



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:11 PM UTC

PDB ID : 8FHU / pdb_00008fhu
Title : Structure of Pyruvate dehydrogenase phosphatase regulatory subunit epitope presented by H2-Dd
Authors : Custodio, J.M.F.; Baker, B.M.
Deposited on : 2022-12-15
Resolution : 1.80 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	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

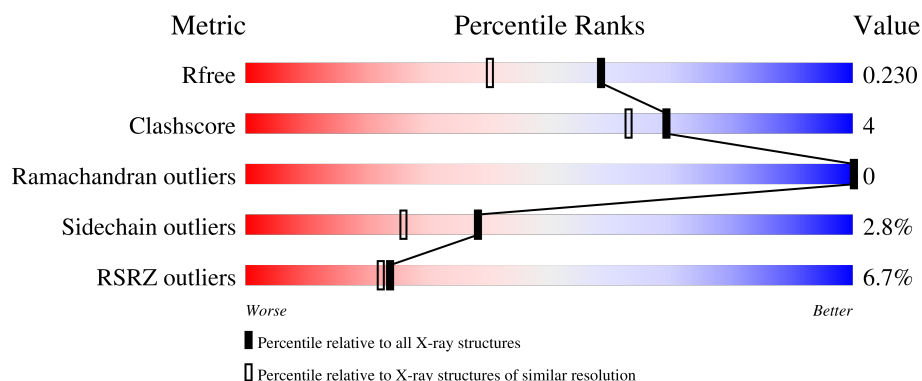
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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}	180053	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	277	3% 86% 10% ..
1	D	277	9% 81% 12% ...
2	B	99	14% 84% 12% ..
2	E	99	3% 85% 12% ...
3	C	9	89% 11%

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Mol	Chain	Length	Quality of chain
3	F	9	 78% 22%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6327 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 H-2 class I histocompatibility antigen, D-D alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	271	Total	C	N	O	S	0	0	0
			2218	1395	399	415	9			
1	D	267	Total	C	N	O	S	0	0	0
			2186	1375	398	404	9			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP P01900
D	1	MET	-	initiating methionine	UNP P01900

- Molecule 2 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	97	Total	C	N	O	S	0	0	0
			799	509	134	149	7			
2	E	98	Total	C	N	O	S	0	0	0
			806	515	137	147	7			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	85	ASP	ALA	variant	UNP P01887
E	85	ASP	ALA	variant	UNP P01887

- Molecule 3 is a protein called Pyruvate dehydrogenase phosphatase regulatory subunit, mitochondrial.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	9	Total	C	N	O	0	0	0
			66	42	12	12			

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	F	9	Total	C	N	O	0	0	0
			66	42	12	12			

- Molecule 4 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	K	0	0
			1	1		
4	D	1	Total	K	0	0
			1	1		

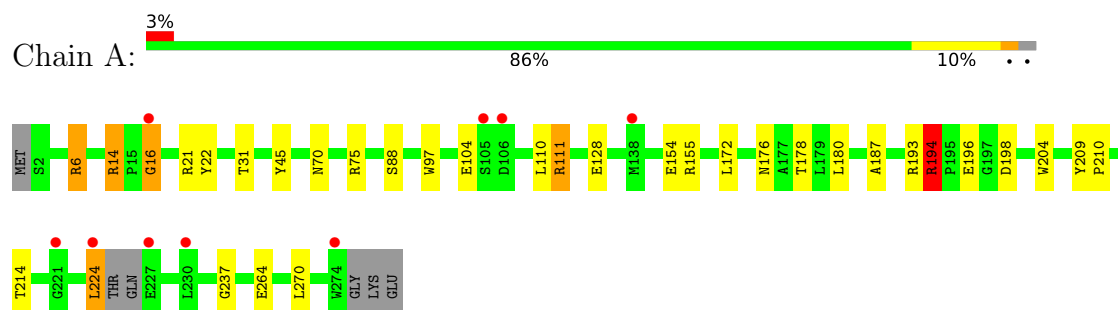
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	79	Total	O	0	0
			79	79		
5	B	22	Total	O	0	0
			22	22		
5	C	6	Total	O	0	0
			6	6		
5	D	55	Total	O	0	0
			55	55		
5	E	21	Total	O	0	0
			21	21		
5	F	1	Total	O	0	0
			1	1		

