



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 7, 2020 – 07:24 AM BST

PDB ID : 6VTH
Title : p53-specific T cell receptor
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Deposited on : 2020-02-12
Resolution : 2.36 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

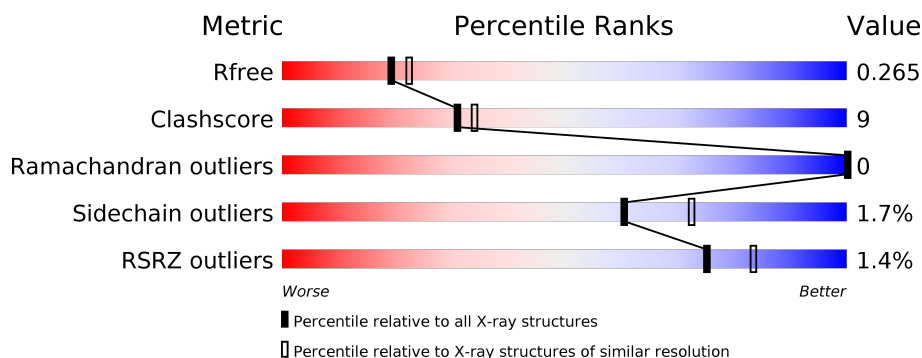
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.36 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}	130704	1164 (2.36-2.36)
Clashscore	141614	1232 (2.36-2.36)
Ramachandran outliers	138981	1211 (2.36-2.36)
Sidechain outliers	138945	1212 (2.36-2.36)
RSRZ outliers	127900	1150 (2.36-2.36)

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 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	<div> <div>74%</div> <div>19%</div> <div>6%</div> </div>
1	D	205	<div> <div>73%</div> <div>21%</div> <div>6%</div> </div>
2	B	246	<div> <div>83%</div> <div>14%</div> <div>•</div> </div>
2	E	246	<div> <div>79%</div> <div>19%</div> <div>•</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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	SO4	B	301	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7046 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 T-cell Receptor 12-6, Alfa chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	192	Total	C	N	O	S	0	0	0
			1464	923	243	289	9			
1	D	193	Total	C	N	O	S	0	0	0
			1465	922	239	295	9			

- Molecule 2 is a protein called TCR 12-6, beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	240	Total	C	N	O	S	0	0	0
			1872	1181	324	359	8			
2	E	242	Total	C	N	O	S	0	0	0
			1893	1194	328	363	8			

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	S	0	0
			5	4	1		

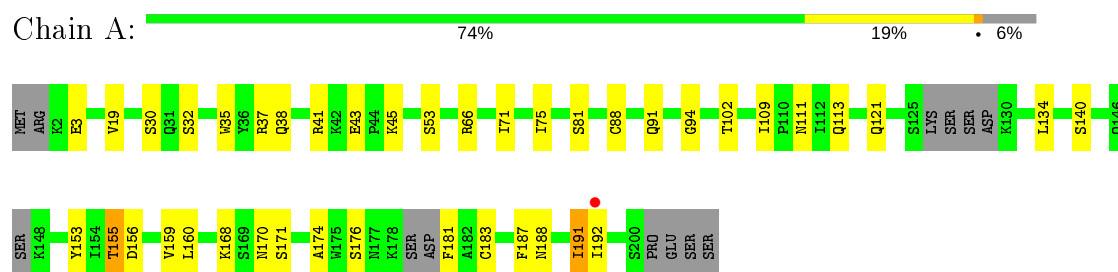
- Molecule 4 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	109	Total	O		0	0
			109	109			
4	B	96	Total	O		0	0
			96	96			
4	D	81	Total	O		0	0
			81	81			
4	E	61	Total	O		0	0
			61	61			

