



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:12 PM UTC

PDB ID : 4IIQ / pdb_00004iiq
Title : Crystal structure of a human MAIT TCR in complex with bovine MR1
Authors : Lopez-Sagaseta, J.; Adams, E.J.
Deposited on : 2012-12-20
Resolution : 2.86 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

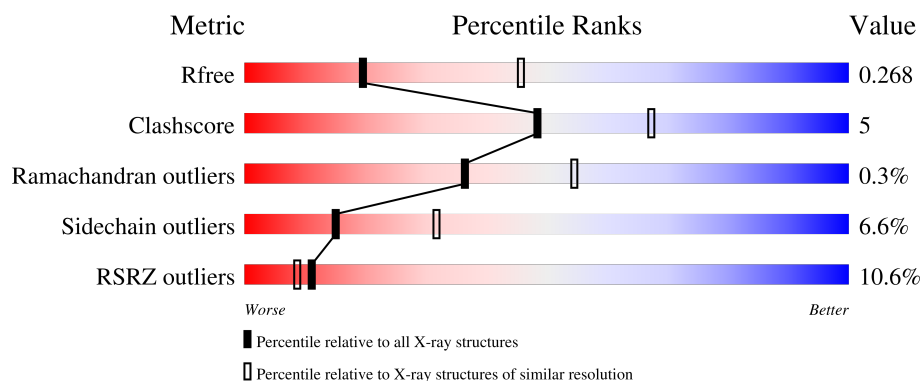
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.86 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	1407 (2.88-2.84)
Clashscore	190562	1446 (2.88-2.84)
Ramachandran outliers	187476	1406 (2.88-2.84)
Sidechain outliers	187428	1407 (2.88-2.84)
RSRZ outliers	180081	1408 (2.88-2.84)

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	205	17% 78% 18% ..
2	B	253	6% 78% 16% ..
3	C	392	9% 74% 16% • 8%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6307 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 Human Mucosal Associated Invariant T cell receptor alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	197	Total	C	N	O	S	0	0	0
			1502	948	242	304	8			

- Molecule 2 is a protein called Human Mucosal Associated Invariant T cell receptor beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	243	Total	C	N	O	S	0	0	0
			1883	1185	323	366	9			

- Molecule 3 is a protein called Beta-2-microglobulin, MHC class I-related protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	360	Total	C	N	O	S	0	0	0
			2900	1857	503	527	13			

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	99	GLY	-	linker	UNP P01888
C	100	GLY	-	linker	UNP P01888
C	101	GLY	-	linker	UNP P01888
C	102	GLY	-	linker	UNP P01888
C	103	SER	-	linker	UNP P01888
C	104	GLY	-	linker	UNP P01888
C	105	GLY	-	linker	UNP P01888
C	106	SER	-	linker	UNP P01888
C	107	GLY	-	linker	UNP P01888
C	108	SER	-	linker	UNP P01888
C	109	GLY	-	linker	UNP P01888
C	110	GLY	-	linker	UNP P01888

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Chain	Residue	Modelled	Actual	Comment	Reference
C	111	GLY	-	linker	UNP P01888
C	112	GLY	-	linker	UNP P01888
C	113	SER	-	linker	UNP P01888

- 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	C	1	Total	O	S	0	0
			5	4	1		

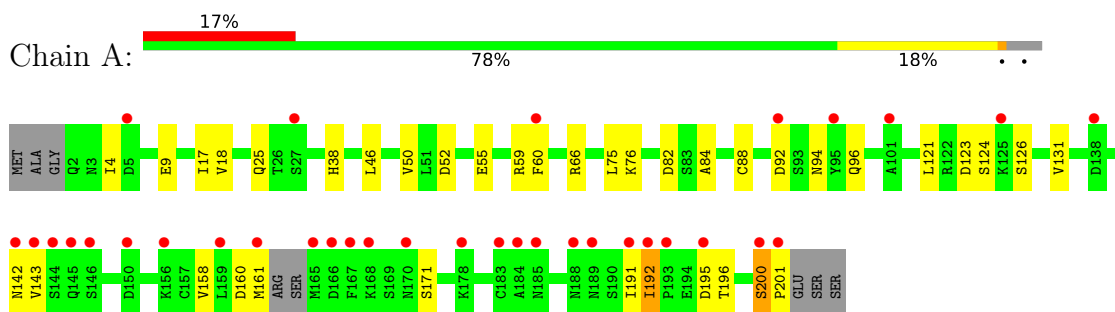
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	3	Total	O	0	0
			3	3		
5	B	1	Total	O	0	0
			1	1		
5	C	3	Total	O	0	0
			3	3		

