



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:33 PM UTC

PDB ID : 4GUP / pdb_00004gup
Title : Structure of MHC-class I related molecule MR1
Authors : Patel, O.; Le Nours, J.; Rossjohn, J.
Deposited on : 2012-08-29
Resolution : 3.20 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

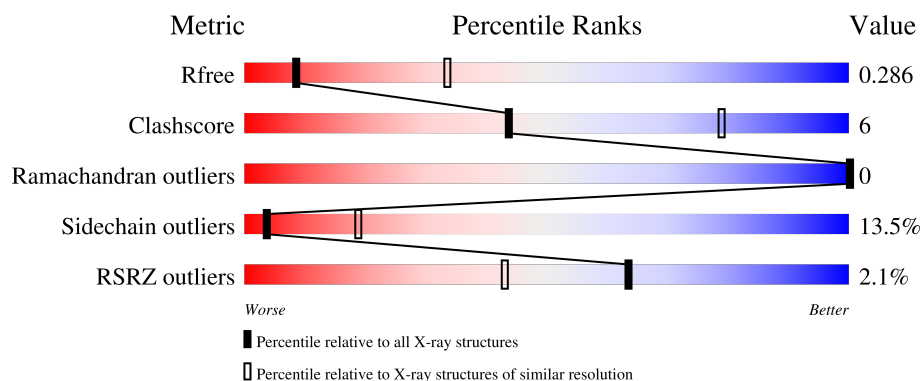
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	271	69% 27% ..
1	C	271	3% 66% 25% . 5%
2	B	99	76% 19% ..
2	D	99	3% 81% 16% ..

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5794 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 Major histocompatibility complex class I-related gene protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	265	Total	C	N	O	S	0	0	0
			2150	1379	363	398	10			
1	C	258	Total	C	N	O	S	0	0	0
			2089	1341	355	383	10			

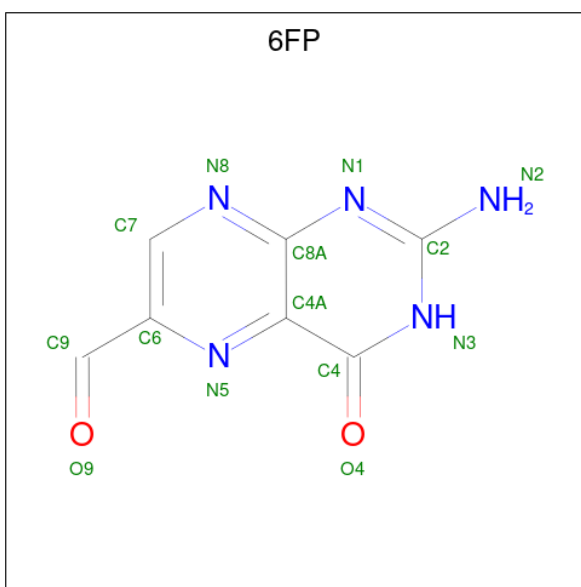
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	expression tag	UNP Q95460
A	261	SER	CYS	engineered mutation	UNP Q95460
C	0	MET	-	expression tag	UNP Q95460
C	261	SER	CYS	engineered mutation	UNP Q95460

- 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
			759	486	130	141	2			
2	D	98	Total	C	N	O	S	0	0	0
			764	488	134	140	2			

- Molecule 3 is 2-amino-4-oxo-3,4-dihydropteridine-6-carbaldehyde (CCD ID: 6FP) (formula: $C_7H_5N_5O_2$).

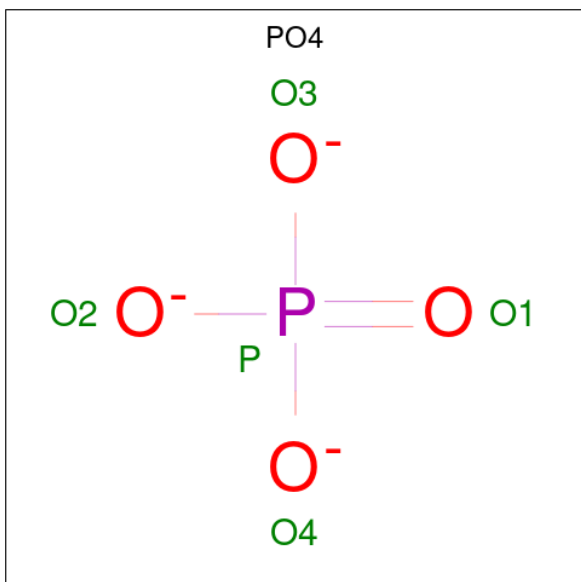


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			13	7	5	1		
3	C	1	Total	C	N	O	0	0
			13	7	5	1		

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Cl	0	0
			1	1		

- Molecule 5 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).

