



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 01:36 AM UTC

PDB ID : 7OW6 / pdb_00007ow6
Title : Crystal structure of a TCR in complex with HLA-A*11:01 bound to KRAS G12D peptide (VVVGADGVGK)
Authors : Karuppiah, V.; Robinson, R.A.
Deposited on : 2021-06-16
Resolution : 2.64 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

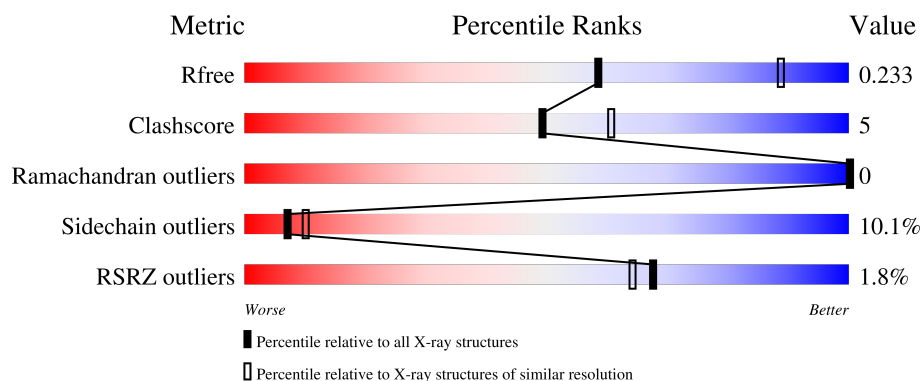
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.64 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	2053 (2.66-2.62)
Clashscore	190562	2097 (2.66-2.62)
Ramachandran outliers	187476	2066 (2.66-2.62)
Sidechain outliers	187428	2066 (2.66-2.62)
RSRZ outliers	180081	2052 (2.66-2.62)

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	 85% 15%
2	B	100	 2% 72% 24% ..
3	C	10	 80% 20%
4	D	203	 5% 60% 25% 11%
5	E	246	 83% 15% ..

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 6536 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
			2251	1399	409	434	9			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	276	PRO	-	expression tag	UNP A0A583ZB34

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	99	Total	C	N	O	S	0	0	0
			828	528	140	157	3			

There is a discrepancy between the modelled and reference sequences:

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

- Molecule 3 is a protein called KRAS G12D peptide (VVVGADGVGK).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	10	Total	C	N	O	0	0	0
			63	39	11	13			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	6	ASP	GLY	engineered mutation	UNP P01111

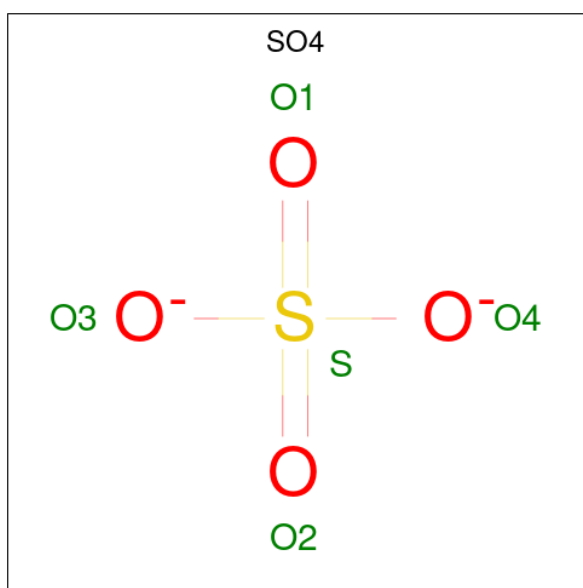
- Molecule 4 is a protein called TCR alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	181	Total	C	N	O	S	0	0	0
			1432	909	231	286	6			

- Molecule 5 is a protein called TCR beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	245	Total	C	N	O	S	0	0	0
			1934	1219	340	365	10			

- Molecule 6 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	E	1	Total	O	S	0	0
			5	4	1		
6	E	1	Total	O	S	0	0
			5	4	1		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	9	Total	O	0	0
			9	9		
7	B	1	Total	O	0	0
			1	1		
7	D	4	Total	O	0	0
			4	4		