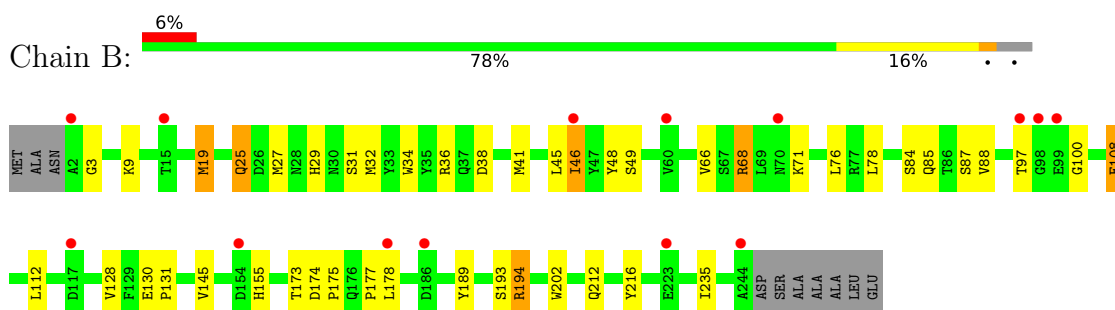
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

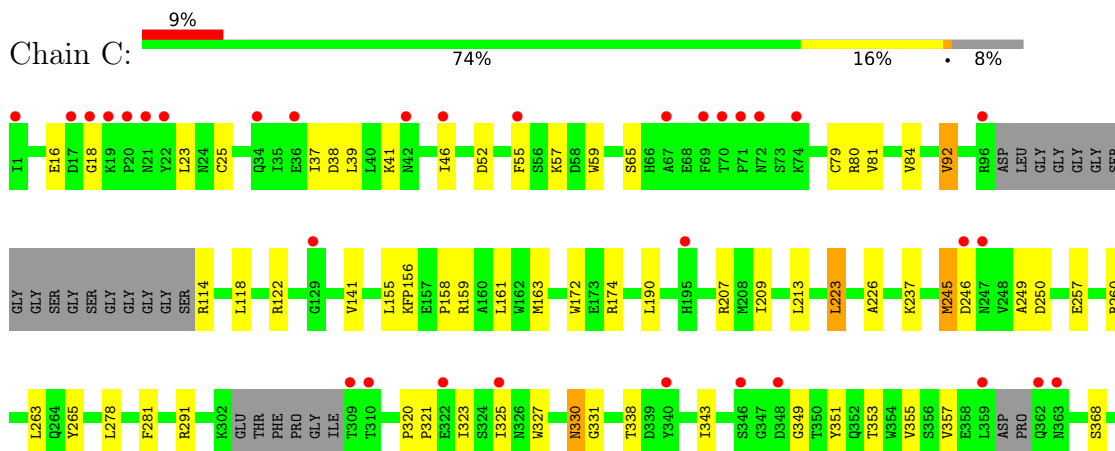
- Molecule 1: Human Mucosal Associated Invariant T cell receptor alpha chain

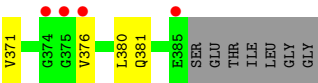


- Molecule 2: Human Mucosal Associated Invariant T cell receptor beta chain



- Molecule 3: Beta-2-microglobulin, MHC class I-related protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	82.05Å 87.41Å 156.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.77 – 2.86 44.77 – 2.86	Depositor EDS
% Data completeness (in resolution range)	99.5 (44.77-2.86) 99.5 (44.77-2.86)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.57 (at 2.86Å)	Xtriage
Refinement program	PHENIX 1.8.1_1168	Depositor
R, R_{free}	0.213 , 0.271 0.216 , 0.268	Depositor DCC
R_{free} test set	1336 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	40.0	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 58.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	6307	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.00% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: KFP, SO4

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	0.30	0/1535	0.71	0/2088
2	B	0.31	0/1934	0.72	3/2639 (0.1%)
3	C	0.30	0/2964	0.70	2/4031 (0.0%)
All	All	0.30	0/6433	0.71	5/8758 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	330	ASN	N-CA-C	6.89	120.92	111.54
3	C	349	GLY	N-CA-C	-5.49	107.09	115.00
2	B	174	ASP	CA-C-N	5.05	124.22	118.97
2	B	174	ASP	C-N-CA	5.05	124.22	118.97
2	B	71	LYS	N-CA-C	-5.02	107.10	113.18

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	1502	0	1379	19	0
2	B	1883	0	1743	23	0
3	C	2900	0	2662	28	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	5	0	0	0	0
4	B	5	0	0	0	0
4	C	5	0	0	0	0
5	A	3	0	0	0	0
5	B	1	0	0	0	0
5	C	3	0	0	0	0
All	All	6307	0	5784	66	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 5.

