



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:11 AM GMT

PDB ID : 2PQ5
Title : Crystal structure of Dual specificity protein phosphatase 13 (DUSP13)
Authors : Ugochukwu, E.; Salah, E.; Savitsky, P.; Barr, A.; Pantic, N.; Niesen, F.; Burgess-Brown, N.; Berridge, G.; Bunkoczi, G.; Uppenberg, J.; Pike, A.C.W.; Sundstrom, M.; Arrowsmith, C.H.; Weigelt, J.; Edwards, A.; von Delft, F.; Knapp, S.; Structural Genomics Consortium (SGC)
Deposited on : 2007-05-01
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

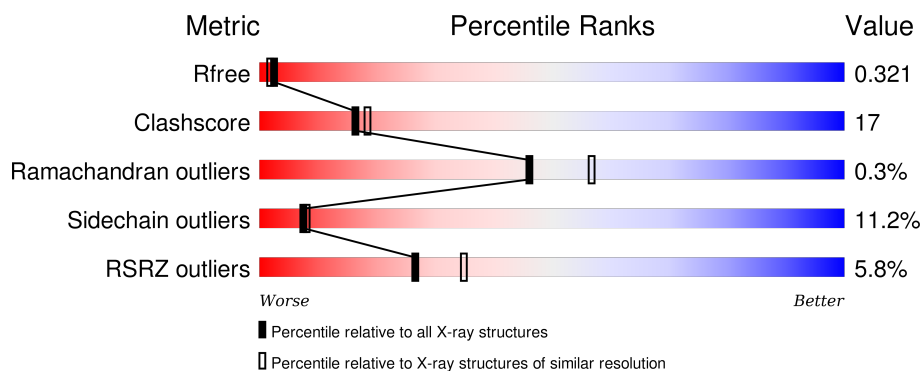
1 Overall quality at a glance ⓘ

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	205	<div> <div>3%</div> <div>52% 27% 18%</div> </div>
1	B	205	<div> <div>3%</div> <div>48% 28% 5% 18%</div> </div>
1	C	205	<div> <div>8%</div> <div>46% 29% 7% 18%</div> </div>
1	D	205	<div> <div>4%</div> <div>47% 30% 18%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5258 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 Dual specificity protein phosphatase 13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	169	Total	C	N	O	S	0	0	0
			1288	832	217	233	6			
1	B	168	Total	C	N	O	S	0	0	0
			1288	833	217	232	6			
1	C	169	Total	C	N	O	S	0	0	0
			1285	828	219	232	6			
1	D	169	Total	C	N	O	S	0	0	0
			1305	841	225	233	6			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	199	ALA	-	CLONING ARTIFACT	UNP Q9UII6
A	200	HIS	-	CLONING ARTIFACT	UNP Q9UII6
A	201	HIS	-	CLONING ARTIFACT	UNP Q9UII6
A	202	HIS	-	CLONING ARTIFACT	UNP Q9UII6
A	203	HIS	-	CLONING ARTIFACT	UNP Q9UII6
A	204	HIS	-	CLONING ARTIFACT	UNP Q9UII6
A	205	HIS	-	CLONING ARTIFACT	UNP Q9UII6
B	199	ALA	-	CLONING ARTIFACT	UNP Q9UII6
B	200	HIS	-	CLONING ARTIFACT	UNP Q9UII6
B	201	HIS	-	CLONING ARTIFACT	UNP Q9UII6
B	202	HIS	-	CLONING ARTIFACT	UNP Q9UII6
B	203	HIS	-	CLONING ARTIFACT	UNP Q9UII6
B	204	HIS	-	CLONING ARTIFACT	UNP Q9UII6
B	205	HIS	-	CLONING ARTIFACT	UNP Q9UII6
C	199	ALA	-	CLONING ARTIFACT	UNP Q9UII6
C	200	HIS	-	CLONING ARTIFACT	UNP Q9UII6
C	201	HIS	-	CLONING ARTIFACT	UNP Q9UII6
C	202	HIS	-	CLONING ARTIFACT	UNP Q9UII6
C	203	HIS	-	CLONING ARTIFACT	UNP Q9UII6
C	204	HIS	-	CLONING ARTIFACT	UNP Q9UII6
C	205	HIS	-	CLONING ARTIFACT	UNP Q9UII6