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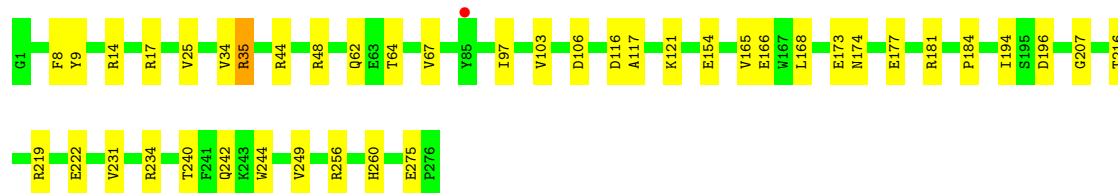
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	E	4	Total	O	0	0
			4	4		

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.

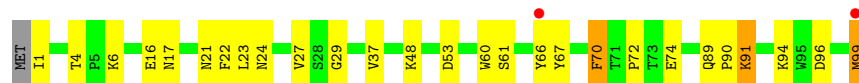
- Molecule 1: MHC class I antigen

Chain A: 



- Molecule 2: Beta-2-microglobulin

Chain B: 



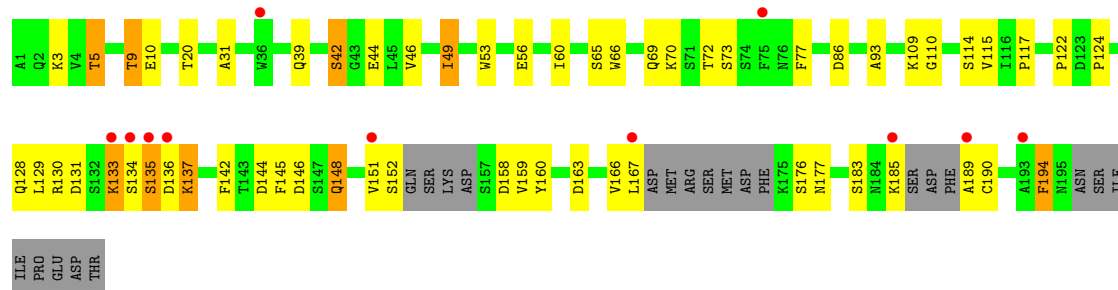
- Molecule 3: KRAS G12D peptide (VVVGADGVGK)

Chain C: 

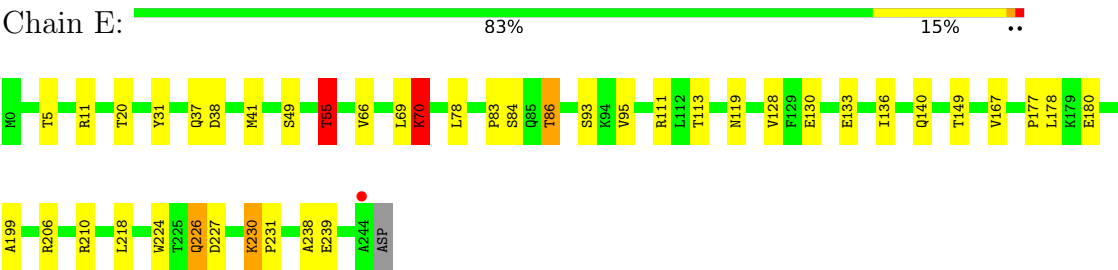


- Molecule 4: TCR alpha

Chain D: 



● Molecule 5: TCR beta



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	144.16Å 144.16Å 119.59Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	64.47 – 2.64 64.47 – 2.64	Depositor EDS
% Data completeness (in resolution range)	100.0 (64.47-2.64) 99.8 (64.47-2.64)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.99 (at 2.65Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.211 , 0.268 (Not available) , 0.233	Depositor DCC
R_{free} test set	1836 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	98.0	Xtriage
Anisotropy	0.026	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 53.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6536	wwPDB-VP
Average B, all atoms (Å ²)	101.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.09% 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

