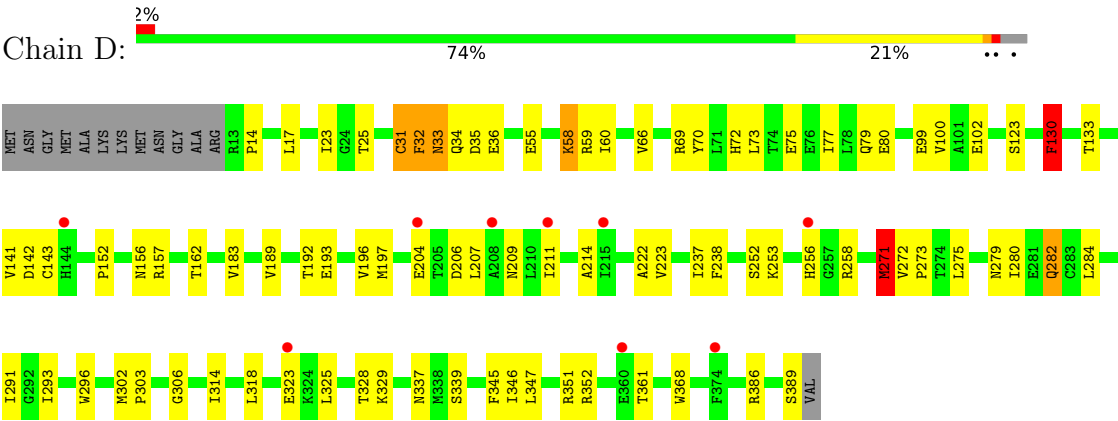
• Molecule 1: EnPKS2



• Molecule 1: EnPKS2



● Molecule 1: EnPKS2



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	95.18Å 112.95Å 154.71Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	91.22 – 2.62 91.22 – 2.62	Depositor EDS
% Data completeness (in resolution range)	99.6 (91.22-2.62) 99.6 (91.22-2.62)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.28 (at 2.62Å)	Xtriage
Refinement program	PHENIX 1.18.1_3865	Depositor
R, R_{free}	0.196 , 0.247 0.196 , 0.248	Depositor DCC
R_{free} test set	1999 reflections (3.95%)	wwPDB-VP
Wilson B-factor (Å ²)	54.4	Xtriage
Anisotropy	0.487	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 38.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11598	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.04% 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, 0.333 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	0.56	3/2927 (0.1%)	0.77	3/3966 (0.1%)
1	B	0.57	5/2945 (0.2%)	0.72	3/3991 (0.1%)
1	C	0.51	2/2930 (0.1%)	1.08	14/3971 (0.4%)
1	D	0.49	3/2957 (0.1%)	0.87	16/4007 (0.4%)
All	All	0.53	13/11759 (0.1%)	0.87	36/15935 (0.2%)

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	C	0	3
1	D	0	2
All	All	0	5

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	131	CYS	CB-SG	-6.43	1.71	1.82
1	D	36	GLU	CD-OE2	6.29	1.32	1.25
1	B	390	VAL	CB-CG2	6.20	1.65	1.52
1	B	116	GLU	CD-OE1	6.11	1.32	1.25
1	A	62	GLU	CD-OE1	6.03	1.32	1.25
1	B	116	GLU	CD-OE2	5.99	1.32	1.25
1	A	62	GLU	CD-OE2	5.92	1.32	1.25
1	D	204	GLU	CD-OE1	5.49	1.31	1.25
1	C	323	GLU	CD-OE2	5.48	1.31	1.25
1	B	116	GLU	CG-CD	5.31	1.59	1.51
1	D	36	GLU	CD-OE1	5.13	1.31	1.25
1	C	323	GLU	CG-CD	-5.13	1.44	1.51
1	B	29	GLU	CG-CD	5.02	1.59	1.51