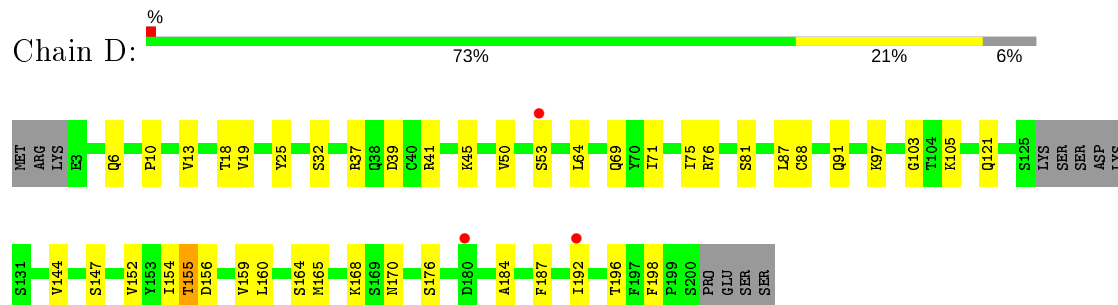
3 Residue-property plots [i](#)

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.

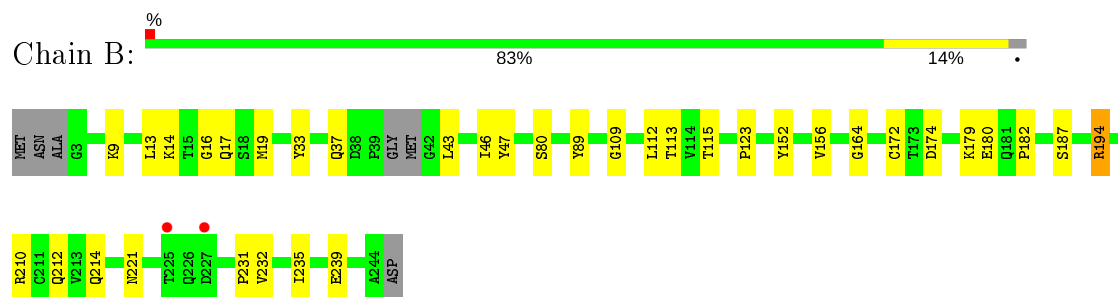
- Molecule 1: T-cell Receptor 12-6, Alfa chain



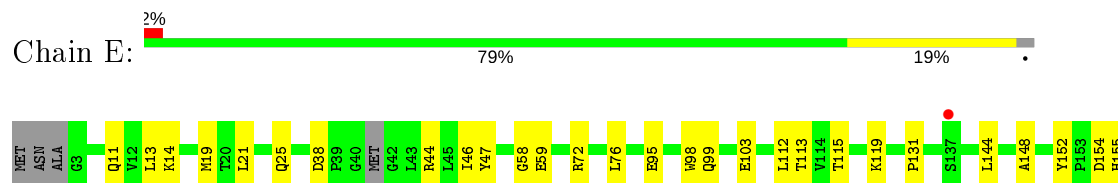
- Molecule 1: T-cell Receptor 12-6, Alfa chain



- Molecule 2: TCR 12-6, beta chain



- Molecule 2: TCR 12-6, beta chain





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	65.61Å 84.95Å 88.75Å 90.00° 93.61° 90.00°	Depositor
Resolution (Å)	20.00 – 2.36 42.48 – 2.36	Depositor EDS
% Data completeness (in resolution range)	91.0 (20.00-2.36) 91.1 (42.48-2.36)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.93 (at 2.37Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.211 , 0.264 0.211 , 0.265	Depositor DCC
R_{free} test set	1787 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	35.2	Xtriage
Anisotropy	0.713	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 45.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7046	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 40.73 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.6191e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: 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	0.40	0/1495	0.76	0/2031
1	D	0.39	0/1498	0.72	0/2040
2	B	0.40	0/1923	0.75	0/2624
2	E	0.37	0/1944	0.72	1/2649 (0.0%)
All	All	0.39	0/6860	0.74	1/9344 (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	A	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	E	194	ARG	CG-CD-NE	5.28	122.90	111.80

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	3	GLU	Peptide

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	1464	0	1359	31	0
1	D	1465	0	1349	32	0
2	B	1872	0	1743	25	0
2	E	1893	0	1774	42	0
3	B	5	0	0	2	0
4	A	109	0	0	9	0
4	B	96	0	0	3	0
4	D	81	0	0	8	0
4	E	61	0	0	7	0
All	All	7046	0	6225	118	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 9.

