



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 14, 2017 – 02:01 AM EST

PDB ID : 5XOT
Title : Crystal structure of pHLA-B35 in complex with TU55 T cell receptor
Authors : Shi, Y.; Qi, J.; Gao, G.F.
Deposited on : unknown
Resolution : 2.79 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20030345
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20030345

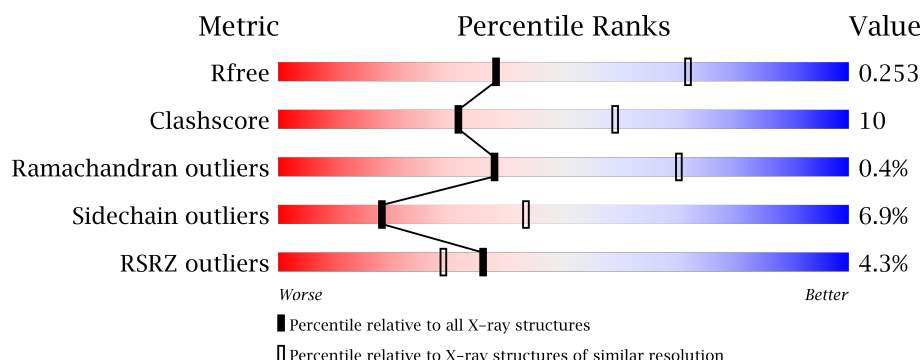
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.79 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}	100719	3276 (2.80-2.76)
Clashscore	112137	3771 (2.80-2.76)
Ramachandran outliers	110173	3707 (2.80-2.76)
Sidechain outliers	110143	3709 (2.80-2.76)
RSRZ outliers	101464	3307 (2.80-2.76)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	<div> <div>4%</div> <div>80%</div> <div>18%</div> <div>.</div> </div>
2	B	99	<div> <div>6%</div> <div>77%</div> <div>21%</div> <div>.</div> </div>
3	C	9	<div> <div>56%</div> <div>22%</div> <div>22%</div> </div>
4	D	204	<div> <div>9%</div> <div>74%</div> <div>23%</div> <div>.</div> </div>
5	E	242	<div> <div>%</div> <div>74%</div> <div>20%</div> <div>6%</div> </div>

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
6	GOL	A	301	-	-	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 6728 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 HLA class I histocompatibility antigen, B-35 alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	276	Total	C	N	O	S	0	0	0
			2255	1404	411	433	7			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	34	ASP	VAL	engineered mutation	UNP P30685

- 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
			829	528	140	158	3			

- Molecule 3 is a protein called An HIV reverse transcriptase epitope.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	9	Total	C	N	O	0	0	0
			71	45	9	17			

- Molecule 4 is a protein called The Delta chain of TU55 TCR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	203	Total	C	N	O	S	0	0	0
			1598	1010	266	313	9			

- Molecule 5 is a protein called The beta chain of TU55 TCR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	242	Total	C	N	O	S	0	0	0
			1924	1207	339	369	9			

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			6	3	3		
6	B	1	Total	C	O	0	0
			6	3	3		
6	B	1	Total	C	O	0	0
			6	3	3		

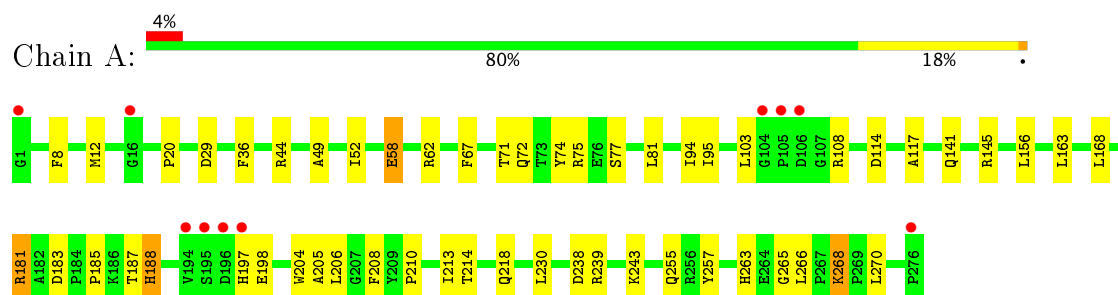
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	11	Total	O	0	0
			11	11		
7	B	1	Total	O	0	0
			1	1		
7	D	7	Total	O	0	0
			7	7		
7	E	14	Total	O	0	0
			14	14		