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	323	GLU	OE1-CD-OE2	-23.89	94.64	123.30
1	C	59	ARG	NE-CZ-NH2	-21.93	109.33	120.30
1	C	59	ARG	NE-CZ-NH1	17.11	128.86	120.30
1	C	323	GLU	CG-CD-OE2	16.79	151.89	118.30
1	C	323	GLU	CG-CD-OE1	-14.31	89.69	118.30
1	D	271	MET	CA-CB-CG	12.99	135.39	113.30
1	C	59	ARG	CD-NE-CZ	11.69	139.96	123.60
1	D	207	LEU	CB-CG-CD2	-10.71	92.79	111.00
1	D	271	MET	CB-CG-SD	9.69	141.47	112.40
1	C	271	MET	CA-CB-CG	7.68	126.35	113.30
1	A	256	HIS	CB-CA-C	-7.57	95.27	110.40
1	C	116	GLU	CA-CB-CG	7.51	129.92	113.40
1	B	40	LEU	CB-CG-CD2	7.34	123.47	111.00
1	A	321	LYS	CD-CE-NZ	-7.12	95.33	111.70
1	D	271	MET	CB-CA-C	-7.07	96.26	110.40
1	D	282	GLN	CA-CB-CG	6.68	128.11	113.40
1	D	59	ARG	NE-CZ-NH1	-6.60	117.00	120.30
1	D	352	ARG	CA-CB-CG	6.59	127.91	113.40
1	B	116	GLU	CA-CB-CG	6.36	127.39	113.40
1	C	271	MET	CB-CA-C	-6.08	98.25	110.40
1	D	207	LEU	CB-CG-CD1	6.07	121.32	111.00
1	C	258	ARG	CG-CD-NE	-5.92	99.37	111.80
1	B	44	VAL	CG1-CB-CG2	5.83	120.23	110.90
1	D	130	PHE	CB-CG-CD2	-5.81	116.73	120.80
1	C	79	GLN	CA-CB-CG	5.80	126.17	113.40
1	C	258	ARG	NE-CZ-NH1	-5.80	117.40	120.30
1	D	36	GLU	OE1-CD-OE2	5.59	130.01	123.30
1	D	284	LEU	CA-CB-CG	5.59	128.16	115.30
1	A	122	LYS	CD-CE-NZ	-5.58	98.86	111.70
1	D	204	GLU	CA-CB-CG	5.46	125.42	113.40
1	D	352	ARG	CG-CD-NE	5.36	123.06	111.80
1	C	235	ARG	CG-CD-NE	5.34	123.02	111.80
1	D	207	LEU	CA-CB-CG	5.25	127.37	115.30
1	C	59	ARG	CG-CD-NE	5.20	122.73	111.80
1	D	58	LYS	CA-CB-CG	5.08	124.59	113.40
1	D	59	ARG	NE-CZ-NH2	5.05	122.82	120.30

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	322	PRO	Peptide
1	C	323	GLU	Sidechain
1	C	59	ARG	Sidechain
1	D	130	PHE	Sidechain
1	D	271	MET	Peptide

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	2872	0	2933	43	1
1	B	2889	0	2949	54	0
1	C	2875	0	2939	62	1
1	D	2900	0	2960	64	0
2	A	25	0	0	2	0
2	B	21	0	0	2	0
2	C	13	0	0	2	0
2	D	3	0	0	0	0
All	All	11598	0	11781	218	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 9.