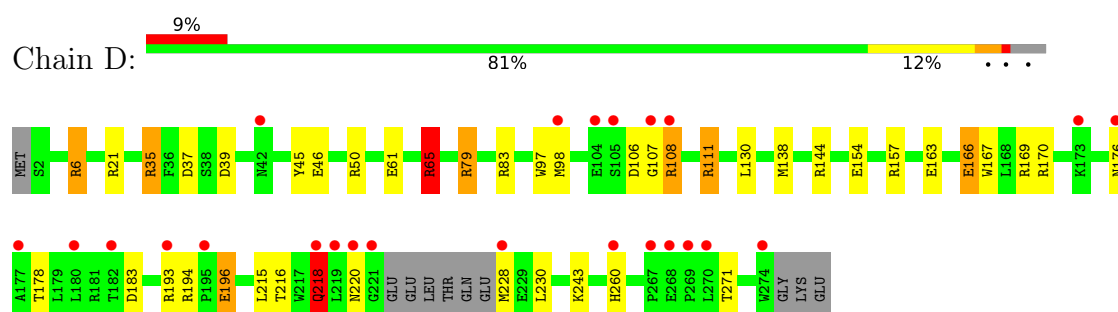
3 Residue-property plots

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.

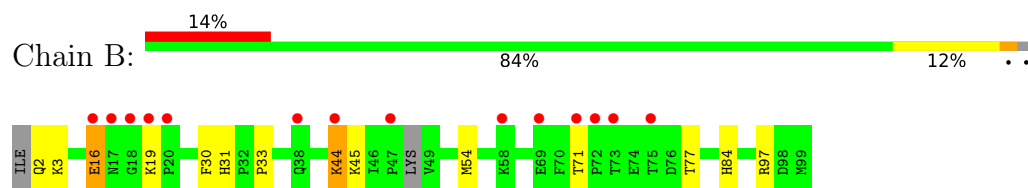
- Molecule 1: H-2 class I histocompatibility antigen, D-D alpha chain



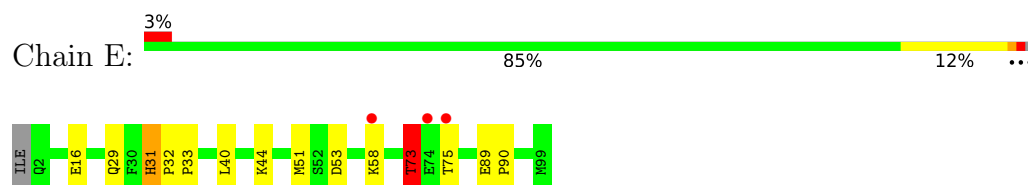
- Molecule 1: H-2 class I histocompatibility antigen, D-D alpha chain



- Molecule 2: Beta-2-microglobulin



- Molecule 2: Beta-2-microglobulin



- Molecule 3: Pyruvate dehydrogenase phosphatase regulatory subunit, mitochondrial

Chain C:  89% 11%



- Molecule 3: Pyruvate dehydrogenase phosphatase regulatory subunit, mitochondrial

Chain F:  78% 22%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	56.41Å 104.56Å 169.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.82 – 1.80 49.82 – 1.80	Depositor EDS
% Data completeness (in resolution range)	73.7 (49.82-1.80) 73.8 (49.82-1.80)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.14 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.8.0403	Depositor
R, R_{free}	0.208 , 0.249 (Not available) , 0.230	Depositor DCC
R_{free} test set	3480 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	20.1	Xtriage
Anisotropy	0.204	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 29.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6327	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.74% 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

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

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	0.80	0/2280	1.41	14/3096 (0.5%)
1	D	0.76	0/2248	1.41	13/3053 (0.4%)
2	B	0.84	1/824 (0.1%)	1.50	10/1119 (0.9%)
2	E	0.82	0/832	1.42	5/1129 (0.4%)
3	C	0.78	0/66	1.29	0/88
3	F	0.88	0/66	1.20	0/88
All	All	0.80	1/6316 (0.0%)	1.42	42/8573 (0.5%)