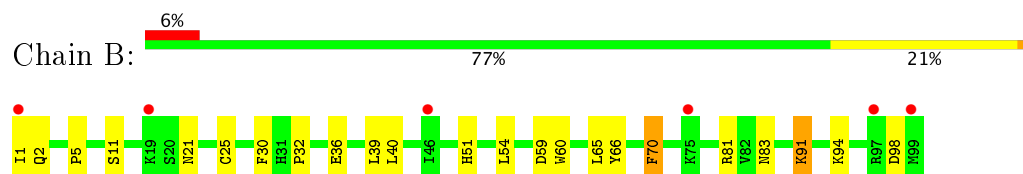
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: HLA class I histocompatibility antigen, B-35 alpha chain



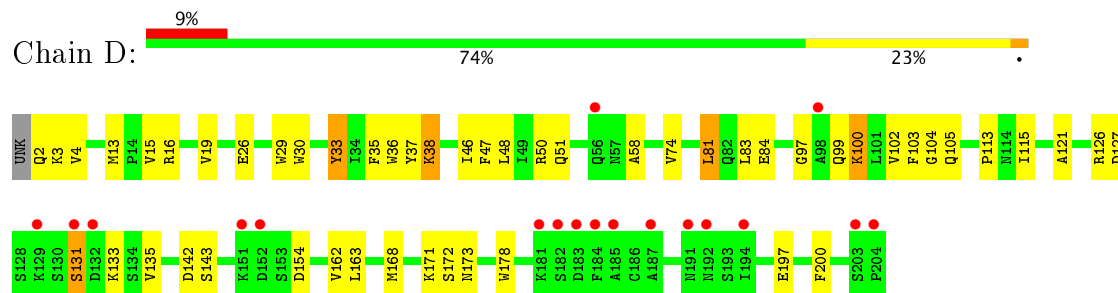
- Molecule 2: Beta-2-microglobulin



- Molecule 3: An HIV reverse transcriptase epitope

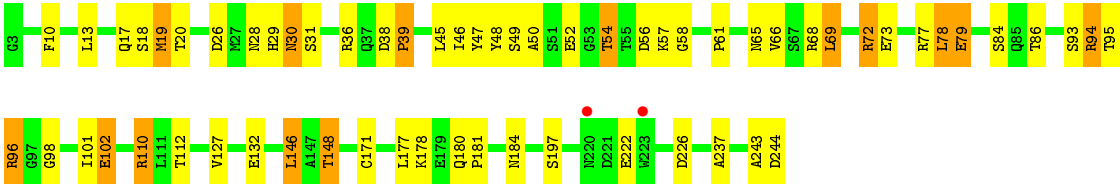


- Molecule 4: The Delta chain of TU55 TCR



- Molecule 5: The beta chain of TU55 TCR





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	195.69Å 59.60Å 92.81Å 90.00° 105.98° 90.00°	Depositor
Resolution (Å)	36.60 – 2.79 42.14 – 2.78	Depositor EDS
% Data completeness (in resolution range)	90.5 (36.60-2.79) 93.5 (42.14-2.78)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.50 (at 2.77Å)	Xtriage
Refinement program	PHENIX 1.5_2	Depositor
R, R_{free}	0.199 , 0.254 0.195 , 0.253	Depositor DCC
R_{free} test set	1255 reflections (5.13%)	DCC
Wilson B-factor (Å ²)	45.7	Xtriage
Anisotropy	0.937	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 45.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6728	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.78% 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: GOL

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	0.22	0/2318	0.42	1/3151 (0.0%)
2	B	0.23	0/852	0.38	0/1152
3	C	0.25	0/71	0.49	0/95
4	D	0.41	0/1633	0.46	0/2210
5	E	0.31	0/1974	0.45	1/2684 (0.0%)
All	All	0.30	0/6848	0.43	2/9292 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	29	ASP	N-CA-C	-6.49	93.48	111.00
5	E	79	GLU	N-CA-C	-5.28	96.75	111.00

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	2255	0	2112	36	0
2	B	829	0	794	14	0
3	C	71	0	72	5	0
4	D	1598	0	1550	48	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	E	1924	0	1835	50	0
6	A	6	0	8	1	0
6	B	12	0	16	2	0
7	A	11	0	0	0	0
7	B	1	0	0	0	0
7	D	7	0	0	0	0
7	E	14	0	0	0	0
All	All	6728	0	6387	134	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 10.