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Chain	Residue	Modelled	Actual	Comment	Reference
D	199	ALA	-	CLONING ARTIFACT	UNP Q9UII6
D	200	HIS	-	CLONING ARTIFACT	UNP Q9UII6
D	201	HIS	-	CLONING ARTIFACT	UNP Q9UII6
D	202	HIS	-	CLONING ARTIFACT	UNP Q9UII6
D	203	HIS	-	CLONING ARTIFACT	UNP Q9UII6
D	204	HIS	-	CLONING ARTIFACT	UNP Q9UII6
D	205	HIS	-	CLONING ARTIFACT	UNP Q9UII6

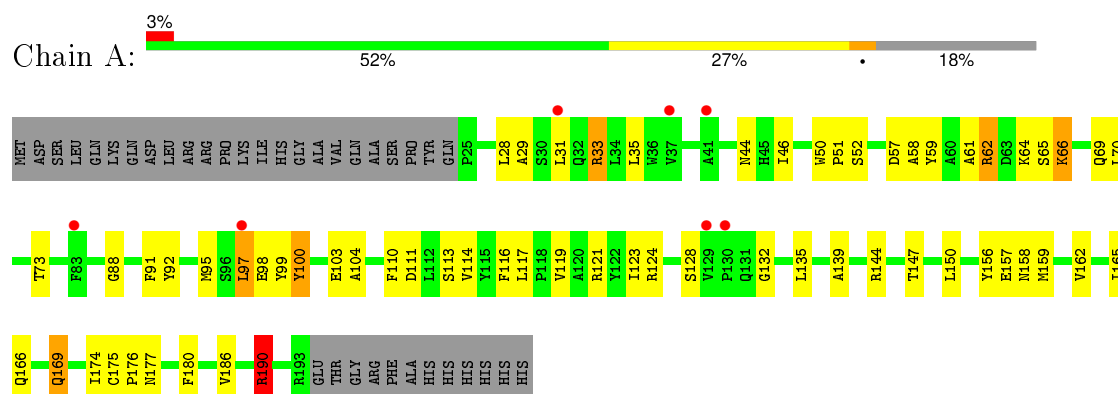
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	32	Total O 32 32	0	0
2	B	25	Total O 25 25	0	0
2	C	20	Total O 20 20	0	0
2	D	15	Total O 15 15	0	0

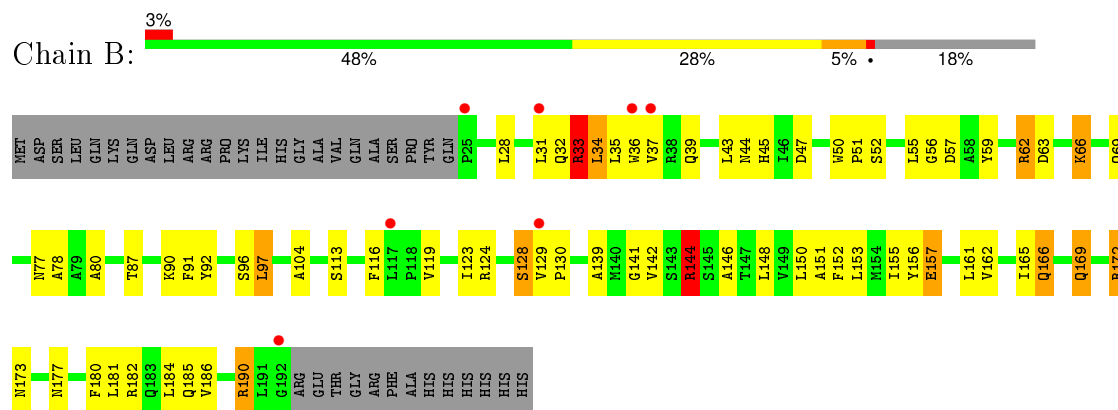
3 Residue-property plots [i](#)

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.

