



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 11:24 PM UTC

PDB ID : 6Z9X / pdb_00006z9x
Title : Human Class I Major Histocompatibility Complex, A02 allele, presenting LLS (t-butyl)Y FGTPT
Authors : Rizkallah, P.J.; Man, S.; Redman, J.E.
Deposited on : 2020-06-04
Resolution : 2.68 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

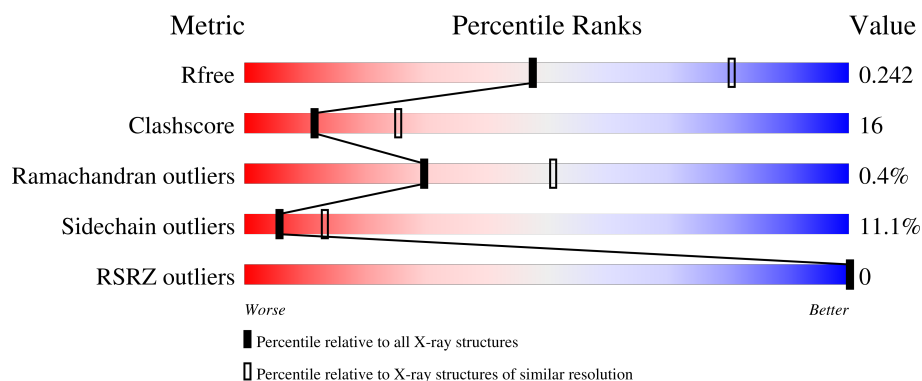
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

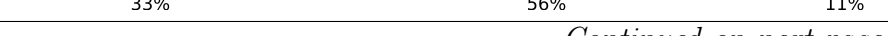
The reported resolution of this entry is 2.68 Å.

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	5070 (2.70-2.66)
Clashscore	190562	5409 (2.70-2.66)
Ramachandran outliers	187476	5324 (2.70-2.66)
Sidechain outliers	187428	5324 (2.70-2.66)
RSRZ outliers	180081	5070 (2.70-2.66)

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	276	
1	D	276	
2	B	100	
2	E	100	
3	C	9	

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Mol	Chain	Length	Quality of chain
3	F	9	33% 44% 22%

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	SO4	B	101	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6412 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 MHC class I antigen.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	276	Total	C	N	O	S	0	0	0
			2254	1408	410	427	9			
1	D	276	Total	C	N	O	S	0	0	0
			2254	1408	410	427	9			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			
2	E	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	MET	-	initiating methionine	UNP P61769
E	0	MET	-	initiating methionine	UNP P61769

- Molecule 3 is a protein called LEU-LEU-SER-TUR-PHE-GLY-THR-PRO-THR.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	9	Total	C	N	O	0	0	0
			75	52	9	14			
3	F	9	Total	C	N	O	0	0	0
			75	52	9	14			

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		

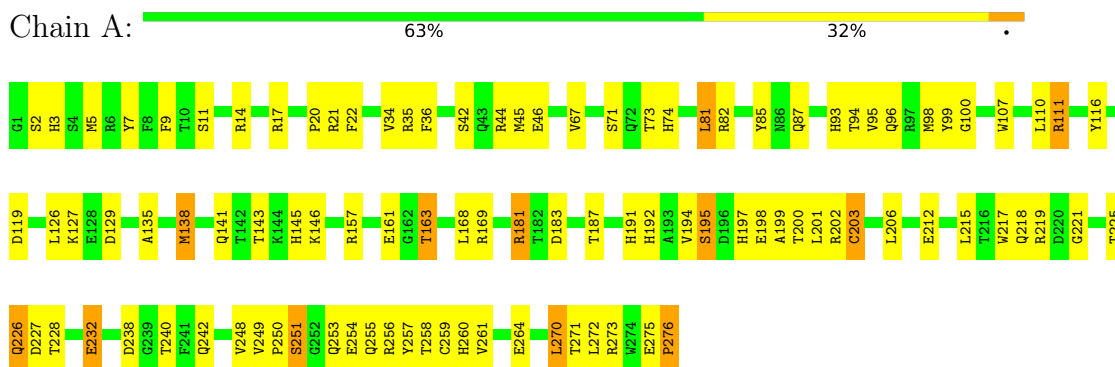
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	18	Total	O	0	0
			18	18		
5	B	8	Total	O	0	0
			8	8		
5	D	19	Total	O	0	0
			19	19		
5	E	15	Total	O	0	0
			15	15		

