



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 10, 2026 – 11:42 pm BST

PDB ID : 9SOK / pdb_00009sok
Title : Structure of three-domain single chain TCR-581 in complex with Peptide-HLA
Authors : Karuppiah, V.
Deposited on : 2025-09-14
Resolution : 2.50 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	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

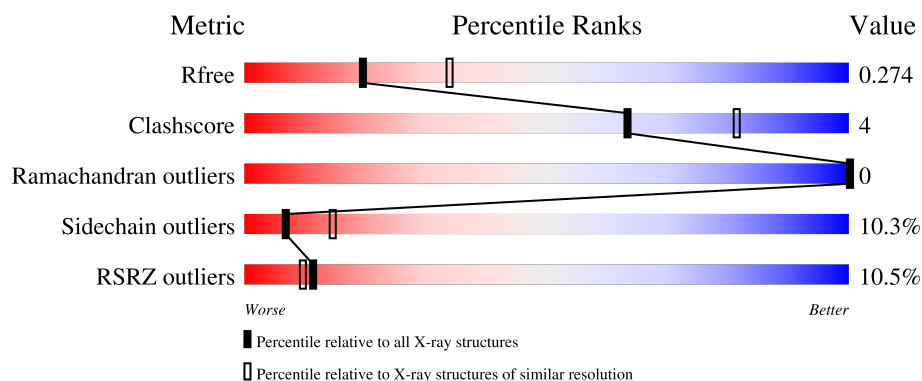
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	4% 84% 12% ..
2	B	100	5% 78% 21% .
3	C	9	89% 11%
4	E	382	16% 71% 19% .. 8%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6054 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	275	Total	C	N	O	S	0	0	0
			2250	1406	409	426	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			

There is a discrepancy between the modelled and reference sequences:

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

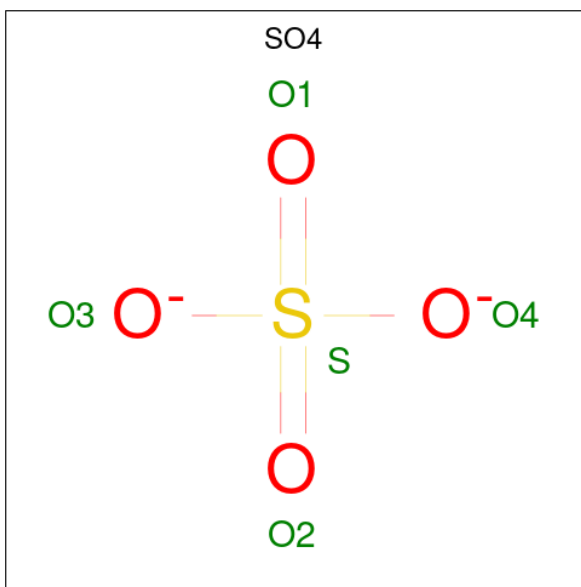
- Molecule 3 is a protein called Piwi-like protein 1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	9	Total	C	N	O	0	0	0
			80	52	13	15			

- Molecule 4 is a protein called TCR (3-domain single chain).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	E	352	Total	C	N	O	S	0	0	0
			2794	1754	492	536	12			

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

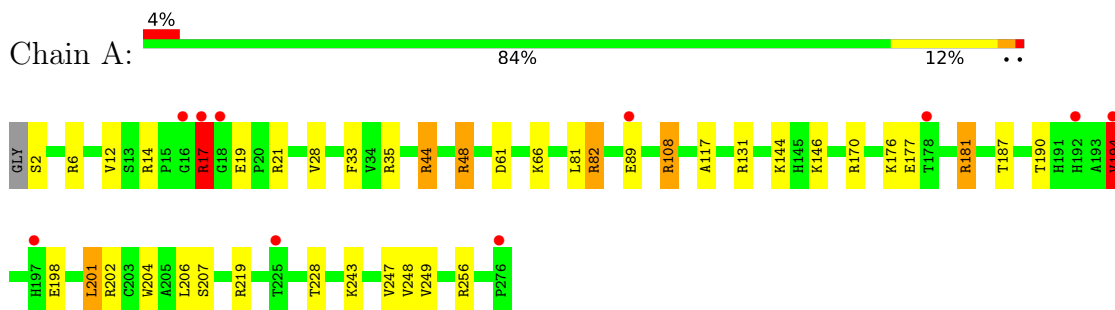
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	31	Total	O	0	0
			31	31		
6	B	10	Total	O	0	0
			10	10		
6	C	5	Total	O	0	0
			5	5		
6	E	32	Total	O	0	0
			32	32		