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	6
1	D	0	11
2	B	0	1
All	All	0	18

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	31	HIS	CE1-NE2	6.64	1.39	1.32

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	111	ARG	CB-CG-CD	9.41	132.96	111.30
2	B	77	THR	CA-CB-OG1	-8.41	96.98	109.60
1	A	75	ARG	NE-CZ-NH1	8.17	129.67	121.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	71	THR	CA-CB-OG1	-7.71	98.04	109.60
1	A	75	ARG	NE-CZ-NH2	-7.64	112.32	119.20
1	D	37	ASP	CA-CB-CG	7.39	119.99	112.60
1	D	178	THR	CA-CB-OG1	-7.34	98.59	109.60
1	D	154	GLU	CB-CG-CD	7.30	125.01	112.60
1	A	6	ARG	CB-CG-CD	7.11	127.65	111.30
1	D	83	ARG	CB-CG-CD	7.04	127.48	111.30
1	D	176	ASN	CA-CB-CG	-7.00	105.60	112.60
2	B	45	LYS	CB-CA-C	6.44	120.19	109.89
1	A	178	THR	CA-CB-OG1	-6.39	100.01	109.60
1	D	218	GLN	CB-CA-C	6.35	121.33	109.71
1	A	16	GLY	N-CA-C	-6.25	98.37	113.18
1	A	75	ARG	CD-NE-CZ	6.18	133.06	124.40
2	B	16	GLU	CA-C-O	-5.98	114.06	120.46
2	B	44	LYS	CB-CG-CD	5.78	124.60	111.30
1	D	218	GLN	CB-CG-CD	5.75	122.37	112.60
1	A	270	LEU	N-CA-CB	-5.74	101.05	110.52
2	E	51	MET	N-CA-C	-5.65	99.08	108.34
2	E	16	GLU	CB-CA-C	5.62	118.80	110.14
2	B	84	HIS	CB-CG-CD2	5.59	138.47	131.20
2	B	30	PHE	CA-CB-CG	5.54	119.34	113.80
1	D	79	ARG	CD-NE-CZ	5.53	132.14	124.40
1	A	196	GLU	CB-CG-CD	5.53	122.00	112.60
1	A	154	GLU	CB-CG-CD	5.50	121.96	112.60
2	E	29	GLN	CB-CG-CD	-5.49	103.28	112.60
2	B	44	LYS	CB-CA-C	5.47	119.38	109.37
2	B	84	HIS	CB-CG-ND1	-5.47	114.50	122.70
2	B	71	THR	CB-CA-C	5.47	119.07	110.71
1	A	14	ARG	CB-CG-CD	5.41	123.75	111.30
1	D	35	ARG	CG-CD-NE	5.37	123.82	112.00
1	A	214	THR	CA-CB-OG1	-5.35	101.58	109.60
1	A	194	ARG	CB-CG-CD	-5.32	99.06	111.30
1	D	79	ARG	CB-CG-CD	5.29	123.47	111.30
1	D	79	ARG	NE-CZ-NH2	-5.19	114.53	119.20
1	D	183	ASP	CB-CA-C	5.16	116.08	110.15
2	E	73	THR	CA-CB-OG1	-5.15	101.88	109.60
2	E	31	HIS	CB-CG-ND1	-5.15	114.98	122.70
1	D	166	GLU	N-CA-CB	5.09	117.39	110.01
1	A	155	ARG	CG-CD-NE	5.05	123.11	112.00

There are no chirality outliers.

All (18) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	14	ARG	Sidechain
1	A	16	GLY	Mainchain
1	A	193	ARG	Sidechain
1	A	194	ARG	Sidechain
1	A	21	ARG	Sidechain
1	A	6	ARG	Sidechain
2	B	97	ARG	Sidechain
1	D	108	ARG	Sidechain
1	D	111	ARG	Sidechain
1	D	144	ARG	Sidechain
1	D	157	ARG	Sidechain
1	D	170	ARG	Sidechain
1	D	193	ARG	Sidechain
1	D	194	ARG	Sidechain
1	D	50	ARG	Sidechain
1	D	6	ARG	Sidechain
1	D	65	ARG	Sidechain
1	D	79	ARG	Sidechain

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	2218	0	2078	14	0
1	D	2186	0	2053	17	0
2	B	799	0	757	3	0
2	E	806	0	773	7	0
3	C	66	0	74	1	0
3	F	66	0	74	3	0
4	A	1	0	0	0	0
4	D	1	0	0	0	0
5	A	79	0	0	0	0
5	B	22	0	0	1	0
5	C	6	0	0	0	0
5	D	55	0	0	1	0
5	E	21	0	0	0	0
5	F	1	0	0	0	0
All	All	6327	0	5809	40	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 4.