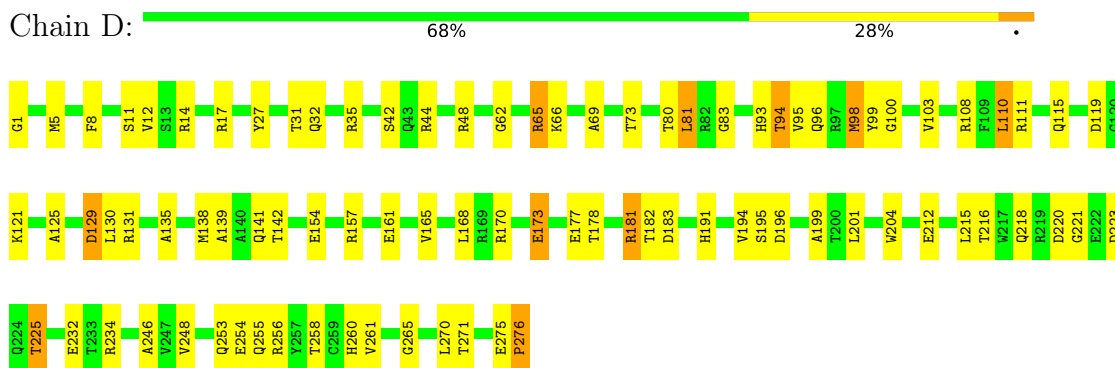
3 Residue-property plots [i](#)

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.

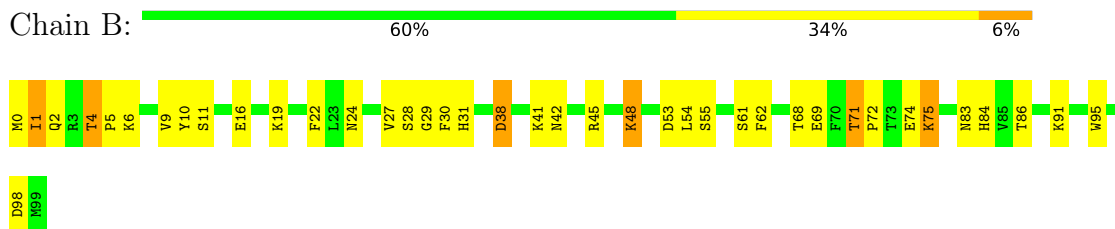
- Molecule 1: MHC class I antigen



- Molecule 1: MHC class I antigen



- Molecule 2: Beta-2-microglobulin

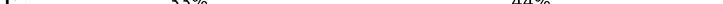


- Molecule 2: Beta-2-microglobulin

MO	I1	Q2	R3	T4	P5	K6	I7	Q8	Q9	Y10	S11	S28	G29	F30	H31	V37	D38	K41	N42	R45	I46	H51	S52	D53	L54	S55	D59	V60	S61	L65	T68	T71	P72	V82	H83	N84	K91	I92	V93	K94	V95	D96	V99
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- Chain C: 33% 56% 11%

L1		QC4	F5	G6	T7	P8	T9
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- Chain F:  33% 44% 22%

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	206.24Å 49.35Å 115.90Å 90.00° 123.85° 90.00°	Depositor
Resolution (Å)	51.34 – 2.68 51.34 – 2.70	Depositor EDS
% Data completeness (in resolution range)	80.2 (51.34-2.68) 80.2 (51.34-2.70)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.54 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.8.0257	Depositor
R, R_{free}	0.186 , 0.241 0.189 , 0.242	Depositor DCC
R_{free} test set	1034 reflections (3.81%)	wwPDB-VP
Wilson B-factor (Å ²)	27.3	Xtriage
Anisotropy	0.122	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 0.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.21$	Xtriage
Estimated twinning fraction	0.388 for -h-2*1,-k,l	Xtriage
Reported twinning fraction	0.600 for H, K, L 0.400 for H+4/2L, -K, -L	Depositor
Outliers	0 of 22144 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6412	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.21% 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: SO4, QCN