All (134) 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)
5:E:30:ASN:ND2	5:E:30:ASN:H	1.75	0.84
4:D:102:VAL:HG22	4:D:102:VAL:O	1.75	0.84
5:E:30:ASN:H	5:E:30:ASN:HD22	1.24	0.83
4:D:99:GLN:HG2	5:E:94:ARG:HH21	1.43	0.82
4:D:135:VAL:HG23	4:D:178:TRP:HB3	1.61	0.81
4:D:102:VAL:CG2	4:D:102:VAL:O	2.30	0.79
5:E:30:ASN:ND2	5:E:30:ASN:N	2.30	0.76
5:E:13:LEU:HD11	5:E:19:MET:HE2	1.67	0.76
1:A:20:PRO:HB2	1:A:75:ARG:HG2	1.70	0.74
4:D:162:VAL:HG22	4:D:173:ASN:ND2	2.02	0.74
5:E:17:GLN:O	5:E:79:GLU:O	2.06	0.73
4:D:4:VAL:HG23	4:D:102:VAL:CG2	2.20	0.72
1:A:263:HIS:CD2	1:A:265:GLY:H	2.06	0.72
5:E:18:SER:HA	5:E:79:GLU:O	1.90	0.71
1:A:230:LEU:HD22	1:A:243:LYS:HE3	1.72	0.70
1:A:49:ALA:O	1:A:52:ILE:HG22	1.93	0.69
4:D:4:VAL:HG23	4:D:102:VAL:HG21	1.75	0.68
4:D:4:VAL:CG2	4:D:102:VAL:CG2	2.72	0.67
5:E:94:ARG:NH1	5:E:96:ARG:HG2	2.13	0.64
4:D:113:PRO:HG3	4:D:162:VAL:HG11	1.81	0.63
4:D:163:LEU:HB3	5:E:171:CYS:HB2	1.79	0.63
5:E:49:SER:HB2	5:E:54:THR:O	1.98	0.63
4:D:113:PRO:HD2	4:D:143:SER:OG	1.99	0.62
1:A:58:GLU:O	1:A:62:ARG:HD3	2.00	0.62
4:D:15:VAL:O	4:D:16:ARG:HB2	2.00	0.61
5:E:46:ILE:HG22	5:E:47:TYR:HD2	1.65	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:50:ALA:HB1	5:E:96:ARG:HH21	1.65	0.60
3:C:4:THR:HG22	4:D:97:GLY:H	1.66	0.59
1:A:77:SER:HB3	3:C:9:LEU:HB2	1.85	0.58
4:D:13:MET:HG3	4:D:19:VAL:HB	1.86	0.58
5:E:69:LEU:HD11	5:E:73:GLU:HB3	1.86	0.58
5:E:46:ILE:HG22	5:E:47:TYR:CD2	2.39	0.58
1:A:238:ASP:O	1:A:239:ARG:HB2	2.04	0.57
4:D:100:LYS:HB3	5:E:45:LEU:HD22	1.85	0.57
5:E:65:ASN:OD1	5:E:77:ARG:HD3	2.04	0.57
5:E:36:ARG:HB2	5:E:46:ILE:HD11	1.87	0.57
1:A:187:THR:O	1:A:188:HIS:HB3	2.04	0.56
4:D:178:TRP:CZ3	5:E:146:LEU:HD21	2.39	0.56
5:E:57:LYS:HB3	5:E:58:GLY:HA3	1.87	0.56
4:D:4:VAL:HG21	4:D:102:VAL:HG22	1.86	0.56
1:A:263:HIS:HD2	1:A:265:GLY:H	1.48	0.56
1:A:268:LYS:HD2	1:A:268:LYS:O	2.06	0.55
4:D:51:GLN:HE22	4:D:58:ALA:H	1.54	0.55
5:E:146:LEU:CD1	5:E:148:THR:HG23	2.37	0.55
2:B:11:SER:OG	6:B:1101:GOL:H31	2.07	0.55
4:D:127:ASP:HB3	4:D:131:SER:HA	1.90	0.54