All (218) 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:31:CYS:SG	2:A:420:HOH:O	2.34	0.85
1:B:31:CYS:SG	2:B:421:HOH:O	2.33	0.84
1:C:200:ARG:H	1:C:213:MET:HE1	1.41	0.83
1:C:56:LYS:HA	1:C:59:ARG:HD2	1.60	0.81
1:A:55:GLU:HG3	1:A:59:ARG:HH12	1.44	0.81
1:A:130:PHE:HE1	1:A:192:THR:HG22	1.51	0.75
1:B:143:CYS:SG	1:B:157:ARG:NH1	2.64	0.70
1:B:40:LEU:HD22	1:B:44:VAL:CG2	2.22	0.69
1:C:107:ALA:HB2	1:C:192:THR:HG21	1.73	0.69
1:D:252:SER:O	1:D:253:LYS:HG2	1.93	0.68
1:B:73:LEU:HG	1:B:196:VAL:HG22	1.75	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:302:MET:HE2	1:C:347:LEU:HA	1.76	0.67
1:B:233:ILE:HG22	1:B:234:GLU:HG3	1.76	0.66
1:D:69:ARG:NH1	1:D:214:ALA:O	2.29	0.65
1:D:130:PHE:HE1	1:D:192:THR:HG22	1.61	0.65
1:A:389:SER:N	2:A:401:HOH:O	2.14	0.65
1:A:253:LYS:HA	1:A:256:HIS:CD2	2.32	0.64
1:D:73:LEU:HG	1:D:196:VAL:HG22	1.79	0.64
1:C:17:LEU:O	2:C:401:HOH:O	2.15	0.64
1:D:302:MET:HE1	1:D:347:LEU:HD23	1.79	0.63
1:C:273:PRO:HA	1:C:310:ILE:HD13	1.81	0.62
1:C:310:ILE:O	1:C:314:ILE:HG12	2.00	0.62
1:B:318:LEU:HB2	1:B:320:LEU:HD13	1.82	0.62
1:A:133:THR:O	1:A:162:THR:HG22	1.99	0.62
1:C:58:LYS:O	1:C:62:GLU:HG3	1.99	0.62
1:D:302:MET:HG2	1:D:346:ILE:HG22	1.83	0.61
1:C:252:SER:O	1:C:253:LYS:HB3	2.00	0.61
1:D:253:LYS:HB2	1:D:256:HIS:NE2	2.16	0.60
1:A:55:GLU:CG	1:A:59:ARG:HH12	2.14	0.60
1:C:133:THR:O	1:C:162:THR:HG22	2.01	0.60
1:A:272:VAL:HG22	1:A:273:PRO:HD3	1.84	0.60
1:D:34:GLN:HE22	1:D:69:ARG:NH1	2.00	0.60
1:B:165:CYS:HB3	1:B:304:HIS:CE1	2.38	0.59
1:B:193:GLU:HG3	1:B:339:SER:HB3	1.84	0.59
1:C:99:GLU:OE2	1:C:197:MET:HB2	2.03	0.59
1:C:237:ILE:HG21	1:C:351:ARG:HD3	1.85	0.58
1:D:31:CYS:C	1:D:32:PHE:HD1	2.07	0.58
1:B:40:LEU:HD22	1:B:44:VAL:HG22	1.84	0.58
1:B:176:LYS:HE2	1:B:240:ILE:O	2.03	0.58
1:C:252:SER:O	1:C:253:LYS:CB	2.51	0.58
1:D:33:ASN:OD1	1:D:35:ASP:N	2.37	0.58
1:D:34:GLN:NE2	1:D:69:ARG:NH1	2.53	0.57
1:D:32:PHE:HD1	1:D:32:PHE:N	2.01	0.57
1:D:33:ASN:OD1	1:D:33:ASN:C	2.41	0.57
1:D:133:THR:O	1:D:162:THR:HG23	2.04	0.57
1:D:23:ILE:HG12	1:D:223:VAL:HG12	1.86	0.57
1:C:236:PRO:HB2	1:C:388:ILE:CD1	2.34	0.56
1:B:53:LEU:HD11	1:B:210:LEU:HD11	1.86	0.56
1:C:238:PHE:HB3	1:C:385:LEU:HD12	1.88	0.56
1:D:130:PHE:HE1	1:D:192:THR:CG2	2.19	0.56
1:B:20:VAL:HG13	1:B:223:VAL:HG13	1.87	0.56
1:A:253:LYS:HD3	1:A:256:HIS:NE2	2.22	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:302:MET:HG2	1:B:346:ILE:HG22	1.88	0.55
1:D:33:ASN:OD1	1:D:34:GLN:N	2.39	0.55