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	1.04	0/2313	1.41	3/3141 (0.1%)
2	B	1.07	0/851	1.32	0/1152
3	C	1.55	0/62	1.86	0/81
4	D	1.11	1/1465 (0.1%)	1.43	5/1990 (0.3%)
5	E	1.10	0/1987	1.42	5/2703 (0.2%)
All	All	1.08	1/6678 (0.0%)	1.41	13/9067 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	93	ALA	C-O	5.19	1.30	1.23

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	135	SER	N-CA-C	-8.97	101.89	112.92
5	E	55	THR	CA-CB-OG1	-6.50	99.84	109.60
5	E	177	PRO	CA-C-O	-6.29	113.87	121.03
1	A	106	ASP	CA-CB-CG	6.26	118.86	112.60
4	D	110	GLY	CA-C-O	-6.06	118.05	122.23
4	D	9	THR	CA-CB-OG1	-5.86	100.82	109.60
4	D	194	PHE	CA-CB-CG	5.69	119.49	113.80
1	A	184	PRO	CB-CA-C	-5.66	105.85	111.17
5	E	70	LYS	CA-C-N	5.35	127.71	120.38
5	E	70	LYS	C-N-CA	5.35	127.71	120.38
5	E	93	SER	CA-C-O	-5.21	115.67	121.51
4	D	5	THR	CB-CA-C	5.16	118.08	110.14
1	A	194	ILE	N-CA-C	-5.03	107.48	111.81