5:E:38:ASP:HB3	5:E:39:PRO:HD2	1.88	0.54
1:A:197:HIS:HB2	1:A:198:GLU:OE1	2.07	0.54
2:B:1:ILE:HG13	2:B:2:GLN:HG3	1.89	0.54
5:E:86:THR:HG23	5:E:112:THR:HA	1.88	0.54
1:A:210:PRO:O	1:A:263:HIS:HE1	1.92	0.53
4:D:38:LYS:HG2	4:D:38:LYS:O	2.09	0.53
4:D:162:VAL:HG22	4:D:173:ASN:HD21	1.70	0.52
4:D:115:ILE:HG13	4:D:142:ASP:HA	1.92	0.52
1:A:114:ASP:OD1	1:A:156:LEU:HD21	2.10	0.51
1:A:163:LEU:HD11	4:D:30:TRP:HB2	1.92	0.51
1:A:62:ARG:HG3	4:D:29:TRP:CZ2	2.46	0.51
5:E:72:ARG:O	5:E:72:ARG:HG3	2.10	0.51
5:E:52:GLU:HA	5:E:68:ARG:HG2	1.93	0.50
5:E:30:ASN:HD21	5:E:95:THR:HA	1.77	0.50
1:A:62:ARG:CG	4:D:29:TRP:CZ2	2.95	0.50
4:D:47:PHE:CZ	5:E:102:GLU:HG2	2.47	0.50
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.46	0.50
5:E:10:PHE:CE1	5:E:110:ARG:HD3	2.47	0.49
2:B:21:ASN:HB3	2:B:70:PHE:CE1	2.48	0.49
2:B:59:ASP:O	2:B:60:TRP:HB2	2.12	0.49
4:D:4:VAL:HG21	4:D:102:VAL:CG2	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:4:VAL:CG2	4:D:102:VAL:HG22	2.41	0.49
4:D:103:PHE:C	4:D:104:GLY:O	2.48	0.49
5:E:177:LEU:HD12	5:E:177:LEU:O	2.11	0.49
1:A:71:THR:O	1:A:75:ARG:HG3	2.12	0.49
5:E:94:ARG:HH12	5:E:96:ARG:HG2	1.77	0.49
4:D:38:LYS:HD2	4:D:48:LEU:HD21	1.96	0.48
2:B:5:PRO:HB3	2:B:30:PHE:HB3	1.95	0.48
4:D:100:LYS:HB3	5:E:45:LEU:CD2	2.43	0.48
4:D:163:LEU:H	4:D:163:LEU:HD23	1.78	0.48
2:B:2:GLN:HG2	2:B:32:PRO:HD3	1.96	0.48
1:A:8:PHE:CZ	6:A:301:GOL:H31	2.48	0.47
5:E:98:GLY:O	5:E:101:ILE:HD11	2.14	0.47
2:B:54:LEU:HD22	6:B:1100:GOL:H32	1.96	0.47
1:A:36:PHE:CD1	1:A:67:PHE:HB3	2.49	0.47
5:E:127:VAL:HG23	5:E:237:ALA:HB3	1.96	0.47
5:E:52:GLU:HA	5:E:68:ARG:CG	2.44	0.47
5:E:146:LEU:HD12	5:E:148:THR:HG23	1.96	0.47
4:D:3:LYS:HD2	4:D:26:GLU:O	2.15	0.47
4:D:99:GLN:HG2	5:E:94:ARG:NH2	2.22	0.46
2:B:40:LEU:HD11	2:B:81:ARG:HB2	1.97	0.46
1:A:77:SER:O	1:A:81:LEU:HD23	2.15	0.46
1:A:185:PRO:HB3	1:A:208:PHE:HB3	1.96	0.46
5:E:20:THR:HG22	5:E:77:ARG:HG3	1.98	0.46
5:E:29:HIS:HD2	5:E:93:SER:OG	1.99	0.46
1:A:103:LEU:HG	1:A:168:LEU:HD23	1.98	0.45
5:E:101:ILE:O	5:E:101:ILE:HG22	2.16	0.45
1:A:141:GLN:O	1:A:145:ARG:HG3	2.17	0.45