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	1.19	1/2320 (0.0%)	1.44	2/3149 (0.1%)
1	D	1.16	3/2320 (0.1%)	1.43	1/3149 (0.0%)
2	B	1.18	0/860	1.32	0/1162
2	E	1.20	1/860 (0.1%)	1.36	0/1162
3	C	1.13	0/59	1.57	0/77
3	F	1.28	0/59	1.65	1/77 (1.3%)
All	All	1.18	5/6478 (0.1%)	1.41	4/8776 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	D	0	3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	194	VAL	C-O	-6.35	1.17	1.24
1	D	99	TYR	C-O	5.88	1.31	1.23
1	A	20	PRO	C-O	-5.53	1.17	1.23
1	D	204	TRP	C-O	5.24	1.30	1.24
2	E	68	THR	C-O	5.17	1.29	1.23

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	250	PRO	CB-CA-C	-7.45	103.44	111.56
1	D	129	ASP	CA-CB-CG	6.90	119.50	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	200	THR	CA-CB-OG1	-6.51	99.84	109.60
3	F	6	GLY	CA-C-O	-5.58	116.45	122.47

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	178	THR	Peptide
1	D	195	SER	Peptide
1	D	225	THR	Peptide

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2254	0	2103	81	1
1	D	2254	0	2103	60	3
2	B	837	0	803	39	1
2	E	837	0	803	24	2
3	C	75	0	62	9	1
3	F	75	0	62	10	0
4	A	5	0	0	0	0
4	B	5	0	0	2	0
4	D	5	0	0	0	0
4	E	5	0	0	1	0
5	A	18	0	0	4	0
5	B	8	0	0	0	0
5	D	19	0	0	2	0
5	E	15	0	0	1	0
All	All	6412	0	5936	196	4

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 16.