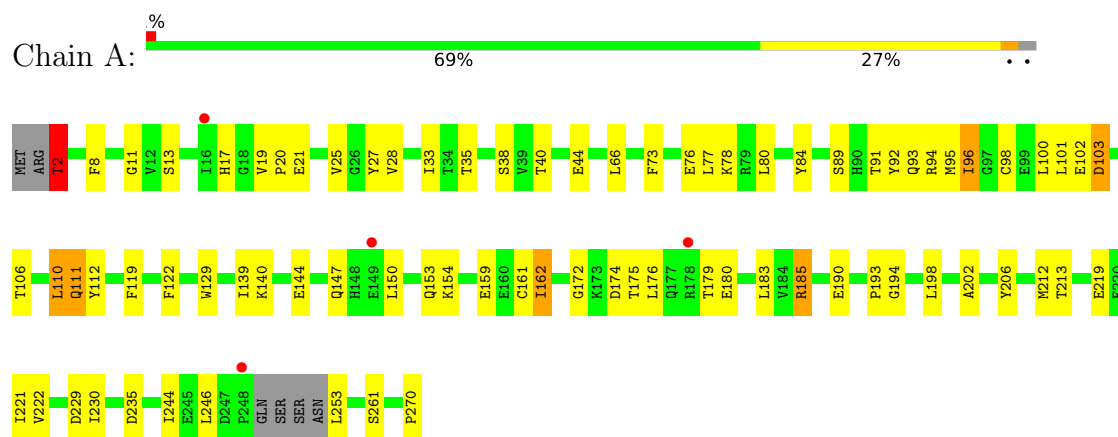


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	P	0	0
			5	4	1		

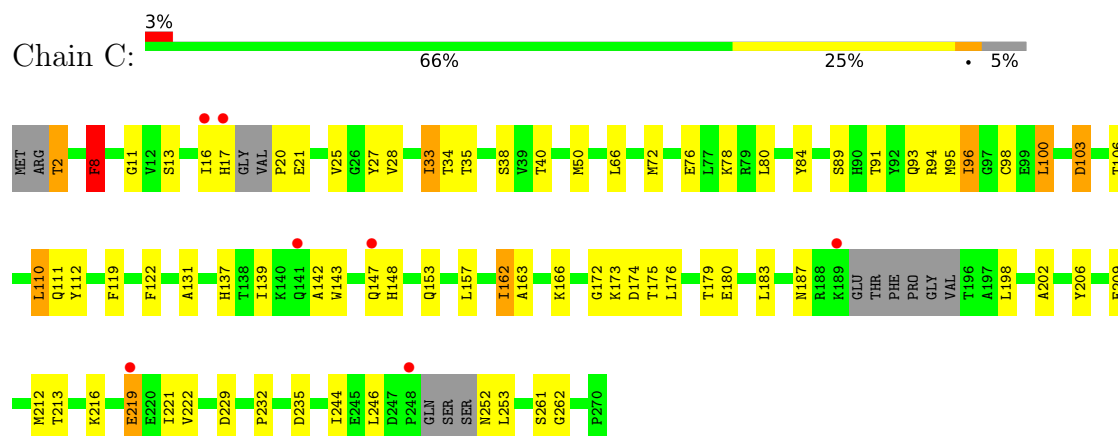
3 Residue-property plots [i](#)

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.

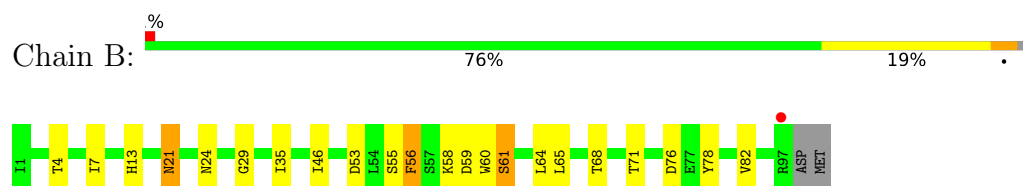
- Molecule 1: Major histocompatibility complex class I-related gene protein