All (66) 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)
2:B:49:SER:OG	2:B:68:ARG:HD2	1.78	0.84
3:C:41:LYS:HB2	3:C:46:ILE:HD11	1.76	0.68
2:B:9:LYS:HE2	2:B:108:GLU:HG3	1.77	0.67
2:B:32:MET:HE2	2:B:68:ARG:NH1	2.12	0.65
3:C:52:ASP:OD2	3:C:159:ARG:NH1	2.30	0.65
3:C:338:THR:HG22	3:C:357:VAL:HG22	1.79	0.63
3:C:343:ILE:HA	3:C:353:THR:HG22	1.81	0.62
1:A:55:GLU:OE2	3:C:260:ARG:NH2	2.33	0.62
1:A:60:PHE:HD1	1:A:75:LEU:HD22	1.66	0.59
3:C:330:ASN:N	3:C:331:GLY:HA2	2.18	0.57
3:C:25:CYS:HB2	3:C:39:LEU:HD21	1.87	0.56
3:C:118:LEU:HB2	3:C:278:LEU:HD13	1.86	0.56
1:A:59:ARG:NH2	1:A:82:ASP:OD2	2.41	0.53
3:C:209:ILE:HG13	3:C:223:LEU:HG	1.90	0.53
2:B:25:GLN:OE1	2:B:29:HIS:N	2.40	0.53
2:B:34:TRP:O	2:B:46:ILE:HG12	2.09	0.52
1:A:142:ASN:HB2	1:A:191:ILE:HB	1.92	0.51
2:B:38:ASP:HB2	2:B:41:MET:HE3	1.92	0.51
3:C:79:CYS:HB3	3:C:92:VAL:HG23	1.93	0.51
3:C:246:ASP:HB3	3:C:249:ALA:H	1.76	0.51
3:C:114:ARG:NH2	3:C:291:ARG:O	2.39	0.50
3:C:321:PRO:HG3	3:C:351:TYR:CE1	2.47	0.50
2:B:173:THR:HG23	2:B:193:SER:HB2	1.93	0.50
1:A:158:VAL:HG23	2:B:175:PRO:HG3	1.96	0.48
2:B:155:HIS:HB3	2:B:216:TYR:HB2	1.95	0.48
2:B:19:MET:HE3	2:B:19:MET:HB3	1.78	0.47
3:C:327:TRP:HD1	3:C:355:VAL:HG12	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:163:MET:HE1	3:C:281:PHE:CG	2.50	0.47
1:A:52:ASP:H	1:A:66:ARG:HH21	1.63	0.47
1:A:124:SER:C	1:A:126:SER:H	2.23	0.47
2:B:100:GLY:HA3	3:C:265:TYR:CD1	2.50	0.47
2:B:3:GLY:HA2	2:B:27:MET:HE3	1.97	0.47
1:A:171:SER:O	2:B:194:ARG:NH2	2.46	0.46
2:B:145:VAL:HG22	2:B:194:ARG:HG3	1.98	0.46
2:B:36:ARG:NH2	2:B:85:GLN:HA	2.31	0.45
3:C:368:SER:HB3	3:C:381:GLN:HA	1.98	0.45
1:A:38:HIS:CD2	1:A:84:ALA:HB2	2.51	0.45
1:A:92:ASP:HB3	1:A:94:ASN:O	2.16	0.45
3:C:245:MET:HE3	3:C:245:MET:HB3	1.87	0.45
1:A:60:PHE:CD1	1:A:75:LEU:HD22	2.51	0.45
3:C:158:PRO:HG3	3:C:172:TRP:CZ2	2.52	0.45
2:B:31:SER:OG	2:B:97:THR:HG23	2.17	0.44
3:C:159:ARG:HA	3:C:159:ARG:HD3	1.86	0.43
1:A:17:ILE:HG13	1:A:76:LYS:HA	2.01	0.43
1:A:160:ASP:OD1	1:A:161:MET:N	2.44	0.43
1:A:18:VAL:HG12	1:A:75:LEU:HB2	2.01	0.43
3:C:122:ARG:HB2	3:C:207:ARG:HB3	2.01	0.43
3:C:59:TRP:CE2	3:C:226:ALA:HB2	2.53	0.43
3:C:325:ILE:HG12	3:C:371:VAL:HG22	2.00	0.43
2:B:131:PRO:HD2	2:B:202:TRP:CZ2	2.54	0.42
1:A:192:ILE:HD13	1:A:192:ILE:HA	1.82	0.42
2:B:45:LEU:HD21	2:B:48:TYR:HB3	2.02	0.42
2:B:87:SER:OG	2:B:88:VAL:N	2.52	0.42
2:B:177:PRO:HB2	2:B:189:TYR:HB3	2.01	0.42
1:A:38:HIS:HD2	1:A:84:ALA:HB2	1.85	0.42
3:C:37:ILE:HG23	3:C:81:VAL:HG22	2.02	0.42
3:C:122:ARG:HD2	3:C:207:ARG:HD3	2.02	0.41
1:A:46:LEU:HD23	1:A:46:LEU:HA	1.90	0.41
1:A:4:ILE:HD11	1:A:88:CYS:SG	2.60	0.41
2:B:36:ARG:NH2	2:B:84:SER:O	2.48	0.41
3:C:38:ASP:OD2	3:C:80:ARG:NH1	2.53	0.41
3:C:257:GLU:HA	3:C:263:LEU:HD11	2.02	0.41
3:C:114:ARG:NH2	3:C:320:PRO:HD3	2.36	0.41
2:B:130:GLU:HA	2:B:131:PRO:HD3	1.97	0.41
2:B:212:GLN:HG3	2:B:235:ILE:HG23	2.03	0.40
1:A:200:SER:OG	1:A:201:PRO:HD3	2.21	0.40