All (196) 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:D:182:THR:HG21	1:D:265:GLY:CA	1.88	1.03
1:A:35:ARG:HD2	2:B:53:ASP:OD2	1.60	1.01
1:A:95:VAL:HG13	1:A:116:TYR:CE1	2.01	0.95
1:A:35:ARG:HD2	2:B:53:ASP:CG	1.94	0.93
1:D:218:GLN:HE21	1:D:221:GLY:HA2	1.34	0.92
2:B:2:GLN:HE21	2:B:86:THR:HG22	1.33	0.91
2:B:9:VAL:N	4:B:101:SO4:O1	2.05	0.88
1:A:218:GLN:HE21	1:A:221:GLY:HA2	1.39	0.86
1:A:192:HIS:NE2	2:B:98:ASP:OD2	2.10	0.84
1:A:73:THR:HG21	3:C:6:GLY:HA3	1.61	0.83
1:A:135:ALA:HB3	1:A:141:GLN:NE2	1.94	0.82
1:D:182:THR:HG21	1:D:265:GLY:HA2	1.61	0.81
1:D:135:ALA:HB3	1:D:141:GLN:HE21	1.45	0.80
1:D:170:ARG:HA	1:D:173:GLU:OE2	1.84	0.78
1:D:182:THR:CG2	1:D:265:GLY:CA	2.62	0.78
1:A:34:VAL:HB	1:A:45:MET:HE2	1.67	0.77
1:A:135:ALA:HB3	1:A:141:GLN:HE21	1.48	0.77
1:D:103:VAL:HG11	1:D:165:VAL:HG13	1.65	0.77
1:A:95:VAL:HG13	1:A:116:TYR:HE1	1.49	0.76
1:A:119:ASP:HB3	2:B:0:MET:HB2	1.67	0.76
1:A:73:THR:HG21	3:C:6:GLY:CA	2.15	0.76
1:A:256:ARG:NH2	5:A:401:HOH:O	2.19	0.75
2:B:0:MET:SD	2:B:1:ILE:O	2.45	0.75
1:A:157:ARG:HG3	1:A:161:GLU:OE2	1.87	0.75
1:D:182:THR:CG2	1:D:265:GLY:HA3	2.17	0.74
1:A:218:GLN:HG3	1:A:260:HIS:CE1	2.23	0.74
2:B:1:ILE:CG2	2:B:2:GLN:N	2.51	0.72
1:D:218:GLN:NE2	1:D:221:GLY:HA2	2.04	0.72
2:E:4:THR:HG22	2:E:5:PRO:HD2	1.71	0.72
3:F:4:QCN:N	3:F:4:QCN:CD1	2.52	0.71
1:A:258:THR:OG1	1:A:260:HIS:HE1	1.72	0.71
2:B:2:GLN:NE2	2:B:86:THR:HG22	2.05	0.71
1:D:258:THR:OG1	1:D:260:HIS:HE1	1.75	0.70
1:A:94:THR:HB	5:A:402:HOH:O	1.92	0.70
1:D:215:LEU:HD23	1:D:261:VAL:HG22	1.72	0.69
1:D:215:LEU:CD2	1:D:261:VAL:HG22	2.23	0.69
1:A:35:ARG:HD2	2:B:53:ASP:OD1	1.91	0.69
2:E:1:ILE:O	2:E:2:GLN:HG2	1.92	0.69
2:E:37:VAL:HG22	2:E:82:VAL:HG22	1.74	0.68
1:A:218:GLN:NE2	1:A:221:GLY:HA2	2.08	0.67
1:A:94:THR:CB	5:A:402:HOH:O	2.42	0.67
3:C:6:GLY:O	3:C:8:PRO:HD3	1.96	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:192:HIS:CE1	2:B:98:ASP:OD2	2.49	0.65
1:A:226:GLN:O	1:A:228:THR:HG23	1.97	0.65
3:F:4:QCN:C4	3:F:4:QCN:OH	2.45	0.64
1:A:42:SER:OG	1:A:46:GLU:OE2	2.13	0.63
2:B:1:ILE:HG23	2:B:2:GLN:N	2.13	0.63
1:A:11:SER:HB2	1:A:74:HIS:HD2	1.65	0.62
2:E:94:LYS:NZ	2:E:94:LYS:HB3	2.15	0.62
1:D:81:LEU:CD1	3:F:9:THR:HG21	2.30	0.62
1:D:81:LEU:HD13	3:F:9:THR:HG21	1.80	0.61
2:E:83:ASN:HD22	2:E:84:HIS:H	1.47	0.61
1:D:32:GLN:NE2	1:D:48:ARG:HG3	2.16	0.61
2:B:4:THR:HG22	2:B:5:PRO:HD2	1.81	0.60
1:A:232:GLU:OE1	2:B:28:SER:OG	2.11	0.60
2:B:1:ILE:HG23	2:B:2:GLN:H	1.66	0.60
2:B:2:GLN:HE21	2:B:86:THR:CG2	2.11	0.60
1:D:35:ARG:HD2	2:E:53:ASP:OD1	2.03	0.59
1:A:218:GLN:NE2	5:A:403:HOH:O	2.36	0.59
3:F:3:SER:HB2	3:F:5:PHE:CD1	2.38	0.58
1:D:157:ARG:HG3	1:D:161:GLU:OE2	2.04	0.57
2:E:11:SER:O	4:E:301:SO4:O4	2.22	0.57
1:D:32:GLN:HE21	1:D:48:ARG:HG3	1.70	0.57
1:D:96:GLN:OE1	2:E:31:HIS:HE1	1.88	0.56