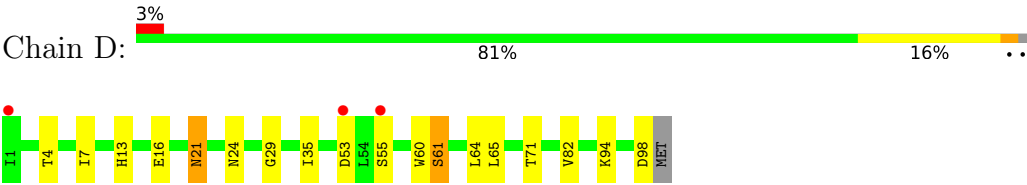
- Molecule 1: Major histocompatibility complex class I-related gene protein



- Molecule 2: Beta-2-microglobulin



- Molecule 2: Beta-2-microglobulin



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	59.15Å 89.78Å 171.34Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	85.67 – 3.20 85.67 – 3.20	Depositor EDS
% Data completeness (in resolution range)	97.7 (85.67-3.20) 97.4 (85.67-3.20)	Depositor EDS
R_{merge}	0.27	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.94 (at 3.19Å)	Xtriage
Refinement program	BUSTER 2.10.0	Depositor
R, R_{free}	0.193 , 0.260 0.213 , 0.286	Depositor DCC
R_{free} test set	764 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å ²)	42.3	Xtriage
Anisotropy	0.334	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 73.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	5794	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.62% 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: CL, 6FP, 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	0.89	1/2218 (0.0%)	1.39	19/3023 (0.6%)
1	C	0.91	0/2153	1.43	27/2932 (0.9%)
2	B	0.80	0/782	1.22	2/1070 (0.2%)
2	D	0.84	0/786	1.18	1/1075 (0.1%)
All	All	0.88	1/5939 (0.0%)	1.36	49/8100 (0.6%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	95	MET	SD-CE	5.28	1.92	1.79

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	8	PHE	CA-CB-CG	9.44	123.24	113.80
2	B	56	PHE	CA-CB-CG	8.20	122.00	113.80
1	A	78	LYS	CA-C-N	6.86	129.36	120.44
1	A	78	LYS	C-N-CA	6.86	129.36	120.44
1	C	28	VAL	N-CA-CB	6.79	119.90	111.41
1	C	179	THR	CA-C-N	6.72	129.65	122.26
1	C	179	THR	C-N-CA	6.72	129.65	122.26
1	C	147	GLN	N-CA-C	6.70	118.24	111.07
1	C	78	LYS	CA-C-N	6.66	129.09	120.44
1	C	78	LYS	C-N-CA	6.66	129.09	120.44
1	A	28	VAL	N-CA-CB	6.55	119.60	111.41
1	C	142	ALA	CA-C-N	6.34	132.54	122.59
1	C	142	ALA	C-N-CA	6.34	132.54	122.59
1	C	235	ASP	CA-CB-CG	6.05	118.65	112.60
1	A	179	THR	CA-C-N	5.99	128.85	122.26
1	A	179	THR	C-N-CA	5.99	128.85	122.26

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	71	THR	CB-CA-C	5.96	116.41	110.33
1	C	209	GLU	CB-CG-CD	5.81	122.47	112.60
1	C	229	ASP	CA-CB-CG	5.71	118.31	112.60
2	D	71	THR	CB-CA-C	5.68	116.13	110.33
1	C	142	ALA	N-CA-C	5.66	117.72	108.55
1	A	235	ASP	CA-CB-CG	5.64	118.24	112.60
1	A	172	GLY	CA-C-N	5.60	131.49	122.14
1	A	172	GLY	C-N-CA	5.60	131.49	122.14
1	C	172	GLY	CA-C-N	5.54	131.39	122.14
1	C	172	GLY	C-N-CA	5.54	131.39	122.14
1	A	103	ASP	CA-CB-CG	5.53	118.13	112.60
1	A	229	ASP	CA-CB-CG	5.50	118.11	112.60
1	A	230	ILE	N-CA-CB	5.40	116.42	110.53
1	A	2	THR	CA-C-N	5.29	129.80	120.87
1	A	2	THR	C-N-CA	5.29	129.80	120.87
1	C	219	GLU	CB-CG-CD	5.23	121.49	112.60
1	C	2	THR	CA-C-N	5.23	128.64	120.75
1	C	2	THR	C-N-CA	5.23	128.64	120.75
1	A	221	ILE	CA-C-N	5.21	129.17	120.62
1	A	221	ILE	C-N-CA	5.21	129.17	120.62
1	C	84	TYR	CA-C-N	-5.18	114.50	123.25
1	C	84	TYR	C-N-CA	-5.18	114.50	123.25
1	C	103	ASP	CA-CB-CG	5.14	117.74	112.60
1	C	157	LEU	CA-C-N	5.11	127.63	120.28
1	C	157	LEU	C-N-CA	5.11	127.63	120.28
1	C	163	ALA	CA-C-N	5.03	127.02	120.28
1	C	163	ALA	C-N-CA	5.03	127.02	120.28
1	A	84	TYR	CA-C-N	-5.01	114.78	123.25
1	A	84	TYR	C-N-CA	-5.01	114.78	123.25
1	A	154	LYS	CA-C-N	5.01	127.00	120.28
1	A	154	LYS	C-N-CA	5.01	127.00	120.28
1	C	221	ILE	CA-C-N	5.01	128.83	120.62
1	C	221	ILE	C-N-CA	5.01	128.83	120.62