1:A:72:HIS:NE2	1:A:102:GLU:HG3	2.21	0.55
1:C:72:HIS:CE1	1:C:102:GLU:HG3	2.42	0.55
1:D:279:ASN:O	1:D:282:GLN:HB3	2.07	0.55
1:B:353:ARG:O	1:B:357:GLU:HG3	2.07	0.55
1:D:237:ILE:HG21	1:D:351:ARG:HD3	1.89	0.54
1:B:40:LEU:CD2	1:B:44:VAL:HG22	2.37	0.54
1:A:189:VAL:O	1:A:222:ALA:HA	2.07	0.54
1:D:206:ASP:OD2	1:D:209:ASN:ND2	2.41	0.54
1:D:32:PHE:N	1:D:32:PHE:CD1	2.76	0.54
1:C:200:ARG:H	1:C:213:MET:CE	2.15	0.54
1:B:127:HIS:CD2	1:B:156:ASN:HB2	2.43	0.53
1:B:53:LEU:HD22	1:B:204:GLU:HG2	1.90	0.53
1:C:281:GLU:HB2	2:C:404:HOH:O	2.07	0.53
1:A:142:ASP:OD2	1:A:157:ARG:HD2	2.07	0.53
1:C:67:ARG:HB2	1:C:333:SER:O	2.08	0.53
1:C:200:ARG:HD2	1:C:209:ASN:HD21	1.73	0.53
1:B:157:ARG:HB2	1:C:245:GLN:OE1	2.10	0.52
1:D:130:PHE:CE2	1:D:141:VAL:HG23	2.44	0.52
1:D:253:LYS:HA	1:D:256:HIS:CD2	2.44	0.52
1:C:95:ARG:HD3	1:C:197:MET:HA	1.92	0.52
1:A:25:THR:HB	1:A:345:PHE:CZ	2.44	0.52
1:B:302:MET:SD	1:B:347:LEU:HD23	2.50	0.52
1:A:117:GLU:OE2	1:A:351:ARG:NH1	2.41	0.51
1:C:272:VAL:HG12	1:C:310:ILE:HD11	1.91	0.51
1:B:291:ILE:HG22	1:B:293:ILE:HG13	1.91	0.51
1:B:53:LEU:HD22	1:B:204:GLU:CG	2.39	0.51
1:D:72:HIS:CE1	1:D:102:GLU:HG3	2.46	0.51
1:D:72:HIS:NE2	1:D:102:GLU:HG3	2.26	0.51
1:D:272:VAL:HG22	1:D:273:PRO:HD3	1.92	0.51
1:D:142:ASP:OD2	1:D:157:ARG:HD2	2.11	0.51
1:B:164:GLY:HA3	1:B:376:PRO:HD2	1.93	0.51
1:A:130:PHE:CE1	1:A:192:THR:HG22	2.39	0.50
1:C:302:MET:CE	1:C:347:LEU:HA	2.41	0.50
1:C:215:ILE:HG22	1:C:337:ASN:HB3	1.94	0.50
1:D:302:MET:HG2	1:D:346:ILE:CG2	2.41	0.50
1:C:288:PHE:HB3	1:C:293:ILE:HB	1.93	0.50
1:C:234:GLU:C	1:C:235:ARG:HD2	2.32	0.50
1:B:100:VAL:HG21	1:C:258:ARG:NH1	2.27	0.50
1:C:189:VAL:O	1:C:222:ALA:HA	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:200:ARG:N	1:C:213:MET:HE1	2.20	0.49
1:A:353:ARG:HD2	1:A:357:GLU:OE1	2.13	0.49
1:B:249:PRO:O	1:B:251:THR:HG23	2.11	0.49
1:C:276:VAL:HG11	1:C:310:ILE:HG23	1.95	0.49
1:A:252:SER:O	1:A:253:LYS:HB2	2.12	0.49
1:A:55:GLU:OE2	1:A:59:ARG:NH1	2.46	0.48
1:A:321:LYS:HB3	1:A:323:GLU:OE1	2.13	0.48
1:D:60:ILE:HD13	1:D:211:ILE:HD13	1.95	0.48
1:A:272:VAL:CG2	1:A:273:PRO:HD3	2.43	0.48
1:C:236:PRO:HB2	1:C:388:ILE:HD13	1.95	0.48
1:D:302:MET:CE	1:D:347:LEU:HD23	2.44	0.48
1:D:314:ILE:O	1:D:318:LEU:HB2	2.13	0.48
1:C:92:LEU:HD13	1:C:263:GLY:HA2	1.95	0.48
1:B:239:GLU:HG3	1:B:388:ILE:HD11	1.96	0.48
1:B:357:GLU:HB3	1:B:359:LYS:HG3	1.95	0.48
1:C:200:ARG:HD2	1:C:209:ASN:ND2	2.28	0.48