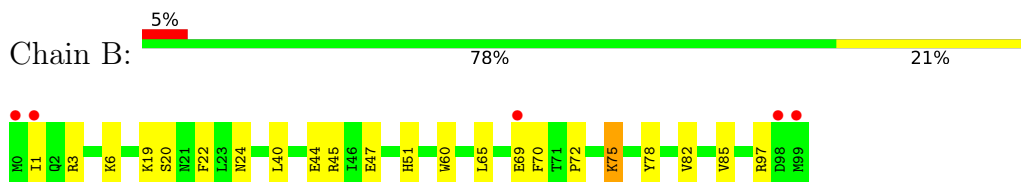
3 Residue-property plots [i](#)

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.

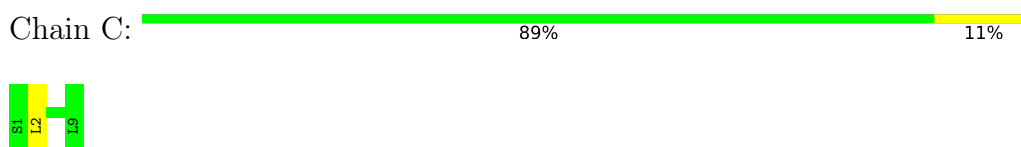
- Molecule 1: MHC class I antigen



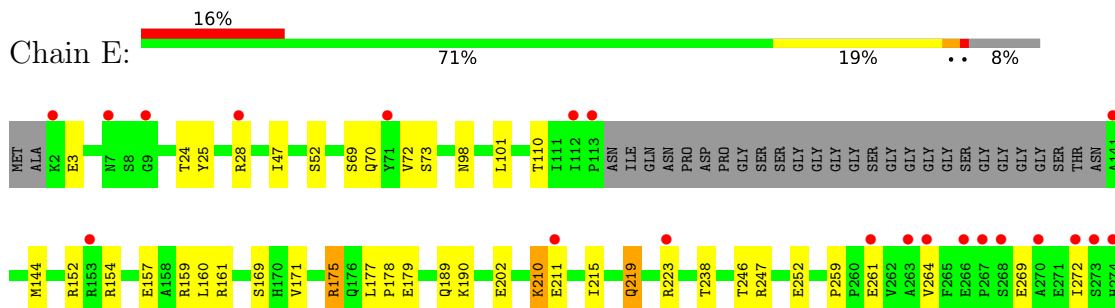
- Molecule 2: Beta-2-microglobulin

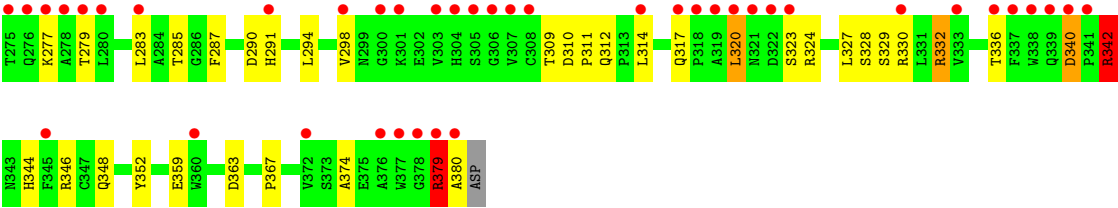


- Molecule 3: Piwi-like protein 1



- Molecule 4: TCR (3-domain single chain)





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	56.76Å 83.70Å 115.37Å 90.00° 99.60° 90.00°	Depositor
Resolution (Å)	67.42 – 2.50 67.42 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.9 (67.42-2.50) 99.9 (67.42-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.86 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.8.0430	Depositor
R, R_{free}	0.208 , 0.262 0.231 , 0.274	Depositor DCC
R_{free} test set	1723 reflections (4.65%)	wwPDB-VP
Wilson B-factor (Å ²)	45.1	Xtriage
Anisotropy	0.582	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 54.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6054	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.93% 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	0.73	0/2316	1.24	4/3144 (0.1%)
2	B	0.71	0/860	1.12	2/1162 (0.2%)
3	C	0.77	0/81	1.31	0/107
4	E	0.74	0/2864	1.17	3/3883 (0.1%)
All	All	0.73	0/6121	1.19	9/8296 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	10
2	B	0	1
4	E	0	9
All	All	0	20