1:A:191:HIS:NE2	1:A:199:ALA:CB	2.68	0.56
1:D:135:ALA:CB	1:D:141:GLN:HE21	2.16	0.56
2:B:16:GLU:OE2	2:B:19:LYS:HD2	2.06	0.55
1:D:35:ARG:HD2	2:E:53:ASP:CG	2.32	0.55
1:A:258:THR:HG22	1:A:273:ARG:HG2	1.88	0.55
3:F:3:SER:HB2	3:F:5:PHE:HD1	1.72	0.54
1:D:275:GLU:O	1:D:276:PRO:O	2.26	0.54
1:D:218:GLN:HE21	1:D:221:GLY:CA	2.13	0.54
1:A:11:SER:HB2	1:A:74:HIS:CD2	2.43	0.53
2:B:5:PRO:HB3	2:B:30:PHE:HB3	1.90	0.53
2:B:48:LYS:O	2:B:68:THR:OG1	2.18	0.53
1:A:119:ASP:HB3	2:B:0:MET:CB	2.39	0.53
1:D:157:ARG:CG	1:D:161:GLU:OE2	2.57	0.53
1:A:36:PHE:CD2	1:A:67:VAL:HG11	2.44	0.52
2:B:54:LEU:HD21	2:B:62:PHE:CE1	2.45	0.52
2:E:38:ASP:N	2:E:38:ASP:OD1	2.42	0.52
2:B:38:ASP:OD2	2:B:45:ARG:HD2	2.09	0.52
1:D:170:ARG:HA	1:D:173:GLU:CD	2.33	0.52
1:A:181:ARG:NH1	1:A:183:ASP:OD2	2.43	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:195:SER:HB2	1:A:198:GLU:HB2	1.92	0.52
1:D:182:THR:HG21	1:D:265:GLY:N	2.24	0.52
1:A:226:GLN:O	1:A:228:THR:N	2.43	0.51
1:D:253:GLN:OE1	1:D:256:ARG:NH2	2.44	0.51
1:A:146:LYS:HE2	3:C:9:THR:O	2.11	0.51
1:D:115:GLN:HE21	1:D:125:ALA:HB2	1.76	0.51
2:E:94:LYS:HG3	5:E:407:HOH:O	2.11	0.51
1:A:201:LEU:HD12	1:A:249:VAL:HG11	1.92	0.50
1:A:206:LEU:HD23	1:A:242:GLN:HG2	1.94	0.49
1:A:95:VAL:HG11	1:A:116:TYR:OH	2.12	0.49
1:A:215:LEU:HD23	1:A:261:VAL:HG22	1.95	0.49
1:A:129:ASP:OD1	1:A:129:ASP:C	2.55	0.49
1:A:96:GLN:OE1	2:B:31:HIS:HE1	1.95	0.49
1:A:275:GLU:O	1:A:276:PRO:O	2.30	0.49
2:E:83:ASN:HD22	2:E:84:HIS:N	2.10	0.49
1:A:219:ARG:HD2	1:A:257:TYR:CZ	2.48	0.49
1:A:111:ARG:HG2	1:A:111:ARG:HH21	1.78	0.49
3:C:1:LEU:HD21	3:C:4:QCN:C3	2.42	0.49
1:D:275:GLU:O	1:D:276:PRO:C	2.56	0.49
1:D:201:LEU:O	1:D:246:ALA:HA	2.13	0.49
3:C:7:THR:O	3:C:7:THR:OG1	2.29	0.48
1:A:107:TRP:HE3	1:A:169:ARG:HG2	1.77	0.48
1:D:129:ASP:O	1:D:131:ARG:NH2	2.46	0.48
1:A:215:LEU:CD2	1:A:261:VAL:HG22	2.43	0.48
1:D:12:VAL:HG22	1:D:94:THR:HG23	1.94	0.48
1:A:9:PHE:CE2	1:A:99:TYR:CE2	3.02	0.48
1:A:203:CYS:HB3	1:A:217:TRP:CZ2	2.49	0.48
2:E:9:VAL:HG13	2:E:95:TRP:HA	1.96	0.48
2:B:27:VAL:HG23	2:B:30:PHE:HE2	1.79	0.48
2:B:1:ILE:HG22	2:B:2:GLN:N	2.29	0.47
2:B:9:VAL:HG12	2:B:95:TRP:HB2	1.96	0.47
2:B:10:TYR:CE1	2:B:24:ASN:HB2	2.49	0.47
1:D:121:LYS:NZ	5:D:405:HOH:O	2.47	0.47
1:D:182:THR:HG22	1:D:265:GLY:HA3	1.92	0.47
1:A:35:ARG:CD	2:B:53:ASP:OD2	2.49	0.47
1:D:181:ARG:O	1:D:181:ARG:HG2	2.13	0.47
1:D:129:ASP:O	1:D:130:LEU:HB2	2.13	0.47
1:D:42:SER:O	1:D:44:ARG:HG2	2.15	0.47
1:D:258:THR:OG1	1:D:260:HIS:CE1	2.62	0.47
2:E:41:LYS:O	2:E:42:ASN:C	2.58	0.46
1:A:35:ARG:C	1:A:45:MET:HE3	2.41	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:11:SER:OG	5:D:401:HOH:O	2.20	0.46
1:D:27:TYR:HA	1:D:31:THR:O	2.14	0.46
1:A:85:TYR:HB2	1:A:87:GLN:OE1	2.16	0.46
1:A:206:LEU:CD2	1:A:242:GLN:HG2	2.46	0.46
1:A:11:SER:HA	1:A:21:ARG:O	2.16	0.46