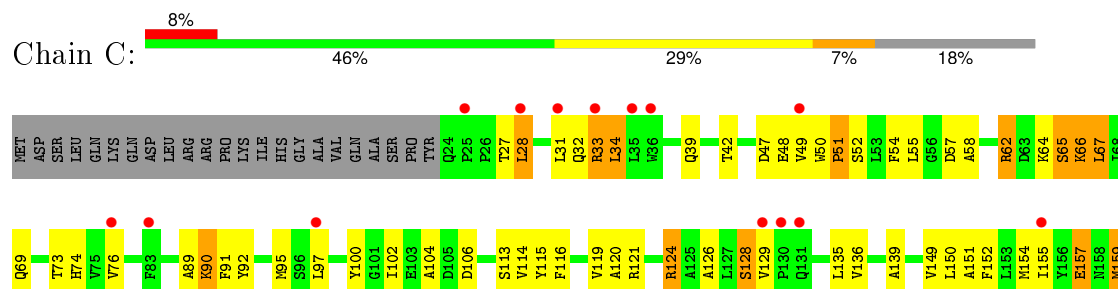
• Molecule 1: Dual specificity protein phosphatase 13

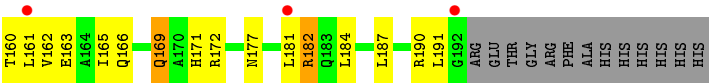


• Molecule 1: Dual specificity protein phosphatase 13

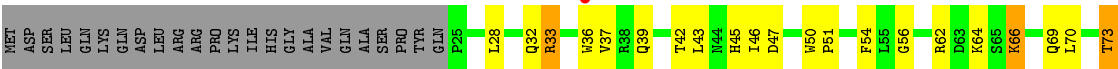


• Molecule 1: Dual specificity protein phosphatase 13





● Molecule 1: Dual specificity protein phosphatase 13



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	73.82Å 73.82Å 303.80Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.30 75.95 – 2.30	Depositor EDS
% Data completeness (in resolution range)	95.2 (20.00-2.30) 97.9 (75.95-2.30)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.04 (at 2.29Å)	Xtriage
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.176 , 0.260 0.177 , 0.321	Depositor DCC
R_{free} test set	16 reflections (0.04%)	DCC
Wilson B-factor (Å ²)	34.3	Xtriage
Anisotropy	0.288	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 83.1	EDS
Estimated twinning fraction	0.460 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.38$, $\langle L^2 \rangle = 0.20$	Xtriage
Outliers	0 of 43981 reflections	Xtriage
F_o, F_c correlation	0.82	EDS
Total number of atoms	5258	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.22% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.35	0/1317	1.26	9/1797 (0.5%)
1	B	0.36	0/1317	1.29	9/1795 (0.5%)
1	C	0.35	0/1312	1.26	8/1788 (0.4%)
1	D	0.35	0/1334	1.23	8/1816 (0.4%)
All	All	0.35	0/5280	1.26	34/7196 (0.5%)

There are no bond length outliers.