1:D:303:PRO:HD2	1:D:328:THR:OG1	2.14	0.48
1:B:72:HIS:NE2	1:B:102:GLU:HG3	2.28	0.47
1:B:306:GLY:HA2	1:B:337:ASN:ND2	2.28	0.47
1:D:361:THR:HA	1:D:389:SER:HA	1.96	0.47
1:B:54:LYS:O	1:B:58:LYS:HG3	2.14	0.47
1:B:238:PHE:CD2	1:B:347:LEU:HD22	2.49	0.47
1:C:49:HIS:O	1:C:51:THR:HG23	2.15	0.47
1:C:383:VAL:HG12	1:C:385:LEU:HD22	1.97	0.47
1:D:123:SER:HA	1:D:152:PRO:HG3	1.96	0.47
1:D:143:CYS:SG	1:D:157:ARG:NH1	2.88	0.47
1:C:256:HIS:HB2	1:C:267:TYR:HB3	1.96	0.47
1:B:326:ARG:CZ	1:B:353:ARG:HD2	2.44	0.46
1:C:123:SER:HA	1:C:152:PRO:HG3	1.97	0.46
1:B:56:LYS:O	1:B:60:ILE:HG12	2.14	0.46
1:D:130:PHE:CE1	1:D:192:THR:HG22	2.45	0.46
1:C:57:PHE:CD1	1:C:210:LEU:HD22	2.50	0.46
1:D:271:MET:HG3	1:D:275:LEU:HD22	1.96	0.46
1:A:127:HIS:CD2	1:A:156:ASN:HB2	2.51	0.46
1:A:351:ARG:HG3	1:A:352:ARG:N	2.30	0.46
1:C:72:HIS:NE2	1:C:102:GLU:HG3	2.30	0.46
1:A:72:HIS:CE1	1:A:102:GLU:HG3	2.51	0.46
1:C:34:GLN:HG2	1:C:66:VAL:O	2.16	0.46
1:D:238:PHE:CD2	1:D:347:LEU:HD22	2.51	0.45
1:A:173:ARG:NH1	1:D:156:ASN:HB3	2.32	0.45
1:C:95:ARG:O	1:C:99:GLU:HG3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:296:TRP:CD1	1:D:318:LEU:CD1	2.99	0.45
1:D:193:GLU:HG3	1:D:339:SER:HB3	1.98	0.45
1:D:291:ILE:HG22	1:D:293:ILE:HG13	1.97	0.45
1:C:43:ARG:HH21	1:C:75:GLU:CD	2.20	0.45
1:D:31:CYS:C	1:D:32:PHE:CD1	2.90	0.45
1:D:55:GLU:OE1	1:D:58:LYS:HD3	2.17	0.45
1:D:99:GLU:OE1	1:D:197:MET:HB2	2.17	0.45
1:C:142:ASP:OD2	1:C:157:ARG:HD2	2.17	0.45
1:A:59:ARG:HG2	1:A:59:ARG:HH11	1.82	0.45
1:C:305:PRO:HA	1:C:311:LEU:HD21	1.98	0.45
1:D:272:VAL:N	1:D:273:PRO:CD	2.80	0.45
1:D:189:VAL:O	1:D:222:ALA:HA	2.17	0.44
1:D:280:ILE:HD13	1:D:314:ILE:HD12	1.98	0.44
1:B:237:ILE:HG21	1:B:351:ARG:HD3	1.98	0.44
1:B:304:HIS:HE1	2:B:419:HOH:O	1.99	0.44
1:C:344:LEU:O	1:C:347:LEU:HB2	2.18	0.44
1:A:237:ILE:HG21	1:A:351:ARG:HD3	1.99	0.44
1:C:193:GLU:HG3	1:C:339:SER:HB3	2.00	0.44
1:B:81:TYR:CD2	1:B:98:ILE:HD11	2.51	0.44
1:A:253:LYS:HD3	1:A:256:HIS:CE1	2.53	0.44
1:B:42:PHE:CZ	1:B:54:LYS:HA	2.53	0.44
1:C:271:MET:HB3	1:C:275:LEU:HG	2.00	0.44
1:C:279:ASN:O	1:C:282:GLN:HB3	2.18	0.44
1:D:77:ILE:O	1:D:80:GLU:HB3	2.18	0.43
1:A:253:LYS:HA	1:A:256:HIS:HD2	1.80	0.43
1:D:31:CYS:HB3	1:D:70:TYR:CD2	2.53	0.43
1:C:192:THR:HG22	1:C:220:ALA:HB1	2.00	0.43
1:A:258:ARG:NH1	1:D:100:VAL:HG21	2.33	0.43
1:C:68:LYS:HB2	1:C:68:LYS:HE3	1.72	0.43
1:A:189:VAL:HB	1:A:223:VAL:HG23	2.00	0.43
1:C:133:THR:O	1:C:197:MET:HE1	2.18	0.43