4:D:100:LYS:HD3	5:E:48:TYR:CE1	2.52	0.45
5:E:243:ALA:O	5:E:244:ASP:HB2	2.16	0.45
4:D:38:LYS:HB3	4:D:48:LEU:HD11	1.99	0.44
1:A:74:TYR:CE2	3:C:6:GLU:HG2	2.52	0.44
5:E:29:HIS:CD2	5:E:93:SER:OG	2.70	0.44
1:A:58:GLU:HG3	1:A:58:GLU:H	1.53	0.44
1:A:95:ILE:HD13	3:C:9:LEU:HD13	1.99	0.44
5:E:36:ARG:NH2	5:E:84:SER:O	2.50	0.44
3:C:1:ILE:HD12	3:C:1:ILE:HA	1.85	0.44
1:A:62:ARG:HG2	4:D:29:TRP:CZ2	2.53	0.43
5:E:78:LEU:O	5:E:79:GLU:HB2	2.18	0.43
1:A:213:ILE:HG12	1:A:214:THR:N	2.33	0.43
1:A:205:ALA:O	1:A:206:LEU:HD23	2.19	0.43
4:D:127:ASP:HB2	4:D:133:LYS:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:121:ALA:HB2	4:D:200:PHE:HB3	2.01	0.43
2:B:51:HIS:HA	2:B:65:LEU:O	2.18	0.43
4:D:74:VAL:O	4:D:74:VAL:HG22	2.19	0.43
2:B:91:LYS:HB3	2:B:91:LYS:HE3	1.46	0.42
4:D:163:LEU:HD23	4:D:172:SER:O	2.20	0.42
5:E:50:ALA:O	5:E:96:ARG:NH2	2.52	0.42
1:A:94:ILE:HD12	1:A:94:ILE:N	2.35	0.42
2:B:94:LYS:HE2	2:B:94:LYS:HB2	1.84	0.42
4:D:100:LYS:HD2	5:E:45:LEU:HD23	2.01	0.42
4:D:37:TYR:CE2	4:D:47:PHE:HB2	2.55	0.42
5:E:26:ASP:HA	5:E:72:ARG:NH1	2.34	0.42
1:A:181:ARG:HG2	1:A:181:ARG:O	2.18	0.41
5:E:178:LYS:HD3	5:E:181:PRO:HA	2.02	0.41
4:D:81:LEU:HA	4:D:81:LEU:HD12	1.88	0.41
1:A:204:TRP:HE3	1:A:206:LEU:HD21	1.86	0.41
2:B:25:CYS:HB2	2:B:39:LEU:HD21	2.02	0.41
4:D:33:TYR:HB3	4:D:35:PHE:CE1	2.56	0.41
2:B:51:HIS:HB3	2:B:66:TYR:CD2	2.56	0.40
4:D:36:TRP:O	4:D:48:LEU:HB2	2.20	0.40
4:D:168:MET:CE	5:E:197:SER:HB3	2.52	0.40
5:E:69:LEU:CD1	5:E:73:GLU:HB3	2.50	0.40
1:A:266:LEU:HD13	1:A:270:LEU:HG	2.04	0.40
1:A:218:GLN:O	1:A:257:TYR:HA	2.21	0.40
1:A:238:ASP:O	1:A:239:ARG:CB	2.69	0.40
5:E:47:TYR:CZ	5:E:61:PRO:HB3	2.57	0.40
4:D:197:GLU:H	4:D:197:GLU:HG3	1.65	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	274/276 (99%)	263 (96%)	10 (4%)	1 (0%)	38	70
2	B	97/99 (98%)	94 (97%)	3 (3%)	0	100	100
3	C	7/9 (78%)	7 (100%)	0	0	100	100
4	D	201/204 (98%)	183 (91%)	17 (8%)	1 (0%)	32	65
5	E	240/242 (99%)	227 (95%)	12 (5%)	1 (0%)	38	70
All	All	819/830 (99%)	774 (94%)	42 (5%)	3 (0%)	38	70