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2251	0	2098	17	0
2	B	828	0	794	14	0
3	C	63	0	69	0	0
4	D	1432	0	1353	23	0
5	E	1934	0	1855	20	0
6	E	10	0	0	0	0
7	A	9	0	0	0	0
7	B	1	0	0	0	0
7	D	4	0	0	0	0
7	E	4	0	0	0	0
All	All	6536	0	6169	69	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 (69) 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:35:ARG:HG2	1:A:48:ARG:HD3	1.64	0.79
5:E:83:PRO:O	5:E:86:THR:HG23	1.88	0.73
5:E:119:ASN:HD21	5:E:226:GLN:CD	2.03	0.66
1:A:177:GLU:OE1	1:A:177:GLU:HA	1.98	0.64
2:B:91:LYS:HA	2:B:91:LYS:HE2	1.80	0.63
5:E:224:TRP:CG	5:E:230:LYS:HG2	2.34	0.62
1:A:17:ARG:HG3	1:A:17:ARG:O	2.05	0.56
4:D:128:GLN:HG2	4:D:130:ARG:NH1	2.22	0.55
1:A:234:ARG:HE	1:A:242:GLN:HE21	1.53	0.55
4:D:69:GLN:HB2	4:D:72:THR:OG1	2.07	0.54
5:E:86:THR:HB	5:E:113:THR:HA	1.90	0.53
2:B:96:ASP:O	2:B:99:MET:HB2	2.09	0.52
5:E:218:LEU:HD12	5:E:231:PRO:HD2	1.92	0.52
4:D:122:PRO:C	4:D:124:PRO:HD3	2.36	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:135:SER:C	4:D:137:LYS:H	2.19	0.51
4:D:145:PHE:CZ	4:D:177:ASN:HB3	2.44	0.51
4:D:135:SER:HB2	4:D:137:LYS:HG3	1.93	0.51
5:E:70:LYS:HZ3	5:E:70:LYS:HB2	1.77	0.50
4:D:160:TYR:HE1	5:E:180:GLU:O	1.94	0.50
1:A:231:VAL:HG13	1:A:244:TRP:CZ2	2.47	0.49
5:E:55:THR:OG1	5:E:66:VAL:HG23	2.12	0.49
4:D:20:THR:HA	4:D:77:PHE:O	2.12	0.49
4:D:42:SER:HB2	4:D:44:GLU:OE1	2.12	0.49
2:B:17:ASN:ND2	2:B:74:GLU:HB2	2.28	0.49
2:B:29:GLY:HA2	2:B:61:SER:OG	2.13	0.49
4:D:39:GLN:HE22	5:E:37:GLN:HE22	1.61	0.49
5:E:78:LEU:HD12	5:E:78:LEU:N	2.28	0.48
4:D:109:LYS:HE3	5:E:41:MET:HA	1.96	0.48
4:D:137:LYS:HE2	4:D:137:LYS:HB2	1.69	0.48
4:D:142:PHE:HB2	4:D:194:PHE:CE2	2.49	0.48
5:E:128:VAL:HG23	5:E:238:ALA:HB3	1.97	0.47
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.50	0.46
2:B:74:GLU:O	2:B:74:GLU:HG2	2.15	0.46
4:D:86:ASP:HB2	4:D:115:VAL:HG21	1.97	0.46
2:B:23:LEU:O	2:B:67:TYR:HA	2.16	0.46
5:E:119:ASN:ND2	5:E:226:GLN:CD	2.73	0.46
4:D:49:ILE:HD13	4:D:66:TRP:CZ3	2.50	0.46
4:D:135:SER:C	4:D:137:LYS:N	2.74	0.46
4:D:133:LYS:O	4:D:134:SER:C	2.59	0.46
1:A:9:TYR:HB2	1:A:97:ILE:HB	1.98	0.46
4:D:159:VAL:HA	4:D:183:SER:HB3	1.98	0.45
5:E:41:MET:HE3	5:E:41:MET:HB3	1.75	0.45
4:D:117:PRO:HG3	4:D:166:VAL:HG13	1.97	0.45
1:A:173:GLU:O	1:A:174:ASN:C	2.60	0.44
2:B:70:PHE:CZ	2:B:72:PRO:HG3	2.53	0.44
2:B:21:ASN:HB3	2:B:70:PHE:CE1	2.51	0.44
5:E:210:ARG:HD3	5:E:239:GLU:OE2	2.18	0.43
1:A:231:VAL:CG1	1:A:244:TRP:CZ2	3.01	0.43
2:B:24:ASN:OD1	2:B:67:TYR:HB3	2.18	0.43
2:B:89:GLN:O	2:B:90:PRO:C	2.62	0.43
1:A:8:PHE:HB2	1:A:25:VAL:CG2	2.49	0.43
1:A:44:ARG:HD2	1:A:64:THR:HG21	2.01	0.43
4:D:129:LEU:HD22	5:E:130:GLU:O	2.19	0.43
4:D:189:ALA:O	4:D:190:CYS:C	2.61	0.43
1:A:116:ASP:N	1:A:116:ASP:OD1	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:216:THR:OG1	1:A:260:HIS:HB2	2.19	0.42
1:A:234:ARG:HE	1:A:242:GLN:NE2	2.17	0.42
4:D:146:ASP:OD2	4:D:148:GLN:HB2	2.19	0.42
5:E:83:PRO:O	5:E:86:THR:CG2	2.63	0.42
5:E:136:ILE:HG23	5:E:199:ALA:HB1	2.01	0.42
1:A:219:ARG:NH2	1:A:256:ARG:HH21	2.18	0.42
2:B:6:LYS:O	2:B:27:VAL:HA	2.20	0.42
2:B:37:VAL:HG11	2:B:66:TYR:CD1	2.56	0.41
5:E:31:TYR:HA	5:E:49:SER:O	2.21	0.41
4:D:31:ALA:HA	4:D:53:TRP:CD2	2.56	0.41
5:E:218:LEU:CD1	5:E:231:PRO:HD2	2.51	0.41
2:B:21:ASN:CG	2:B:22:PHE:H	2.28	0.40
1:A:103:VAL:HG13	1:A:168:LEU:HD23	2.03	0.40
1:A:207:GLY:HA2	1:A:240:THR:HB	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	274/276 (99%)	263 (96%)	11 (4%)	0	100	100
2	B	97/100 (97%)	92 (95%)	5 (5%)	0	100	100
3	C	8/10 (80%)	6 (75%)	2 (25%)	0	100	100
4	D	173/203 (85%)	147 (85%)	26 (15%)	0	100	100
5	E	243/246 (99%)	225 (93%)	18 (7%)	0	100	100
All	All	795/835 (95%)	733 (92%)	62 (8%)	0	100	100

There are no Ramachandran outliers to report.