2:E:37:VAL:HG22	2:E:82:VAL:HG13	1.97	0.46
1:D:170:ARG:CA	1:D:173:GLU:OE2	2.62	0.46
1:A:273:ARG:O	1:A:275:GLU:HG3	2.15	0.46
1:A:36:PHE:CD2	1:A:67:VAL:CG1	2.99	0.45
1:D:65:ARG:HA	1:D:65:ARG:HD3	1.73	0.45
1:A:275:GLU:O	1:A:276:PRO:C	2.58	0.45
1:A:238:ASP:OD1	1:A:240:THR:OG1	2.20	0.45
2:B:83:ASN:HD22	2:B:84:HIS:H	1.63	0.45
1:D:139:ALA:O	1:D:142:THR:HB	2.17	0.45
1:A:218:GLN:HE21	1:A:221:GLY:CA	2.21	0.45
2:B:48:LYS:HA	2:B:48:LYS:HE3	1.97	0.45
2:E:7:ILE:HB	2:E:93:VAL:HG21	1.99	0.45
1:A:5:MET:O	1:A:100:GLY:HA3	2.17	0.44
1:D:234:ARG:NH1	2:E:10:TYR:CD2	2.86	0.44
1:A:81:LEU:HD12	1:A:81:LEU:HA	1.80	0.44
1:D:1:GLY:HA3	1:D:110:LEU:CD1	2.47	0.44
1:D:81:LEU:HD12	1:D:81:LEU:HA	1.79	0.44
1:A:93:HIS:HD2	1:A:119:ASP:OD2	2.01	0.44
1:A:95:VAL:HG13	1:A:116:TYR:CZ	2.51	0.44
1:A:191:HIS:NE2	1:A:199:ALA:HB1	2.32	0.44
2:B:41:LYS:O	2:B:42:ASN:C	2.60	0.44
2:E:71:THR:HA	2:E:72:PRO:HD2	1.89	0.44
1:A:22:PHE:CD2	1:A:71:SER:HB3	2.53	0.43
1:A:146:LYS:HE3	3:C:8:PRO:O	2.18	0.43
1:A:187:THR:HB	1:A:272:LEU:CD1	2.47	0.43
1:A:98:MET:HE3	1:A:98:MET:HB3	1.81	0.43
1:A:42:SER:O	1:A:44:ARG:HG2	2.18	0.43
1:D:66:LYS:NZ	3:F:2:LEU:O	2.47	0.43
1:D:93:HIS:HD2	1:D:119:ASP:OD2	2.01	0.43
1:A:127:LYS:HB3	1:A:127:LYS:HE2	1.86	0.43
1:A:253:GLN:HB3	1:A:256:ARG:HD3	2.00	0.43
2:E:51:HIS:HA	2:E:65:LEU:O	2.19	0.43
2:B:54:LEU:HD21	2:B:62:PHE:CD1	2.54	0.42
1:D:5:MET:O	1:D:100:GLY:HA3	2.19	0.42
1:D:81:LEU:HD13	3:F:9:THR:CG2	2.47	0.42
1:A:270:LEU:HD23	1:A:270:LEU:HA	1.88	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:9:VAL:O	4:B:101:SO4:O1	2.37	0.42
1:D:8:PHE:CE2	1:D:98:MET:HG3	2.54	0.42
1:D:5:MET:HB2	1:D:168:LEU:HD13	2.02	0.42
1:D:216:THR:HG21	1:D:223:ASP:OD1	2.19	0.42
1:D:220:ASP:OD2	1:D:256:ARG:NH1	2.37	0.42
2:E:29:GLY:HA2	2:E:61:SER:OG	2.19	0.42
1:A:2:SER:C	1:A:3:HIS:CG	2.97	0.42
1:D:35:ARG:HD2	2:E:53:ASP:OD2	2.21	0.41
1:A:258:THR:OG1	1:A:260:HIS:CE1	2.62	0.41
2:B:29:GLY:HA2	2:B:61:SER:OG	2.19	0.41
2:B:71:THR:HA	2:B:72:PRO:HD2	1.87	0.41
1:D:81:LEU:HD11	3:F:9:THR:HG21	2.02	0.41
3:F:5:PHE:CD1	3:F:5:PHE:N	2.89	0.41
1:A:163:THR:HG21	3:C:1:LEU:HD23	2.03	0.41
1:D:191:HIS:NE2	1:D:199:ALA:CB	2.84	0.41
1:A:7:TYR:O	1:A:98:MET:HA	2.20	0.41
1:A:197:HIS:O	1:A:198:GLU:HG3	2.20	0.41
1:A:202:ARG:HD2	2:B:98:ASP:OD1	2.20	0.41
1:A:143:THR:HG23	3:C:9:THR:HA	2.02	0.41
2:B:11:SER:HA	2:B:22:PHE:O	2.20	0.41
1:D:69:ALA:O	1:D:73:THR:HG23	2.20	0.41
1:A:5:MET:HB2	1:A:168:LEU:HD13	2.02	0.41
1:A:258:THR:CG2	1:A:273:ARG:NH1	2.84	0.41
2:B:1:ILE:O	2:B:2:GLN:HG3	2.21	0.41
1:D:181:ARG:NH1	1:D:183:ASP:OD2	2.52	0.40
2:E:94:LYS:HB3	2:E:94:LYS:HZ3	1.85	0.40
2:B:9:VAL:HA	2:B:24:ASN:O	2.22	0.40
2:E:59:ASP:OD1	2:E:59:ASP:C	2.64	0.40
1:D:182:THR:CG2	1:D:265:GLY:HA2	2.40	0.40
1:A:138:MET:HE3	1:A:138:MET:HB3	1.86	0.40
2:E:46:ILE:HG21	2:E:68:THR:HG21	2.04	0.40