All (40) 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:THR:HG21	1:A:209:TYR:OH	1.62	0.99
1:D:196:GLU:N	1:D:196:GLU:OE1	2.03	0.92
2:E:73:THR:HG22	2:E:75:THR:H	1.49	0.77
1:D:106:ASP:CB	1:D:108:ARG:HH21	1.98	0.77
1:A:111:ARG:HD3	1:A:128:GLU:HG3	1.74	0.68
1:A:194:ARG:HD3	1:A:198:ASP:OD1	1.93	0.68
1:D:111:ARG:NH2	5:D:401:HOH:O	2.29	0.65
1:D:216:THR:HG22	1:D:260:HIS:HB2	1.81	0.63
1:D:216:THR:CG2	1:D:260:HIS:HB2	2.28	0.63
1:A:176:ASN:HB2	1:A:180:LEU:HD13	1.82	0.61
1:D:35:ARG:CZ	2:E:53:ASP:HB2	2.35	0.56
1:D:163:GLU:HB3	3:F:1:ILE:HD11	1.87	0.56
1:A:97:TRP:CH2	3:C:3:PRO:HG2	2.40	0.56
1:A:172:LEU:O	1:A:176:ASN:HB3	2.09	0.53
1:D:97:TRP:CH2	3:F:3:PRO:HG2	2.45	0.51
2:B:16:GLU:HB2	2:B:19:LYS:CB	2.40	0.51
1:A:224:LEU:C	1:A:224:LEU:HD12	2.37	0.50
1:D:230:LEU:HD11	1:D:243:LYS:HE3	1.94	0.49
1:A:172:LEU:HD22	1:A:180:LEU:HD11	1.96	0.48
2:E:32:PRO:HB2	2:E:33:PRO:HD2	1.95	0.47
1:A:104:GLU:HG2	1:A:110:LEU:HD12	1.97	0.47
1:D:218:GLN:HG3	1:D:220:ASN:CG	2.40	0.46
1:A:22:TYR:OH	1:A:70:ASN:HB3	2.17	0.45
1:D:107:GLY:HA3	1:D:169:ARG:HH21	1.81	0.45
2:E:73:THR:HG22	2:E:75:THR:N	2.25	0.45
1:D:6:ARG:HD2	1:D:98:MET:SD	2.57	0.44
2:E:40:LEU:HA	2:E:44:LYS:O	2.18	0.44
1:D:61:GLU:HG3	1:D:65:ARG:NH2	2.34	0.42
1:D:228:MET:HE3	1:D:230:LEU:HD13	2.01	0.42
1:A:237:GLY:HA3	5:B:106:HOH:O	2.19	0.42
2:E:31:HIS:ND1	2:E:32:PRO:HA	2.35	0.42
2:B:33:PRO:O	2:B:54:MET:HE1	2.19	0.42
2:B:16:GLU:HA	2:B:16:GLU:OE2	2.20	0.41
1:A:224:LEU:C	1:A:224:LEU:CD1	2.93	0.41
1:A:210:PRO:HG2	1:A:264:GLU:OE2	2.19	0.41
2:E:89:GLU:HG2	2:E:90:PRO:HD2	2.02	0.41
1:D:65:ARG:CZ	1:D:65:ARG:HB2	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:187:ALA:HA	1:A:204:TRP:O	2.21	0.40
1:D:167:TRP:CD1	3:F:1:ILE:HG13	2.56	0.40
1:D:21:ARG:HD2	1:D:39:ASP:OD2	2.21	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	267/277 (96%)	262 (98%)	5 (2%)	0	100	100
1	D	263/277 (95%)	257 (98%)	6 (2%)	0	100	100
2	B	93/99 (94%)	92 (99%)	1 (1%)	0	100	100
2	E	96/99 (97%)	93 (97%)	3 (3%)	0	100	100
3	C	7/9 (78%)	7 (100%)	0	0	100	100
3	F	7/9 (78%)	7 (100%)	0	0	100	100
All	All	733/770 (95%)	718 (98%)	15 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	225/232 (97%)	222 (99%)	3 (1%)	61	54
1	D	221/232 (95%)	211 (96%)	10 (4%)	24	12
2	B	91/94 (97%)	88 (97%)	3 (3%)	33	21
2	E	91/94 (97%)	89 (98%)	2 (2%)	45	34
3	C	7/7 (100%)	7 (100%)	0	100	100
3	F	7/7 (100%)	7 (100%)	0	100	100
All	All	642/666 (96%)	624 (97%)	18 (3%)	38	26