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	233/233 (100%)	219 (94%)	14 (6%)	17	29
2	B	94/95 (99%)	85 (90%)	9 (10%)	8	12
3	C	6/6 (100%)	4 (67%)	2 (33%)	0	0
4	D	162/184 (88%)	136 (84%)	26 (16%)	2	2
5	E	210/211 (100%)	190 (90%)	20 (10%)	8	12
All	All	705/729 (97%)	634 (90%)	71 (10%)	7	10

All (71) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	ARG
1	A	34	VAL
1	A	35	ARG
1	A	62	GLN
1	A	67	VAL
1	A	121	LYS
1	A	154	GLU
1	A	165	VAL
1	A	166	GLU
1	A	181	ARG
1	A	196	ASP
1	A	222	GLU
1	A	249	VAL
1	A	275	GLU
2	B	1	ILE
2	B	4	THR
2	B	16	GLU
2	B	48	LYS
2	B	53	ASP
2	B	70	PHE
2	B	91	LYS
2	B	94	LYS
2	B	99	MET

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Mol	Chain	Res	Type
3	C	3	VAL
3	C	8	VAL
4	D	3	LYS
4	D	5	THR
4	D	9	THR
4	D	10	GLU
4	D	42	SER
4	D	46	VAL
4	D	49	ILE
4	D	56	GLU
4	D	60	ILE
4	D	65	SER
4	D	70	LYS
4	D	73	SER
4	D	114	SER
4	D	131	ASP
4	D	133	LYS
4	D	136	ASP
4	D	137	LYS
4	D	144	ASP
4	D	148	GLN
4	D	151	VAL
4	D	152	SER
4	D	158	ASP
4	D	163	ASP
4	D	167	LEU
4	D	176	SER
4	D	185	LYS
5	E	5	THR
5	E	11	ARG
5	E	20	THR
5	E	38	ASP
5	E	55	THR
5	E	69	LEU
5	E	70	LYS
5	E	84	SER
5	E	86	THR
5	E	95	VAL
5	E	111	ARG
5	E	133	GLU
5	E	140	GLN
5	E	149	THR

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Mol	Chain	Res	Type
5	E	167	VAL
5	E	178	LEU
5	E	206	ARG
5	E	226	GLN
5	E	227	ASP
5	E	230	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	GLN
1	A	62	GLN
1	A	180	GLN
1	A	192	HIS
1	A	218	GLN
1	A	242	GLN
1	A	262	GLN
2	B	51	HIS
4	D	57	GLN
4	D	58	ASN
4	D	67	ASN
4	D	118	ASN
4	D	148	GLN
4	D	192	ASN
5	E	28	ASN
5	E	37	GLN
5	E	73	ASN
5	E	119	ASN
5	E	203	GLN

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	$\# Z > 2$
6	SO4	E	302	-	4,4,4	0.33	0	6,6,6	0.12	0
6	SO4	E	301	-	4,4,4	0.33	0	6,6,6	0.21	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 [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%)	-0.14	1 (0%) 88 87	72, 93, 122, 153	0
2	B	99/100 (99%)	0.17	2 (2%) 65 61	83, 118, 145, 157	0
3	C	10/10 (100%)	-0.01	0 100 100	74, 77, 87, 89	0
4	D	181/203 (89%)	0.41	11 (6%) 27 22	30, 113, 159, 180	0
5	E	245/246 (99%)	-0.09	1 (0%) 88 87	69, 92, 121, 138	0
All	All	811/835 (97%)	0.04	15 (1%) 67 64	30, 97, 140, 180	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	D	134	SER	7.7
5	E	244	ALA	4.3
4	D	189	ALA	3.9
4	D	135	SER	3.2
4	D	185	LYS	3.2
4	D	133	LYS	3.0
2	B	66	TYR	2.8
4	D	36	TRP	2.4
4	D	136	ASP	2.4
4	D	167	LEU	2.4
1	A	85	TYR	2.4
4	D	75	PHE	2.4
4	D	151	VAL	2.3
2	B	99	MET	2.2
4	D	193	ALA	2.1

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
6	SO4	E	301	5/5	0.82	0.12	122,128,137,138	0
6	SO4	E	302	5/5	0.86	0.08	115,121,121,124	0

6.5 Other polymers [i](#)

There are no such residues in this entry.