1:C:249:PRO:O	1:C:251:THR:HG23	2.19	0.43
1:A:368:TRP:CH2	1:A:386:ARG:HD2	2.53	0.43
1:B:120:GLN:NE2	1:B:233:ILE:HG23	2.33	0.43
1:B:301:TRP:HE3	1:B:303:PRO:HG3	1.84	0.43
1:A:79:GLN:HE21	1:A:79:GLN:HB3	1.50	0.42
1:B:166:HIS:HB3	1:B:375:GLY:N	2.34	0.42
1:C:74:THR:HB	1:C:76:GLU:OE1	2.19	0.42
1:C:322:PRO:CA	1:C:323:GLU:HG3	2.48	0.42
1:D:34:GLN:HG2	1:D:66:VAL:O	2.19	0.42
1:D:75:GLU:O	1:D:79:GLN:HG3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:130:PHE:CE1	1:D:192:THR:CG2	3.01	0.42
1:C:300:PHE:CE2	1:C:369:GLY:HA3	2.54	0.42
1:B:25:THR:HB	1:B:345:PHE:CZ	2.54	0.42
1:D:31:CYS:HB3	1:D:70:TYR:CE2	2.54	0.42
1:D:306:GLY:HA2	1:D:337:ASN:ND2	2.34	0.42
1:D:325:LEU:O	1:D:329:LYS:HG3	2.19	0.42
1:A:49:HIS:O	1:A:51:THR:HG23	2.20	0.42
1:A:291:ILE:HG22	1:A:293:ILE:HG12	2.01	0.42
1:A:306:GLY:HA2	1:A:337:ASN:ND2	2.35	0.42
1:D:323:GLU:H	1:D:323:GLU:HG2	1.59	0.42
1:B:40:LEU:HD22	1:B:44:VAL:HG21	2.01	0.42
1:B:142:ASP:OD2	1:B:157:ARG:HD2	2.20	0.42
1:D:206:ASP:CG	1:D:209:ASN:HD22	2.22	0.42
1:A:23:ILE:HG12	1:A:223:VAL:HG12	2.02	0.41
1:A:70:TYR:O	1:A:218:ASP:HB2	2.20	0.41
1:A:323:GLU:H	1:A:323:GLU:CD	2.23	0.41
1:B:127:HIS:HB2	1:B:187:VAL:HG13	2.01	0.41
1:C:277:ALA:HB2	1:C:313:GLU:HB3	2.02	0.41
1:C:133:THR:HA	1:C:162:THR:HA	2.02	0.41
1:A:238:PHE:CD2	1:A:347:LEU:HD22	2.55	0.41
1:C:199:PHE:HA	1:C:213:MET:CE	2.50	0.41
1:C:207:LEU:HD12	1:C:207:LEU:HA	1.86	0.41
1:D:25:THR:HB	1:D:345:PHE:CZ	2.55	0.41
1:B:360:GLU:O	1:B:390:VAL:N	2.48	0.41
1:B:297:ASN:OD1	1:B:321:LYS:HG3	2.21	0.41
1:B:325:LEU:O	1:B:329:LYS:HG3	2.20	0.41
1:C:297:ASN:HB3	1:C:321:LYS:HE2	2.02	0.41
1:A:100:VAL:HG21	1:D:258:ARG:NH1	2.35	0.41
1:A:351:ARG:CG	1:A:352:ARG:N	2.83	0.41
1:B:64:SER:HB2	1:B:215:ILE:HD13	2.03	0.41
1:B:241:TYR:CZ	1:B:291:ILE:HD11	2.56	0.41
1:D:14:PRO:HG3	1:D:183:VAL:HA	2.02	0.41
1:D:237:ILE:HG21	1:D:351:ARG:CD	2.51	0.41
1:A:356:LYS:HD3	1:A:356:LYS:C	2.41	0.40
1:B:32:PHE:CE1	1:B:71:LEU:HB2	2.56	0.40
1:B:68:LYS:HE3	1:B:70:TYR:OH	2.21	0.40
1:B:317:LYS:HD2	1:B:317:LYS:HA	1.82	0.40
1:B:288:PHE:HB3	1:B:293:ILE:HB	2.03	0.40
1:A:143:CYS:SG	1:A:157:ARG:NH1	2.94	0.40
1:C:299:ILE:HA	1:C:368:TRP:O	2.22	0.40
1:B:72:HIS:CE1	1:B:102:GLU:HG3	2.57	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:354:SER:HA	1:B:357:GLU:OE1	2.21	0.40
1:D:368:TRP:CH2	1:D:386:ARG:HD2	2.57	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:356:LYS:NZ	1:C:232:SER:OG[4_456]	2.01	0.19