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	249	VAL	N-CA-CB	-8.20	106.17	111.83
1	A	33	PHE	CA-CB-CG	7.32	121.12	113.80
2	B	51	HIS	CA-CB-CG	6.41	120.21	113.80
2	B	70	PHE	CA-CB-CG	5.86	119.66	113.80
1	A	194	VAL	N-CA-C	-5.67	106.94	111.81
4	E	189	GLN	N-CA-C	-5.31	102.34	110.14
4	E	287	PHE	CA-CB-CG	5.16	118.96	113.80
1	A	28	VAL	N-CA-CB	5.11	118.34	111.90
4	E	261	GLU	CB-CA-C	5.04	118.07	109.75

There are no chirality outliers.

All (20) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	108	ARG	Sidechain
1	A	131	ARG	Sidechain
1	A	17	ARG	Sidechain
1	A	170	ARG	Sidechain
1	A	181	ARG	Sidechain
1	A	21	ARG	Sidechain
1	A	44	ARG	Sidechain
1	A	48	ARG	Sidechain
1	A	6	ARG	Sidechain
1	A	82	ARG	Sidechain
2	B	97	ARG	Sidechain
4	E	154	ARG	Sidechain
4	E	159	ARG	Sidechain
4	E	175	ARG	Sidechain
4	E	223	ARG	Sidechain
4	E	324	ARG	Sidechain
4	E	330	ARG	Sidechain
4	E	332	ARG	Sidechain
4	E	342	ARG	Sidechain
4	E	379	ARG	Sidechain