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	190	ARG	CD-NE-CZ	25.28	158.99	123.60
1	B	190	ARG	CD-NE-CZ	22.05	154.47	123.60
1	C	190	ARG	CD-NE-CZ	20.53	152.34	123.60
1	D	190	ARG	CD-NE-CZ	19.36	150.70	123.60
1	B	33	ARG	CD-NE-CZ	15.75	145.65	123.60
1	A	62	ARG	NE-CZ-NH2	15.30	127.95	120.30
1	C	62	ARG	NE-CZ-NH1	14.79	127.69	120.30
1	D	62	ARG	NE-CZ-NH1	14.72	127.66	120.30
1	C	62	ARG	CD-NE-CZ	12.91	141.68	123.60
1	D	62	ARG	CD-NE-CZ	12.79	141.51	123.60
1	B	62	ARG	CD-NE-CZ	12.68	141.34	123.60
1	B	62	ARG	NE-CZ-NH2	12.65	126.63	120.30
1	C	190	ARG	NE-CZ-NH2	11.75	126.17	120.30
1	B	190	ARG	NE-CZ-NH1	11.02	125.81	120.30
1	A	62	ARG	CD-NE-CZ	10.61	138.45	123.60
1	D	190	ARG	NE-CZ-NH2	10.04	125.32	120.30
1	D	182	ARG	NE-CZ-NH1	9.62	125.11	120.30
1	B	33	ARG	CG-CD-NE	8.10	128.82	111.80
1	A	100	TYR	CB-CG-CD1	7.89	125.73	121.00
1	A	190	ARG	NE-CZ-NH1	7.62	124.11	120.30
1	A	62	ARG	NE-CZ-NH1	-6.88	116.86	120.30
1	B	144	ARG	NE-CZ-NH1	6.39	123.50	120.30
1	C	190	ARG	NH1-CZ-NH2	-5.87	112.95	119.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	33	ARG	NE-CZ-NH1	5.81	123.21	120.30
1	B	190	ARG	NH1-CZ-NH2	-5.80	113.02	119.40
1	C	33	ARG	NE-CZ-NH1	5.69	123.15	120.30
1	A	180	PHE	CB-CG-CD1	5.67	124.77	120.80
1	C	121	ARG	NE-CZ-NH1	5.38	122.99	120.30
1	D	124	ARG	CD-NE-CZ	5.35	131.09	123.60
1	A	100	TYR	CB-CG-CD2	-5.27	117.84	121.00
1	A	190	ARG	NH1-CZ-NH2	-5.22	113.66	119.40
1	D	133	ARG	CD-NE-CZ	5.15	130.82	123.60
1	C	34	LEU	CA-CB-CG	5.04	126.91	115.30
1	B	62	ARG	NE-CZ-NH1	-5.04	117.78	120.30

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	1288	0	1260	34	0
1	B	1288	0	1271	47	0
1	C	1285	0	1267	53	0
1	D	1305	0	1298	43	0
2	A	32	0	0	3	0
2	B	25	0	0	3	0
2	C	20	0	0	0	0
2	D	15	0	0	2	0
All	All	5258	0	5096	171	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 17.