There are no symmetry-related clashes.

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	193/205 (94%)	183 (95%)	10 (5%)	0	100	100
2	B	241/253 (95%)	232 (96%)	9 (4%)	0	100	100
3	C	351/392 (90%)	335 (95%)	14 (4%)	2 (1%)	21	38
All	All	785/850 (92%)	750 (96%)	33 (4%)	2 (0%)	36	54

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	18	GLY
3	C	376	VAL

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	162/181 (90%)	150 (93%)	12 (7%)	13	26
2	B	200/215 (93%)	188 (94%)	12 (6%)	17	36
3	C	294/341 (86%)	275 (94%)	19 (6%)	15	32
All	All	656/737 (89%)	613 (93%)	43 (7%)	15	32

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	GLU

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Mol	Chain	Res	Type
1	A	25	GLN
1	A	50	VAL
1	A	96	GLN
1	A	121	LEU
1	A	123	ASP
1	A	131	VAL
1	A	143	VAL
1	A	192	ILE
1	A	195	ASP
1	A	196	THR
1	A	200	SER
2	B	19	MET
2	B	25	GLN
2	B	46	ILE
2	B	66	VAL
2	B	68	ARG
2	B	76	LEU
2	B	78	LEU
2	B	108	GLU
2	B	112	LEU
2	B	128	VAL
2	B	178	LEU
2	B	194	ARG
3	C	16	GLU
3	C	23	LEU
3	C	55	PHE
3	C	57	LYS
3	C	65	SER
3	C	84	VAL
3	C	92	VAL
3	C	141	VAL
3	C	155	LEU
3	C	161	LEU
3	C	174	ARG
3	C	190	LEU
3	C	213	LEU
3	C	223	LEU
3	C	237	LYS
3	C	245	MET
3	C	250	ASP
3	C	323	ILE
3	C	380	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	ASN
1	A	113	ASN
2	B	30	ASN
2	B	214	GLN
3	C	50	GLN
3	C	203	HIS
3	C	206	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](#)

1 non-standard protein/DNA/RNA residue is 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	$\# Z > 2$
3	KFP	C	156	3	22,23,24	1.79	5 (22%)	20,30,32	1.94	7 (35%)

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	KFP	C	156	3	-	1/10/11/13	0/2/2/2

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	156	KFP	CAF-C6	-5.04	1.40	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	156	KFP	C4A-C4	-3.69	1.40	1.47
3	C	156	KFP	C7-N8	2.92	1.40	1.34
3	C	156	KFP	CAF-NAL	-2.37	1.35	1.45
3	C	156	KFP	C4-N3	2.13	1.41	1.37