All (4) 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)
3:C:4:QCN:OH	1:D:154:GLU:OE2[2_555]	2.13	0.07
1:D:83:GLY:O	2:E:45:ARG:NH2[4_556]	2.16	0.04
1:A:251:SER:O	1:D:253:GLN:NE2[2_546]	2.17	0.03
2:B:75:LYS:NZ	2:E:96:ASP:OD2[2_556]	2.19	0.01

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	274/276 (99%)	251 (92%)	22 (8%)	1 (0%)	30	51
1	D	274/276 (99%)	255 (93%)	17 (6%)	2 (1%)	18	37
2	B	98/100 (98%)	90 (92%)	8 (8%)	0	100	100
2	E	98/100 (98%)	91 (93%)	7 (7%)	0	100	100
3	C	6/9 (67%)	4 (67%)	2 (33%)	0	100	100
3	F	6/9 (67%)	6 (100%)	0	0	100	100
All	All	756/770 (98%)	697 (92%)	56 (7%)	3 (0%)	30	51

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	177	GLU
1	A	227	ASP
1	D	62	GLY

5.3.2 Protein sidechains

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	232/232 (100%)	205 (88%)	27 (12%)	5	12
1	D	232/232 (100%)	208 (90%)	24 (10%)	7	15
2	B	95/95 (100%)	84 (88%)	11 (12%)	5	12
2	E	95/95 (100%)	86 (90%)	9 (10%)	8	18

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	C	7/7 (100%)	6 (86%)	1 (14%)	3	7
3	F	7/7 (100%)	5 (71%)	2 (29%)	0	1
All	All	668/668 (100%)	594 (89%)	74 (11%)	6	13