All (171) 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:162:VAL:O	1:A:166:GLN:HG3	1.86	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:142:VAL:HG22	2:B:214:HOH:O	1.86	0.75
1:D:161:LEU:O	1:D:165:ILE:HG13	1.88	0.74
1:C:126:ALA:O	1:C:129:VAL:HB	1.88	0.74
1:B:151:ALA:O	1:B:155:ILE:HG13	1.89	0.72
1:A:58:ALA:O	1:A:61:ALA:HB3	1.90	0.70
1:A:31:LEU:O	1:A:35:LEU:HG	1.91	0.70
1:B:31:LEU:O	1:B:35:LEU:HG	1.90	0.70
1:B:78:ALA:O	1:B:104:ALA:HB3	1.92	0.70
1:A:119:VAL:O	1:A:123:ILE:HG13	1.93	0.69
1:C:160:THR:OG1	1:C:163:GLU:HG3	1.93	0.69
1:B:182:ARG:O	1:B:186:VAL:HG23	1.93	0.69
1:B:144:ARG:HB2	2:B:207:HOH:O	1.93	0.68
1:B:124:ARG:O	1:B:128:SER:HB3	1.93	0.67
1:D:111:ASP:OD2	1:D:114:VAL:HG12	1.95	0.66
1:C:28:LEU:O	1:C:32:GLN:HG3	1.96	0.66
1:B:119:VAL:O	1:B:123:ILE:HG13	1.97	0.65
1:D:74:HIS:HA	1:D:98:GLU:O	1.95	0.64
1:D:64:LYS:HD3	1:D:95:MET:SD	2.36	0.64
1:C:58:ALA:O	1:C:62:ARG:HG3	1.98	0.63
1:C:162:VAL:O	1:C:166:GLN:HG3	1.98	0.63
1:A:64:LYS:HE3	1:A:91:PHE:CE2	2.34	0.63
1:C:92:TYR:HA	1:C:95:MET:SD	2.38	0.62
1:C:182:ARG:HG3	1:C:182:ARG:O	2.00	0.61
1:C:65:SER:O	1:C:69:GLN:HG3	2.01	0.61
1:B:165:ILE:O	1:B:169:GLN:HB3	2.00	0.61
1:A:59:TYR:CE2	1:D:39:GLN:HB2	2.35	0.61
1:D:66:LYS:HA	1:D:69:GLN:OE1	2.01	0.61
1:B:66:LYS:O	1:B:69:GLN:HB2	2.01	0.61
1:C:187:LEU:O	1:C:191:LEU:HG	2.01	0.60
1:B:59:TYR:CE2	1:C:39:GLN:HB2	2.36	0.60
1:B:162:VAL:O	1:B:166:GLN:HG3	2.01	0.60
1:A:186:VAL:O	1:A:190:ARG:HG3	2.02	0.60
1:A:113:SER:HA	1:A:116:PHE:CG	2.37	0.59
1:C:124:ARG:O	1:C:128:SER:HB3	2.02	0.59
1:A:66:LYS:O	1:A:70:LEU:HG	2.04	0.58
1:C:57:ASP:HB2	1:C:139:ALA:O	2.03	0.58
1:D:54:PHE:O	1:D:135:LEU:HD12	2.03	0.58
1:B:44:ASN:OD1	1:C:42:THR:HG22	2.05	0.57
1:D:180:PHE:O	1:D:184:LEU:HG	2.03	0.57
1:A:97:LEU:HD12	1:A:98:GLU:O	2.04	0.57
1:D:121:ARG:HG3	1:D:156:TYR:OH	2.03	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:151:ALA:O	1:C:155:ILE:HG13	2.05	0.57
1:D:36:TRP:HB2	2:D:210:HOH:O	2.03	0.57
1:C:100:TYR:CE2	1:C:119:VAL:HG22	2.40	0.56
1:B:32:GLN:O	1:B:35:LEU:HB2	2.05	0.56
1:A:147:THR:OG1	1:A:176:PRO:HB3	2.05	0.56
1:B:50:TRP:CZ2	1:B:157:GLU:HG3	2.41	0.56
1:B:113:SER:O	1:B:116:PHE:HB2	2.06	0.56
1:D:104:ALA:HA	1:D:110:PHE:CZ	2.41	0.55
1:C:154:MET:SD	1:C:161:LEU:HA	2.47	0.55
1:C:159:MET:HA	1:C:163:GLU:OE1	2.07	0.54
1:A:44:ASN:OD1	1:D:42:THR:HG22	2.07	0.54
1:B:161:LEU:O	1:B:165:ILE:HG13	2.06	0.54
1:D:142:VAL:HG22	2:D:206:HOH:O	2.08	0.54
1:C:150:LEU:O	1:C:154:MET:HG3	2.07	0.54
1:A:92:TYR:O	1:A:95:MET:HB2	2.08	0.54
1:D:66:LYS:O	1:D:70:LEU:HG	2.07	0.53