There are no chirality outliers.

There are no planarity outliers.

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	2150	0	1997	30	0
1	C	2089	0	1938	28	0
2	B	759	0	682	11	0
2	D	764	0	695	7	0
3	A	13	0	4	1	0
3	C	13	0	4	1	0
4	A	1	0	0	0	0
5	A	5	0	0	0	0
All	All	5794	0	5320	65	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 6.

All (65) 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:C:202:ALA:HB3	1:C:212:MET:HE1	1.61	0.83
1:A:202:ALA:HB3	1:A:212:MET:HE1	1.63	0.81
1:A:13:SER:HB2	1:A:89:SER:HA	1.65	0.79
1:C:13:SER:HB2	1:C:89:SER:HA	1.65	0.77
1:A:270:PRO:HB3	1:C:262:GLY:HA3	1.70	0.72
1:A:98:CYS:SG	1:A:162:ILE:HD12	2.31	0.70
1:C:98:CYS:SG	1:C:162:ILE:HD12	2.31	0.70
1:A:77:LEU:HD23	1:A:92:TYR:HB2	1.74	0.68
1:C:20:PRO:HB2	1:C:38:SER:OG	1.99	0.62
1:A:20:PRO:HB2	1:A:38:SER:OG	2.00	0.62
2:B:24:ASN:HB3	2:B:65:LEU:HD11	1.81	0.62
2:D:24:ASN:HB3	2:D:65:LEU:HD11	1.83	0.60
1:C:94:ARG:HH12	3:C:301:6FP:H4	1.49	0.59
1:A:110:LEU:HB3	1:A:122:PHE:HB3	1.87	0.57
1:A:94:ARG:HH12	3:A:301:6FP:H4	1.50	0.57
1:A:98:CYS:HG	1:A:161:CYS:HG	0.64	0.57
1:C:33:ILE:HG23	1:C:34:THR:HG23	1.87	0.57
1:A:111:GLN:HB3	2:B:60:TRP:HH2	1.71	0.56
1:C:110:LEU:HB3	1:C:122:PHE:HB3	1.87	0.56
1:A:11:GLY:HA2	1:A:21:GLU:O	2.05	0.56
1:C:11:GLY:HA2	1:C:21:GLU:O	2.07	0.55
1:A:111:GLN:HB3	2:B:60:TRP:CH2	2.43	0.54
1:A:129:TRP:HB2	1:A:140:LYS:HG3	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:185:ARG:HD3	1:C:183:LEU:HD21	1.92	0.52
2:D:7:ILE:HG12	2:D:82:VAL:HG21	1.90	0.52
2:D:29:GLY:HA2	2:D:61:SER:HB2	1.93	0.50
2:B:7:ILE:HG12	2:B:82:VAL:HG21	1.93	0.50
1:A:8:PHE:HB2	2:B:56:PHE:CZ	2.47	0.50
1:C:180:GLU:HB3	1:C:206:TYR:HB3	1.94	0.49
1:A:180:GLU:HB3	1:A:206:TYR:HB3	1.95	0.49
2:D:13:HIS:HB2	2:D:21:ASN:HD21	1.79	0.48
1:C:93:GLN:OE1	2:D:60:TRP:HB3	2.14	0.48
1:C:106:THR:HB	1:C:162:ILE:HD13	1.95	0.48
2:B:29:GLY:HA2	2:B:61:SER:HB2	1.96	0.47
1:C:94:ARG:HG3	1:C:112:TYR:CE2	2.49	0.47
1:C:131:ALA:HB3	1:C:137:HIS:CE1	2.50	0.47
1:A:94:ARG:HG3	1:A:112:TYR:CE2	2.49	0.47
2:B:13:HIS:HB2	2:B:21:ASN:HD21	1.79	0.47
1:C:33:ILE:HG12	1:C:50:MET:SD	2.55	0.47
1:A:106:THR:HB	1:A:162:ILE:HD13	1.96	0.46
1:C:80:LEU:HD21	1:C:119:PHE:CZ	2.51	0.46