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	D	131	SER
1	A	188	HIS
5	E	39	PRO

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	234/234 (100%)	225 (96%)	9 (4%)	38	71
2	B	94/94 (100%)	89 (95%)	5 (5%)	26	57
3	C	8/8 (100%)	6 (75%)	2 (25%)	1	2
4	D	180/180 (100%)	167 (93%)	13 (7%)	17	41
5	E	210/210 (100%)	189 (90%)	21 (10%)	9	24
All	All	726/726 (100%)	676 (93%)	50 (7%)	18	43

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	MET
1	A	44	ARG
1	A	58	GLU
1	A	72	GLN
1	A	108	ARG

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Mol	Chain	Res	Type
1	A	181	ARG
1	A	183	ASP
1	A	255	GLN
1	A	268	LYS
2	B	36	GLU
2	B	70	PHE
2	B	83	ASN
2	B	91	LYS
2	B	98	ASP
3	C	1	ILE
3	C	9	LEU
4	D	2	GLN
4	D	33	TYR
4	D	38	LYS
4	D	46	ILE
4	D	50	ARG
4	D	81	LEU
4	D	83	LEU
4	D	84	GLU
4	D	100	LYS
4	D	105	GLN
4	D	126	ARG
4	D	154	ASP
4	D	171	LYS
5	E	19	MET
5	E	28	ASN
5	E	30	ASN
5	E	31	SER
5	E	54	THR
5	E	56	ASP
5	E	66	VAL
5	E	69	LEU
5	E	72	ARG
5	E	78	LEU
5	E	94	ARG
5	E	96	ARG
5	E	102	GLU
5	E	110	ARG
5	E	132	GLU
5	E	146	LEU
5	E	148	THR
5	E	180	GLN

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Mol	Chain	Res	Type
5	E	184	ASN
5	E	222	GLU
5	E	226	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	70	ASN
1	A	80	ASN
1	A	93	HIS
1	A	180	GLN
1	A	263	HIS
2	B	83	ASN
2	B	89	GLN
4	D	51	GLN
4	D	116	GLN
4	D	173	ASN
5	E	11	GLN
5	E	28	ASN
5	E	29	HIS
5	E	30	ASN
5	E	62	ASN
5	E	118	ASN
5	E	213	GLN
5	E	220	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 carbohydrates in this entry.

5.6 Ligand geometry

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
6	GOL	A	301	-	5,5,5	0.32	0	5,5,5	0.25	0
6	GOL	B	1100	-	5,5,5	0.32	0	5,5,5	0.33	0
6	GOL	B	1101	-	5,5,5	0.34	0	5,5,5	0.23	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
6	GOL	A	301	-	-	0/4/4/4	0/0/0/0
6	GOL	B	1100	-	-	0/4/4/4	0/0/0/0
6	GOL	B	1101	-	-	0/4/4/4	0/0/0/0

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.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	301	GOL	1	0
6	B	1100	GOL	1	0
6	B	1101	GOL	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 ⓘ

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	276/276 (100%)	0.19	10 (3%) 43 36	36, 52, 82, 119	0
2	B	99/99 (100%)	0.45	6 (6%) 22 16	41, 62, 93, 123	0
3	C	9/9 (100%)	0.61	0 100 100	38, 46, 59, 63	0
4	D	203/204 (99%)	0.49	18 (8%) 10 6	36, 55, 109, 150	0
5	E	242/242 (100%)	0.12	2 (0%) 86 83	31, 51, 86, 126	0
All	All	829/830 (99%)	0.28	36 (4%) 36 29	31, 53, 95, 150	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	D	204	PRO	6.6
4	D	131	SER	5.4
1	A	105	PRO	4.3
4	D	182	SER	4.2
2	B	99	MET	3.9
4	D	132	ASP	3.6
4	D	181	LYS	3.5
4	D	152	ASP	3.5
4	D	183	ASP	3.5
4	D	151	LYS	3.4
4	D	98	ALA	3.3
4	D	203	SER	3.3
5	E	220	ASN	3.2
1	A	104	GLY	3.1
1	A	1	GLY	3.0
4	D	187	ALA	2.8
1	A	106	ASP	2.8
4	D	191	ASN	2.7
5	E	223	TRP	2.7
4	D	194	ILE	2.5

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Mol	Chain	Res	Type	RSRZ
2	B	46	ILE	2.4
4	D	129	LYS	2.4
1	A	276	PRO	2.4
2	B	97	ARG	2.4
1	A	16	GLY	2.4
4	D	185	ALA	2.4
1	A	194	VAL	2.3
4	D	56	GLN	2.3
2	B	75	LYS	2.2
2	B	1	ILE	2.2
4	D	184	PHE	2.2
4	D	192	ASN	2.2
2	B	19	LYS	2.1
1	A	197	HIS	2.0
1	A	195	SER	2.0
1	A	196	ASP	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
6	GOL	A	301	6/6	0.87	0.26	6.04	65,69,70,85	0
6	GOL	B	1101	6/6	0.93	0.20	0.02	64,67,73,75	0
6	GOL	B	1100	6/6	0.96	0.15	-1.02	59,64,72,72	0

6.5 Other polymers

There are no such residues in this entry.