5.3 Torsion angles

5.3.1 Protein backbone

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	372/390 (95%)	360 (97%)	11 (3%)	1 (0%)	41	62
1	B	374/390 (96%)	358 (96%)	15 (4%)	1 (0%)	41	62
1	C	372/390 (95%)	351 (94%)	18 (5%)	3 (1%)	19	36
1	D	375/390 (96%)	356 (95%)	18 (5%)	1 (0%)	41	62
All	All	1493/1560 (96%)	1425 (95%)	62 (4%)	6 (0%)	34	55

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	323	GLU
1	D	17	LEU
1	B	250	ASN
1	C	253	LYS
1	C	322	PRO
1	A	164	GLY

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	314/326 (96%)	314 (100%)	0	100	100
1	B	316/326 (97%)	316 (100%)	0	100	100
1	C	315/326 (97%)	314 (100%)	1 (0%)	92	97
1	D	317/326 (97%)	314 (99%)	3 (1%)	78	90
All	All	1262/1304 (97%)	1258 (100%)	4 (0%)	92	97

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	31	CYS
1	D	31	CYS
1	D	32	PHE
1	D	33	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	79	GLN
1	A	144	HIS
1	A	209	ASN
1	B	120	GLN
1	B	144	HIS
1	B	282	GLN
1	B	304	HIS
1	C	34	GLN
1	C	120	GLN
1	C	181	ASN
1	C	209	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 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 [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	374/390 (95%)	-0.18	0 100 100	38, 50, 66, 87	0
1	B	376/390 (96%)	-0.11	4 (1%) 80 78	37, 51, 72, 106	0
1	C	374/390 (95%)	0.04	3 (0%) 86 84	40, 61, 87, 129	0
1	D	377/390 (96%)	0.06	9 (2%) 59 53	41, 64, 90, 123	0
All	All	1501/1560 (96%)	-0.05	16 (1%) 80 78	37, 55, 83, 129	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	215	ILE	3.6
1	B	15	HIS	3.6
1	D	374	PHE	3.5
1	C	390	VAL	3.2
1	D	360	GLU	3.2
1	D	204	GLU	2.7
1	C	271	MET	2.6
1	D	144	HIS	2.5
1	B	357	GLU	2.3
1	C	51	THR	2.2
1	D	256	HIS	2.2
1	D	208	ALA	2.2
1	B	67	ARG	2.1
1	D	211	ILE	2.0
1	B	210	LEU	2.0
1	D	323	GLU	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.