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

Mol	Chain	Res	Type
1	A	14	ARG
1	A	17	ARG
1	A	81	LEU
1	A	82	ARG
1	A	110	LEU
1	A	111	ARG
1	A	126	LEU
1	A	138	MET
1	A	145	HIS
1	A	163	THR
1	A	181	ARG
1	A	194	VAL
1	A	195	SER
1	A	203	CYS
1	A	212	GLU
1	A	225	THR
1	A	226	GLN
1	A	232	GLU
1	A	248	VAL
1	A	251	SER
1	A	254	GLU
1	A	255	GLN
1	A	259	CYS
1	A	264	GLU
1	A	270	LEU
1	A	271	THR
1	A	276	PRO
2	B	1	ILE
2	B	4	THR
2	B	6	LYS
2	B	38	ASP
2	B	48	LYS
2	B	55	SER
2	B	69	GLU

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Mol	Chain	Res	Type
2	B	71	THR
2	B	74	GLU
2	B	75	LYS
2	B	91	LYS
3	C	7	THR
1	D	14	ARG
1	D	17	ARG
1	D	65	ARG
1	D	80	THR
1	D	81	LEU
1	D	94	THR
1	D	95	VAL
1	D	98	MET
1	D	108	ARG
1	D	110	LEU
1	D	111	ARG
1	D	138	MET
1	D	173	GLU
1	D	181	ARG
1	D	196	ASP
1	D	212	GLU
1	D	225	THR
1	D	232	GLU
1	D	248	VAL
1	D	254	GLU
1	D	255	GLN
1	D	270	LEU
1	D	271	THR
1	D	276	PRO
2	E	0	MET
2	E	1	ILE
2	E	4	THR
2	E	28	SER
2	E	38	ASP
2	E	55	SER
2	E	71	THR
2	E	91	LYS
2	E	94	LYS
3	F	5	PHE
3	F	9	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	GLN
1	A	74	HIS
1	A	86	ASN
1	A	87	GLN
1	A	93	HIS
1	A	141	GLN
1	A	151	HIS
1	A	180	GLN
1	A	218	GLN
1	A	255	GLN
1	A	260	HIS
2	B	2	GLN
2	B	13	HIS
2	B	31	HIS
2	B	83	ASN
2	B	89	GLN
1	D	32	GLN
1	D	43	GLN
1	D	74	HIS
1	D	86	ASN
1	D	93	HIS
1	D	115	GLN
1	D	141	GLN
1	D	151	HIS
1	D	197	HIS
1	D	218	GLN
1	D	260	HIS
2	E	13	HIS
2	E	31	HIS
2	E	83	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	SO4	D	301	-	4,4,4	0.25	0	6,6,6	0.14	0
4	SO4	E	301	-	4,4,4	0.35	0	6,6,6	0.14	0
4	SO4	B	101	-	4,4,4	0.25	0	6,6,6	0.17	0
4	SO4	A	301	-	4,4,4	0.24	0	6,6,6	0.16	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	E	301	SO4	1	0
4	B	101	SO4	2	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	276/276 (100%)	-1.64	0 100 100	13, 22, 33, 40	0
1	D	276/276 (100%)	-1.60	0 100 100	17, 27, 37, 45	0
2	B	100/100 (100%)	-1.68	0 100 100	14, 18, 25, 31	0
2	E	100/100 (100%)	-1.67	0 100 100	13, 20, 30, 32	0
3	C	8/9 (88%)	-1.52	0 100 100	20, 24, 26, 28	0
3	F	8/9 (88%)	-1.42	0 100 100	24, 26, 27, 28	0
All	All	768/770 (99%)	-1.63	0 100 100	13, 23, 34, 45	0

There are no RSRZ outliers to report.

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	SO4	B	101	5/5	0.99	0.07	27,28,28,29	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	SO4	D	301	5/5	0.99	0.03	39,43,46,48	0
4	SO4	A	301	5/5	1.00	0.03	26,26,29,31	0
4	SO4	E	301	5/5	1.00	0.05	25,26,28,28	0

6.5 Other polymers [i](#)

There are no such residues in this entry.