1:A:183:LEU:HB3	1:C:183:LEU:HD22	1.97	0.45
2:B:46:ILE:HG21	2:B:68:THR:HG21	1.99	0.45
1:C:8:PHE:HD1	1:C:95:MET:HG3	1.80	0.45
1:A:80:LEU:HD21	1:A:119:PHE:CZ	2.51	0.45
1:C:198:LEU:HB2	1:C:244:ILE:HG22	1.99	0.45
1:C:25:VAL:HG22	1:C:35:THR:HG22	1.99	0.44
1:A:93:GLN:OE1	2:B:60:TRP:HB3	2.17	0.44
1:A:25:VAL:HG22	1:A:35:THR:HG22	1.99	0.44
1:A:144:GLU:HA	1:A:150:LEU:HD11	1.99	0.44
1:C:33:ILE:HG23	1:C:50:MET:HE2	2.00	0.43
1:A:110:LEU:HD13	1:A:129:TRP:CH2	2.54	0.43
1:A:27:TYR:OH	2:B:55:SER:HB2	2.18	0.43
1:A:96:ILE:HG23	1:A:110:LEU:HG	2.01	0.42
1:C:27:TYR:OH	2:D:55:SER:HB2	2.19	0.42
1:C:94:ARG:HG3	1:C:112:TYR:CZ	2.54	0.42
1:C:96:ILE:HG23	1:C:110:LEU:HG	2.00	0.42
1:C:8:PHE:CD1	1:C:95:MET:HG3	2.54	0.42
1:A:198:LEU:HB2	1:A:244:ILE:HG22	2.01	0.41
1:C:232:PRO:HG2	2:D:65:LEU:HD22	2.02	0.41
1:A:19:VAL:CB	1:A:20:PRO:HA	2.51	0.41
1:A:193:PRO:HA	1:A:194:GLY:HA2	1.79	0.41
2:B:76:ASP:HB3	2:B:78:TYR:CE2	2.56	0.41
1:C:100:LEU:HD11	1:C:166:LYS:HE3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2:THR:HG23	1:A:101:LEU:HA	2.03	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	261/271 (96%)	239 (92%)	22 (8%)	0	100	100
1	C	250/271 (92%)	232 (93%)	18 (7%)	0	100	100
2	B	95/99 (96%)	92 (97%)	3 (3%)	0	100	100
2	D	96/99 (97%)	94 (98%)	2 (2%)	0	100	100
All	All	702/740 (95%)	657 (94%)	45 (6%)	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	225/241 (93%)	194 (86%)	31 (14%)	3	18
1	C	218/241 (90%)	185 (85%)	33 (15%)	3	14
2	B	79/94 (84%)	71 (90%)	8 (10%)	7	30
2	D	79/94 (84%)	70 (89%)	9 (11%)	5	24

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	601/670 (90%)	520 (86%)	81 (14%)	4 19

All (81) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	THR
1	A	17	HIS
1	A	33	ILE
1	A	40	THR
1	A	44	GLU
1	A	66	LEU
1	A	73	PHE
1	A	76	GLU
1	A	91	THR
1	A	96	ILE
1	A	100	LEU
1	A	102	GLU
1	A	103	ASP
1	A	110	LEU
1	A	111	GLN
1	A	139	ILE
1	A	147	GLN
1	A	153	GLN
1	A	159	GLU
1	A	162	ILE
1	A	174	ASP
1	A	175	THR
1	A	176	LEU
1	A	185	ARG
1	A	190	GLU
1	A	213	THR
1	A	219	GLU
1	A	222	VAL
1	A	246	LEU
1	A	253	LEU
1	A	261	SER
2	B	4	THR
2	B	21	ASN
2	B	35	ILE
2	B	53	ASP
2	B	58	LYS
2	B	59	ASP