1:C:49:VAL:HA	1:C:171:HIS:CD2	2.42	0.53
1:A:165:ILE:O	1:A:169:GLN:HB3	2.09	0.53
1:B:63:ASP:OD2	1:B:66:LYS:HE3	2.09	0.53
1:A:114:VAL:HA	2:A:226:HOH:O	2.08	0.53
1:D:50:TRP:CH2	1:D:157:GLU:HG3	2.44	0.53
1:C:76:VAL:HA	1:C:100:TYR:O	2.09	0.53
1:D:92:TYR:O	1:D:95:MET:HB2	2.10	0.52
1:C:66:LYS:O	1:C:69:GLN:HB2	2.09	0.52
1:D:102:ILE:HG12	1:D:115:TYR:CD2	2.45	0.52
1:C:154:MET:CE	1:C:187:LEU:HD23	2.40	0.52
1:D:33:ARG:O	1:D:37:VAL:HB	2.10	0.52
1:B:39:GLN:HG2	1:B:173:ASN:OD1	2.10	0.52
1:A:66:LYS:O	1:A:69:GLN:HB2	2.10	0.51
1:A:124:ARG:HD3	1:A:156:TYR:CD2	2.46	0.51
1:D:74:HIS:CD2	1:D:98:GLU:HB2	2.45	0.51
1:C:55:LEU:HA	1:C:136:VAL:O	2.09	0.51
1:B:181:LEU:O	1:B:184:LEU:HB2	2.10	0.51
1:A:144:ARG:HB2	2:A:215:HOH:O	2.11	0.51
1:C:50:TRP:CZ2	1:C:157:GLU:HG3	2.46	0.50
1:D:66:LYS:O	1:D:69:GLN:HB2	2.11	0.50
1:B:57:ASP:HB2	1:B:139:ALA:O	2.11	0.50
1:B:35:LEU:CD1	1:B:181:LEU:HD11	2.42	0.50
1:D:50:TRP:CD2	1:D:51:PRO:HD2	2.46	0.50
1:C:116:PHE:O	1:C:120:ALA:HB2	2.12	0.50
1:C:50:TRP:CD2	1:C:51:PRO:HD2	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:74:HIS:ND1	1:C:126:ALA:HB1	2.27	0.49
1:A:113:SER:HA	1:A:116:PHE:CD2	2.47	0.49
1:C:48:GLU:HG3	1:C:54:PHE:CZ	2.48	0.49
1:D:73:THR:OG1	1:D:133:ARG:O	2.30	0.49
1:B:43:LEU:O	1:C:42:THR:HA	2.12	0.49
1:D:113:SER:HA	1:D:116:PHE:CD1	2.48	0.49
1:C:92:TYR:O	1:C:95:MET:HB2	2.13	0.49
1:A:51:PRO:O	1:A:52:SER:HB2	2.13	0.49
1:B:113:SER:HA	1:B:116:PHE:CG	2.48	0.48
1:B:146:ALA:O	1:B:150:LEU:HG	2.13	0.48
1:A:104:ALA:HA	1:A:110:PHE:CZ	2.48	0.48
1:B:63:ASP:CG	1:B:66:LYS:HB2	2.33	0.48
1:D:45:HIS:O	1:D:56:GLY:HA2	2.13	0.48
1:B:124:ARG:HD3	1:B:156:TYR:CG	2.50	0.47
1:C:114:VAL:HG13	1:C:115:TYR:CD1	2.50	0.47
1:A:103:GLU:OE2	1:C:27:THR:HG22	2.14	0.47
1:C:181:LEU:O	1:C:184:LEU:HB2	2.15	0.47
1:D:153:LEU:O	1:D:157:GLU:HB2	2.15	0.47
1:A:46:ILE:CD1	1:A:135:LEU:HD13	2.45	0.47
1:D:115:TYR:C	1:D:118:PRO:HD2	2.35	0.46
1:B:182:ARG:HA	1:B:185:GLN:OE1	2.15	0.46
1:B:36:TRP:HB2	2:B:228:HOH:O	2.14	0.46
1:C:149:VAL:O	1:C:152:PHE:HB3	2.15	0.46
1:A:174:ILE:O	1:A:175:CYS:HB2	2.14	0.46
1:D:117:LEU:HD11	1:D:121:ARG:NH1	2.31	0.46
1:B:51:PRO:O	1:B:52:SER:HB2	2.15	0.46
1:D:84:GLN:O	1:D:86:ASP:N	2.49	0.46
1:A:88:GLY:HA2	1:A:99:TYR:CE2	2.51	0.46
1:C:48:GLU:HG3	1:C:54:PHE:CE2	2.52	0.45
1:D:178:SER:O	1:D:182:ARG:HB2	2.15	0.45
1:B:113:SER:HA	1:B:116:PHE:CD1	2.50	0.45
1:D:104:ALA:O	1:D:144:ARG:NH1	2.50	0.45
1:A:111:ASP:O	1:A:114:VAL:HG12	2.16	0.45
1:D:43:LEU:HD11	1:D:172:ARG:CZ	2.46	0.45
1:B:129:VAL:HG13	1:B:130:PRO:HD2	1.99	0.45
1:B:33:ARG:O	1:B:37:VAL:HG23	2.17	0.45