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	45	TYR
1	A	88	SER
1	A	224	LEU
2	B	2	GLN
2	B	3	LYS
2	B	44	LYS
1	D	45	TYR
1	D	46	GLU
1	D	65	ARG
1	D	130	LEU
1	D	138	MET
1	D	166	GLU
1	D	196	GLU
1	D	215	LEU
1	D	218	GLN
1	D	271	THR
2	E	58	LYS
2	E	73	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	ASN
1	A	115	GLN
1	A	255	GLN
2	E	6	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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

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	271/277 (97%)	0.02	9 (3%) 49 49	12, 23, 42, 70	0
1	D	267/277 (96%)	0.49	24 (8%) 15 13	13, 31, 60, 77	0
2	B	97/99 (97%)	0.71	14 (14%) 6 5	16, 35, 66, 77	0
2	E	98/99 (98%)	0.18	3 (3%) 51 51	16, 27, 47, 55	0
3	C	9/9 (100%)	-0.23	0 100 100	14, 17, 21, 28	0
3	F	9/9 (100%)	-0.21	0 100 100	22, 24, 27, 36	0
All	All	751/770 (97%)	0.29	50 (6%) 24 22	12, 27, 55, 77	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	224	LEU	4.5
1	D	267	PRO	3.9
1	D	195	PRO	3.7
2	B	17	ASN	3.5
1	D	108	ARG	3.5
2	B	44	LYS	3.5
2	B	18	GLY	3.4
1	A	16	GLY	3.4
2	B	20	PRO	3.2
2	B	47	PRO	3.2
2	B	19	LYS	3.1
2	E	75	THR	3.1
1	D	269	PRO	3.0
1	A	105	SER	2.9
1	D	42	ASN	2.9
1	D	268	GLU	2.9
1	D	107	GLY	2.9
1	D	177	ALA	2.9
1	D	260	HIS	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	173	LYS	2.8
2	E	58	LYS	2.8
2	B	73	THR	2.8
1	D	228	MET	2.6
1	D	218	GLN	2.6
1	D	220	ASN	2.6
2	B	75	THR	2.6
1	A	106	ASP	2.6
1	D	105	SER	2.4
1	D	219	LEU	2.4
1	A	230	LEU	2.4
1	D	180	LEU	2.3
1	D	270	LEU	2.3
1	D	98	MET	2.3
1	D	193	ARG	2.2
2	E	74	GLU	2.2
2	B	58	LYS	2.2
1	D	104	GLU	2.2
2	B	71	THR	2.2
2	B	72	PRO	2.2
1	D	274	TRP	2.1
1	A	138	MET	2.1
2	B	16	GLU	2.1
2	B	38	GLN	2.1
1	A	274	TRP	2.1
2	B	69	GLU	2.1
1	A	221	GLY	2.1
1	D	176	ASN	2.1
1	D	182	THR	2.0
1	D	221	GLY	2.0
1	A	227	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 oligosaccharides 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. 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	B-factors(Å ²)	Q<0.9
4	K	D	301	1/1	0.90	0.31	80,80,80,80	0
4	K	A	301	1/1	0.97	0.30	39,39,39,39	0

6.5 Other polymers [i](#)

There are no such residues in this entry.