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Mol	Chain	Res	Type
2	B	61	SER
2	B	64	LEU
1	C	2	THR
1	C	8	PHE
1	C	16	ILE
1	C	17	HIS
1	C	33	ILE
1	C	40	THR
1	C	66	LEU
1	C	72	MET
1	C	76	GLU
1	C	91	THR
1	C	96	ILE
1	C	100	LEU
1	C	103	ASP
1	C	110	LEU
1	C	111	GLN
1	C	139	ILE
1	C	143	TRP
1	C	148	HIS
1	C	153	GLN
1	C	162	ILE
1	C	173	LYS
1	C	174	ASP
1	C	175	THR
1	C	176	LEU
1	C	187	ASN
1	C	213	THR
1	C	216	LYS
1	C	219	GLU
1	C	222	VAL
1	C	246	LEU
1	C	252	ASN
1	C	253	LEU
1	C	261	SER
2	D	4	THR
2	D	16	GLU
2	D	21	ASN
2	D	35	ILE
2	D	53	ASP
2	D	61	SER
2	D	64	LEU

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Mol	Chain	Res	Type
2	D	94	LYS
2	D	98	ASP

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

Mol	Chain	Res	Type
1	A	71	GLN
1	A	81	GLN
1	A	117	GLN
1	A	123	ASN
1	A	137	HIS
1	A	217	ASN
1	C	117	GLN
1	C	123	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 1 is monoatomic - leaving 3 for Mogul analysis.

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
3	6FP	C	301	1	14,14,15	2.01	4 (28%)	16,20,21	1.62	3 (18%)
3	6FP	A	301	1	14,14,15	2.10	4 (28%)	16,20,21	1.84	4 (25%)
5	PO4	A	303	-	4,4,4	1.70	1 (25%)	6,6,6	0.58	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
3	6FP	C	301	1	-	-	0/2/2/2
3	6FP	A	301	1	-	-	0/2/2/2

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	301	6FP	C9-C6	-5.09	1.38	1.50
3	C	301	6FP	C9-C6	-4.85	1.39	1.50
3	C	301	6FP	C4A-C4	-3.87	1.39	1.47
3	A	301	6FP	C7-N8	3.86	1.42	1.34
3	A	301	6FP	C4A-C4	-3.28	1.41	1.47
5	A	303	PO4	P-O1	2.93	1.57	1.50
3	C	301	6FP	C7-N8	2.59	1.39	1.34
3	C	301	6FP	C4-N3	2.35	1.41	1.37
3	A	301	6FP	C4-N3	2.26	1.41	1.37

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	301	6FP	C4A-N5-C6	3.64	121.66	118.28
3	A	301	6FP	C6-C7-N8	-3.42	118.67	123.20
3	C	301	6FP	C2-N3-C4	-3.08	119.47	125.11
3	C	301	6FP	C6-C7-N8	-2.83	119.45	123.20
3	A	301	6FP	C2-N3-C4	-2.63	120.29	125.11
3	C	301	6FP	C4A-N5-C6	2.40	120.51	118.28
3	A	301	6FP	C9-C6-N5	2.07	119.69	116.56

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	301	6FP	1	0
3	A	301	6FP	1	0

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	265/271 (97%)	0.09	4 (1%) 72 52	11, 34, 63, 80	0
1	C	258/271 (95%)	0.15	7 (2%) 56 36	10, 33, 62, 85	0
2	B	97/99 (97%)	0.11	1 (1%) 79 63	20, 44, 71, 100	0
2	D	98/99 (98%)	0.18	3 (3%) 51 32	22, 43, 72, 74	0
All	All	718/740 (97%)	0.13	15 (2%) 63 43	10, 36, 66, 100	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	178	ARG	4.2
2	D	53	ASP	3.9
1	C	189	LYS	3.4
2	D	55	SER	3.0
1	A	248	PRO	2.5
1	C	16	ILE	2.5
1	C	141	GLN	2.4
1	C	219	GLU	2.3
1	C	248	PRO	2.3
1	A	16	ILE	2.2
1	C	147	GLN	2.2
2	B	97	ARG	2.1
1	A	149	GLU	2.1
1	C	17	HIS	2.0
2	D	1	ILE	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($\text{\AA}^2$)	Q<0.9
5	PO4	A	303	5/5	0.79	0.18	90,91,92,92	0
3	6FP	C	301	13/14	0.93	0.11	48,52,55,56	0
4	CL	A	302	1/1	0.94	0.07	51,51,51,51	0
3	6FP	A	301	13/14	0.94	0.10	35,37,40,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.