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2250	0	2097	13	0
2	B	837	0	803	10	0
3	C	80	0	82	1	0
4	E	2794	0	2679	27	0
5	A	10	0	0	0	0
5	C	5	0	0	0	0
6	A	31	0	0	0	0
6	B	10	0	0	0	0
6	C	5	0	0	0	0
6	E	32	0	0	0	0
All	All	6054	0	5661	47	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (47) 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)
4:E:346:ARG:NH2	4:E:348:GLN:OE1	2.26	0.67
4:E:161:ARG:HG2	4:E:215:ILE:CD1	2.35	0.56
4:E:320:LEU:HB3	4:E:323:SER:HB2	1.89	0.55
2:B:19:LYS:O	2:B:72:PRO:HD2	2.08	0.54
4:E:178:PRO:O	4:E:179:GLU:HB2	2.10	0.51
1:A:219:ARG:HH11	1:A:256:ARG:HH22	1.59	0.50
1:A:146:LYS:HD2	4:E:238:THR:HG23	1.94	0.49
4:E:310:ASP:HB2	4:E:327:LEU:HD12	1.94	0.49
4:E:340:ASP:C	4:E:342:ARG:H	2.20	0.49
1:A:17:ARG:NH2	1:A:19:GLU:HB3	2.28	0.48
4:E:157:GLU:HG3	4:E:219:GLN:HE21	1.79	0.48
1:A:187:THR:HA	1:A:204:TRP:O	2.14	0.48
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.49	0.47
1:A:44:ARG:NH2	1:A:61:ASP:OD1	2.48	0.47
4:E:379:ARG:O	4:E:380:ALA:C	2.57	0.46
4:E:25:TYR:CZ	4:E:70:GLN:HA	2.50	0.46
4:E:259:PRO:HD3	4:E:367:PRO:HB3	1.97	0.46
4:E:340:ASP:N	4:E:340:ASP:OD1	2.49	0.46
1:A:117:ALA:HB2	2:B:60:TRP:CD2	2.51	0.46
4:E:160:LEU:HD22	4:E:246:THR:HG21	1.98	0.46
4:E:264:VAL:HG23	4:E:374:ALA:HB3	1.98	0.45
4:E:298:VAL:HA	4:E:344:HIS:O	2.17	0.45
2:B:40:LEU:O	2:B:78:TYR:HA	2.16	0.45
2:B:69:GLU:CD	2:B:69:GLU:H	2.24	0.45
2:B:24:ASN:HB3	2:B:65:LEU:HD11	1.98	0.44
4:E:359:GLU:H	4:E:359:GLU:HG2	1.54	0.44
2:B:75:LYS:H	2:B:75:LYS:HG3	1.28	0.43
4:E:175:ARG:HD3	4:E:177:LEU:HG	2.00	0.43
4:E:379:ARG:H	4:E:379:ARG:HG3	1.60	0.43
1:A:201:LEU:HD12	1:A:201:LEU:HA	1.81	0.43
4:E:320:LEU:HD22	4:E:320:LEU:HA	1.93	0.43
4:E:190:LYS:HD3	4:E:210:LYS:HA	2.01	0.43
4:E:294:LEU:C	4:E:294:LEU:HD23	2.44	0.42
1:A:194:VAL:HG13	1:A:198:GLU:HG3	2.01	0.42
2:B:40:LEU:HA	2:B:44:GLU:O	2.19	0.42
4:E:310:ASP:O	4:E:311:PRO:C	2.61	0.42
4:E:309:THR:HG23	4:E:329:SER:HB2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:190:THR:OG1	1:A:202:ARG:HB3	2.19	0.42
4:E:283:LEU:HG	4:E:285:THR:HG23	2.01	0.41
4:E:98:ASN:OD1	4:E:98:ASN:C	2.63	0.41
4:E:161:ARG:HG2	4:E:215:ILE:HD13	2.03	0.41
1:A:66:LYS:HE3	3:C:2:LEU:O	2.21	0.41
1:A:219:ARG:NH1	1:A:256:ARG:HH22	2.19	0.41
1:A:48:ARG:HD3	1:A:48:ARG:HA	1.90	0.40
2:B:22:PHE:CE2	2:B:69:GLU:HG3	2.56	0.40
2:B:22:PHE:CZ	2:B:69:GLU:HG3	2.57	0.40
4:E:291:HIS:HB3	4:E:352:TYR:HB2	2.02	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	273/276 (99%)	264 (97%)	9 (3%)	0	100	100
2	B	98/100 (98%)	96 (98%)	2 (2%)	0	100	100
3	C	7/9 (78%)	7 (100%)	0	0	100	100
4	E	348/382 (91%)	330 (95%)	18 (5%)	0	100	100
All	All	726/767 (95%)	697 (96%)	29 (4%)	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	232/232 (100%)	211 (91%)	21 (9%)	9	19
2	B	95/95 (100%)	86 (90%)	9 (10%)	8	17
3	C	9/9 (100%)	9 (100%)	0	100	100
4	E	304/320 (95%)	268 (88%)	36 (12%)	5	11
All	All	640/656 (98%)	574 (90%)	66 (10%)	7	14

All (66) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	SER
1	A	12	VAL
1	A	14	ARG
1	A	17	ARG
1	A	35	ARG
1	A	81	LEU
1	A	82	ARG
1	A	89	GLU
1	A	108	ARG
1	A	144	LYS
1	A	176	LYS
1	A	177	GLU
1	A	181	ARG
1	A	194	VAL
1	A	201	LEU
1	A	206	LEU
1	A	207	SER
1	A	228	THR
1	A	243	LYS
1	A	247	VAL
1	A	248	VAL
2	B	1	ILE
2	B	3	ARG
2	B	6	LYS
2	B	20	SER
2	B	45	ARG
2	B	47	GLU
2	B	75	LYS
2	B	82	VAL
2	B	85	VAL

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Mol	Chain	Res	Type
4	E	3	GLU
4	E	24	THR
4	E	28	ARG
4	E	47	ILE
4	E	52	SER
4	E	69	SER
4	E	72	VAL
4	E	73	SER
4	E	101	LEU
4	E	110	THR
4	E	144	MET
4	E	152	ARG
4	E	169	SER
4	E	171	VAL
4	E	202	GLU
4	E	210	LYS
4	E	211	GLU
4	E	219	GLN
4	E	247	ARG
4	E	252	GLU
4	E	269	GLU
4	E	272	ILE
4	E	277	LYS
4	E	279	THR
4	E	290	ASP
4	E	312	GLN
4	E	314	LEU
4	E	317	GLN
4	E	320	LEU
4	E	328	SER
4	E	332	ARG
4	E	336	THR
4	E	340	ASP
4	E	342	ARG
4	E	363	ASP
4	E	379	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	224	GLN
4	E	64	GLN