1:B:152:PHE:CE2	1:B:153:LEU:HD23	2.52	0.45
1:C:114:VAL:HG13	1:C:115:TYR:CE1	2.52	0.44
1:C:106:ASP:OD1	1:C:177:ASN:ND2	2.50	0.44
1:B:177:ASN:O	1:B:180:PHE:N	2.50	0.44
1:D:82:LYS:HA	1:D:86:ASP:CG	2.38	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:162:VAL:O	1:D:166:GLN:NE2	2.50	0.44
1:B:45:HIS:O	1:B:56:GLY:HA2	2.18	0.44
1:A:29:ALA:O	1:A:33:ARG:HB2	2.17	0.44
1:C:28:LEU:O	1:C:31:LEU:HB2	2.17	0.43
1:B:77:ASN:ND2	1:B:80:ALA:HA	2.33	0.43
1:A:50:TRP:CZ2	1:A:157:GLU:HG3	2.53	0.43
1:D:88:GLY:HA2	1:D:99:TYR:CE2	2.53	0.43
1:A:57:ASP:HB2	1:A:139:ALA:O	2.18	0.43
1:C:102:ILE:O	1:C:104:ALA:N	2.52	0.43
1:D:95:MET:CB	1:D:97:LEU:HD22	2.49	0.43
1:C:64:LYS:O	1:C:67:LEU:HB2	2.19	0.43
1:D:47:ASP:OD2	1:D:172:ARG:NH1	2.51	0.43
1:B:62:ARG:HA	1:B:91:PHE:CZ	2.54	0.43
1:C:62:ARG:HA	1:C:91:PHE:CE2	2.54	0.43
1:B:56:GLY:O	1:B:141:GLY:HA3	2.19	0.43
1:D:106:ASP:OD1	1:D:143:SER:OG	2.32	0.43
1:C:89:ALA:HB3	1:C:90:LYS:NZ	2.33	0.42
1:C:113:SER:HA	1:C:116:PHE:CG	2.54	0.42
1:C:50:TRP:CG	1:C:51:PRO:HD2	2.55	0.42
1:B:47:ASP:CB	1:B:172:ARG:HB2	2.49	0.42
1:B:124:ARG:HD3	1:B:156:TYR:CD2	2.54	0.42
1:C:54:PHE:O	1:C:135:LEU:HD12	2.19	0.42
1:C:51:PRO:O	1:C:52:SER:HB2	2.19	0.42
1:B:62:ARG:HA	1:B:91:PHE:CE2	2.54	0.42
1:C:161:LEU:HD11	1:C:184:LEU:HD13	2.02	0.41
1:B:152:PHE:CD2	1:B:153:LEU:HD23	2.55	0.41
1:A:177:ASN:HB2	2:A:211:HOH:O	2.19	0.41
1:D:122:TYR:O	1:D:126:ALA:HB2	2.20	0.41
1:D:153:LEU:O	1:D:157:GLU:N	2.51	0.41
1:C:113:SER:HA	1:C:116:PHE:CD1	2.56	0.41
1:A:113:SER:HA	1:A:116:PHE:CD1	2.56	0.41
1:B:92:TYR:CD2	1:B:97:LEU:HD21	2.56	0.41
1:B:34:LEU:HD23	1:B:165:ILE:CG2	2.50	0.41
1:D:46:ILE:HD11	1:D:135:LEU:HD13	2.03	0.41
1:C:102:ILE:HD11	1:C:119:VAL:HG21	2.03	0.41
1:D:113:SER:HA	1:D:116:PHE:CE1	2.56	0.41
1:C:73:THR:O	1:C:97:LEU:HA	2.21	0.41
1:D:154:MET:HA	1:D:159:MET:O	2.21	0.41
1:A:62:ARG:HD3	1:A:91:PHE:CE1	2.55	0.41
1:C:73:THR:O	1:C:97:LEU:HB2	2.20	0.41
1:A:117:LEU:HG	1:A:121:ARG:NH1	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:50:TRP:CD2	1:A:51:PRO:HD2	2.56	0.40
1:B:55:LEU:HD11	1:B:146:ALA:HA	2.03	0.40
1:C:64:LYS:HA	1:C:95:MET:HE1	2.04	0.40
1:D:77:ASN:ND2	1:D:85:VAL:O	2.54	0.40
1:C:165:ILE:O	1:C:169:GLN:HB3	2.21	0.40
1:B:119:VAL:HG11	1:B:148:LEU:HD13	2.04	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	167/205 (82%)	153 (92%)	13 (8%)	1 (1%)	30	36
1	B	166/205 (81%)	147 (89%)	18 (11%)	1 (1%)	30	36
1	C	167/205 (82%)	153 (92%)	14 (8%)	0	100	100
1	D	167/205 (82%)	152 (91%)	15 (9%)	0	100	100
All	All	667/820 (81%)	605 (91%)	60 (9%)	2 (0%)	46	57