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	156	KFP	CAF-NAL-CAJ	4.05	127.45	113.20
3	C	156	KFP	C6-C7-N8	-3.85	119.45	123.14
3	C	156	KFP	C2-N3-C4	-2.86	119.87	125.11
3	C	156	KFP	C4A-C4-N3	2.54	118.92	114.07
3	C	156	KFP	C7-C6-N5	-2.44	119.29	120.87
3	C	156	KFP	CAF-C6-N5	2.22	120.76	116.56
3	C	156	KFP	N3-C2-N1	-2.11	119.46	123.32

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	156	KFP	CAH-CAJ-NAL-CAF

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

3 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	B	301	-	4,4,4	0.24	0	6,6,6	0.10	0
4	SO4	C	401	-	4,4,4	0.23	0	6,6,6	0.19	0
4	SO4	A	301	-	4,4,4	0.23	0	6,6,6	0.09	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.

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	197/205 (96%)	0.76	34 (17%) 4 3	37, 64, 109, 128	0
2	B	243/253 (96%)	0.57	14 (5%) 29 21	39, 67, 104, 122	0
3	C	359/392 (91%)	0.70	37 (10%) 12 10	37, 67, 115, 140	0
All	All	799/850 (94%)	0.67	85 (10%) 11 9	37, 66, 109, 140	0

All (85) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	201	PRO	5.8
2	B	98	GLY	5.5
1	A	166	ASP	5.0
3	C	71	PRO	4.6
1	A	144	SER	4.6
3	C	96	ARG	4.5
3	C	247	ASN	4.3
3	C	385	GLU	4.3
3	C	17	ASP	4.1
2	B	2	ALA	3.9
3	C	21	ASN	3.9
3	C	129	GLY	3.6
3	C	363	ASN	3.6
3	C	309	THR	3.5
1	A	138	ASP	3.5
1	A	189	ASN	3.4
1	A	200	SER	3.3
3	C	36	GLU	3.3
1	A	167	PHE	3.2
1	A	150	ASP	3.2
2	B	117	ASP	3.2
1	A	170	ASN	3.1
3	C	325	ILE	3.1

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Mol	Chain	Res	Type	RSRZ
3	C	359	LEU	2.9
2	B	154	ASP	2.9
3	C	70	THR	2.9
3	C	362	GLN	2.8
2	B	178	LEU	2.8
1	A	165	MET	2.8
1	A	159	LEU	2.8
3	C	34	GLN	2.7
1	A	161	MET	2.7
2	B	223	GLU	2.7
3	C	19	LYS	2.7
3	C	195	HIS	2.7
3	C	69	PHE	2.6
1	A	185	ASN	2.6
1	A	195	ASP	2.6
3	C	374	GLY	2.6
3	C	74	LYS	2.6
1	A	168	LYS	2.5
1	A	143	VAL	2.5
2	B	97	THR	2.5
1	A	142	ASN	2.5
3	C	72	ASN	2.5
3	C	346	SER	2.5
3	C	375	GLY	2.5
1	A	184	ALA	2.5
1	A	178	LYS	2.5
1	A	188	ASN	2.5
1	A	92	ASP	2.4
1	A	60	PHE	2.4
1	A	192	ILE	2.4
1	A	101	ALA	2.4
1	A	145	GLN	2.4
1	A	146	SER	2.4
3	C	20	PRO	2.4
3	C	22	TYR	2.3
3	C	340	TYR	2.3
1	A	95	TYR	2.3
3	C	348	ASP	2.3
3	C	310	THR	2.3
2	B	60	VAL	2.3
2	B	70	ASN	2.2
3	C	18	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	183	CYS	2.2
2	B	244	ALA	2.2
3	C	46	ILE	2.2
3	C	246	ASP	2.2
3	C	376	VAL	2.1
3	C	55	PHE	2.1
1	A	27	SER	2.1
2	B	46	ILE	2.1
2	B	99	GLU	2.1
3	C	322	GLU	2.1
1	A	125	LYS	2.1
1	A	193	PRO	2.1
3	C	42	ASN	2.0
1	A	5	ASP	2.0
1	A	191	ILE	2.0
2	B	186	ASP	2.0
3	C	67	ALA	2.0
1	A	156	LYS	2.0
3	C	1	ILE	2.0
2	B	15	THR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
3	KFP	C	156	22/23	0.84	0.19	57,67,79,83	0

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
4	SO4	B	301	5/5	0.90	0.12	95,97,100,102	0
4	SO4	A	301	5/5	0.91	0.11	86,90,95,96	0
4	SO4	C	401	5/5	0.95	0.11	61,62,63,64	0

6.5 Other polymers [i](#)

There are no such residues in this entry.