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Mol	Chain	Res	Type
4	E	219	GLN
4	E	276	GLN
4	E	291	HIS

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](#)

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	$\# Z > 2$
5	SO4	A	302	-	4,4,4	0.44	0	6,6,6	0.05	0
5	SO4	A	301	-	4,4,4	0.32	0	6,6,6	0.11	0
5	SO4	C	101	-	4,4,4	0.38	0	6,6,6	0.22	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	275/276 (99%)	0.38	10 (3%) 46 41	27, 42, 76, 106	0
2	B	100/100 (100%)	0.51	5 (5%) 34 30	31, 51, 84, 98	0
3	C	9/9 (100%)	0.06	0 100 100	30, 34, 43, 44	0
4	E	352/382 (92%)	0.94	62 (17%) 4 3	30, 55, 124, 143	0
All	All	736/767 (95%)	0.67	77 (10%) 11 10	27, 49, 105, 143	0

All (77) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	18	GLY	5.4
4	E	380	ALA	4.9
1	A	197	HIS	4.7
4	E	141	ALA	4.6
4	E	322	ASP	4.4
4	E	211	GLU	4.3
4	E	319	ALA	4.1
4	E	377	TRP	4.1
4	E	376	ALA	4.0
4	E	318	PRO	3.9
4	E	279	THR	3.9
4	E	274	HIS	3.9
4	E	338	TRP	3.8
4	E	333	VAL	3.8
4	E	321	ASN	3.8
4	E	278	ALA	3.7
4	E	275	THR	3.7
4	E	71	TYR	3.6
4	E	267	PRO	3.6
4	E	304	HIS	3.5
4	E	303	VAL	3.4

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Mol	Chain	Res	Type	RSRZ
4	E	314	LEU	3.4
2	B	1	ILE	3.4
4	E	272	ILE	3.4
4	E	112	ILE	3.4
2	B	99	MET	3.4
4	E	277	LYS	3.4
4	E	261	GLU	3.4
4	E	2	LYS	3.3
4	E	330	ARG	3.3
4	E	291	HIS	3.2
1	A	16	GLY	3.2
4	E	379	ARG	3.2
2	B	0	MET	3.2
4	E	307	VAL	3.2
4	E	320	LEU	3.2
4	E	113	PRO	3.1
1	A	178	THR	3.1
4	E	317	GLN	3.0
4	E	336	THR	3.0
1	A	17	ARG	2.9
1	A	194	VAL	2.9
4	E	341	PRO	2.9
1	A	192	HIS	2.8
1	A	225	THR	2.8
4	E	28	ARG	2.8
4	E	372	VAL	2.7
4	E	273	SER	2.7
1	A	276	PRO	2.6
4	E	223	ARG	2.6
4	E	270	ALA	2.6
2	B	98	ASP	2.6
4	E	280	LEU	2.6
4	E	300	GLY	2.5
4	E	283	LEU	2.5
4	E	305	SER	2.5
4	E	378	GLY	2.5
4	E	339	GLN	2.4
4	E	345	PHE	2.4
4	E	263	ALA	2.4
4	E	308	CYS	2.3
4	E	337	PHE	2.3
4	E	298	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
4	E	360	TRP	2.3
2	B	69	GLU	2.3
1	A	89	GLU	2.3
4	E	266	GLU	2.3
4	E	306	GLY	2.2
4	E	340	ASP	2.2
4	E	9	GLY	2.2
4	E	264	VAL	2.1
4	E	153	ARG	2.1
4	E	323	SER	2.1
4	E	301	LYS	2.1
4	E	276	GLN	2.1
4	E	7	ASN	2.0
4	E	268	SER	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	SO4	A	302	5/5	0.82	0.17	87,100,117,126	0
5	SO4	A	301	5/5	0.83	0.14	78,97,102,112	0
5	SO4	C	101	5/5	0.87	0.18	56,60,69,86	0

6.5 Other polymers [i](#)

There are no such residues in this entry.