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	172	ARG
1	A	132	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	131/171 (77%)	118 (90%)	13 (10%)	10	11
1	B	132/171 (77%)	118 (89%)	14 (11%)	8	9
1	C	131/171 (77%)	115 (88%)	16 (12%)	6	6
1	D	135/171 (79%)	119 (88%)	16 (12%)	6	7
All	All	529/684 (77%)	470 (89%)	59 (11%)	7	8

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	LEU
1	A	33	ARG
1	A	65	SER
1	A	66	LYS
1	A	73	THR
1	A	97	LEU
1	A	100	TYR
1	A	128	SER
1	A	150	LEU
1	A	158	ASN
1	A	159	MET
1	A	169	GLN
1	A	190	ARG
1	B	28	LEU
1	B	33	ARG
1	B	34	LEU
1	B	66	LYS
1	B	87	THR
1	B	90	LYS
1	B	96	SER
1	B	97	LEU
1	B	128	SER
1	B	144	ARG
1	B	157	GLU
1	B	166	GLN
1	B	169	GLN
1	B	190	ARG
1	C	28	LEU
1	C	33	ARG
1	C	34	LEU

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Mol	Chain	Res	Type
1	C	47	ASP
1	C	51	PRO
1	C	65	SER
1	C	66	LYS
1	C	67	LEU
1	C	90	LYS
1	C	124	ARG
1	C	128	SER
1	C	157	GLU
1	C	159	MET
1	C	169	GLN
1	C	172	ARG
1	C	182	ARG
1	D	28	LEU
1	D	32	GLN
1	D	66	LYS
1	D	73	THR
1	D	82	LYS
1	D	107	ASN
1	D	114	VAL
1	D	128	SER
1	D	133	ARG
1	D	143	SER
1	D	157	GLU
1	D	158	ASN
1	D	159	MET
1	D	169	GLN
1	D	182	ARG
1	D	190	ARG

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

Mol	Chain	Res	Type
1	A	74	HIS

5.3.3 RNA ⓘ

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](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	169/205 (82%)	0.45	7 (4%) 41 50	20, 33, 57, 87	0
1	B	168/205 (81%)	0.57	7 (4%) 40 49	21, 34, 59, 85	0
1	C	169/205 (82%)	0.61	17 (10%) 9 13	20, 34, 59, 67	0
1	D	169/205 (82%)	0.49	8 (4%) 35 44	21, 34, 58, 70	0
All	All	675/820 (82%)	0.53	39 (5%) 26 35	20, 34, 59, 87	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	129	VAL	4.6
1	C	28	LEU	3.9
1	D	36	TRP	3.3
1	C	33	ARG	3.2
1	B	37	VAL	2.9
1	D	129	VAL	2.9
1	B	192	GLY	2.9
1	B	36	TRP	2.9
1	B	25	PRO	2.8
1	C	49	VAL	2.8
1	C	31	LEU	2.8
1	A	129	VAL	2.8
1	C	155	ILE	2.8
1	C	76	VAL	2.7
1	D	75	VAL	2.7
1	C	97	LEU	2.6
1	A	37	VAL	2.5
1	D	97	LEU	2.5
1	B	31	LEU	2.5
1	C	35	LEU	2.4
1	B	129	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	171	HIS	2.4
1	C	192	GLY	2.4
1	C	25	PRO	2.4
1	A	41	ALA	2.4
1	C	161	LEU	2.3
1	B	117	LEU	2.3
1	C	36	TRP	2.3
1	A	130	PRO	2.2
1	D	132	GLY	2.2
1	D	130	PRO	2.2
1	A	97	LEU	2.2
1	C	131	GLN	2.1
1	A	83	PHE	2.1
1	C	83	PHE	2.1
1	C	130	PRO	2.1
1	C	181	LEU	2.0
1	A	31	LEU	2.0
1	D	117	LEU	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 carbohydrates 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.