All (118) 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:155:THR:HG21	2:B:194:ARG:HH12	1.37	0.88
1:A:168:LYS:NZ	4:A:303:HOH:O	2.12	0.80
2:E:210:ARG:NH1	2:E:212:GLN:OE1	2.18	0.77
2:E:59:GLU:OE1	4:E:301:HOH:O	2.02	0.77
2:E:164:GLY:O	4:E:302:HOH:O	2.04	0.76
2:E:219:SER:HA	4:E:313:HOH:O	1.87	0.75
1:D:76:ARG:NH1	4:D:301:HOH:O	2.20	0.74
1:D:45:LYS:NZ	4:D:304:HOH:O	2.20	0.73
1:A:191:ILE:O	1:A:191:ILE:HG13	1.90	0.72
2:E:239:GLU:OE1	4:E:304:HOH:O	2.09	0.70
2:E:103:GLU:OE1	4:E:303:HOH:O	2.09	0.70
2:B:180:GLU:OE1	4:B:401:HOH:O	2.08	0.70
2:B:174:ASP:OD1	4:B:403:HOH:O	2.10	0.69
1:A:155:THR:HG21	2:B:194:ARG:NH1	2.07	0.69
1:D:18:THR:OG1	4:D:301:HOH:O	2.12	0.68
1:A:102:THR:O	4:A:302:HOH:O	2.12	0.67
1:D:196:THR:O	4:D:302:HOH:O	2.13	0.67
1:D:41:ARG:NH2	2:E:154:ASP:O	2.28	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:94:GLY:O	4:A:301:HOH:O	2.12	0.66
1:A:155:THR:CG2	2:B:194:ARG:HH12	2.08	0.66
1:A:41:ARG:NH1	4:A:306:HOH:O	2.28	0.65
1:A:37:ARG:NH2	1:A:81:SER:O	2.31	0.64
1:D:155:THR:HG21	2:E:194:ARG:HH12	1.63	0.64
1:D:50:VAL:HG12	1:D:64:LEU:HB2	1.78	0.63
2:E:119:LYS:HD3	2:E:226:GLN:NE2	2.14	0.62
2:B:210:ARG:NH1	2:B:212:GLN:OE1	2.34	0.61
2:E:224:TRP:HB2	2:E:230:LYS:HD2	1.84	0.59
1:D:121:GLN:O	4:D:303:HOH:O	2.17	0.59
1:D:160:LEU:HD23	4:D:313:HOH:O	2.03	0.58
2:E:115:THR:HG21	2:E:152:TYR:HE1	1.68	0.58
1:A:111:ASN:ND2	1:A:113:GLN:OE1	2.37	0.57
2:B:179:LYS:HE3	2:B:187:SER:HB3	1.87	0.57
1:A:188:ASN:N	4:A:307:HOH:O	2.38	0.56
1:A:111:ASN:O	4:A:304:HOH:O	2.18	0.56
1:D:13:VAL:HG21	1:D:19:VAL:HB	1.87	0.56
1:D:187:PHE:HB2	1:D:192:ILE:HD11	1.88	0.55
2:B:37:GLN:NE2	4:B:402:HOH:O	2.09	0.55
2:E:21:LEU:HD12	2:E:76:LEU:HD23	1.88	0.55
1:D:155:THR:HG22	1:D:156:ASP:O	2.06	0.55
2:B:210:ARG:HG3	2:B:239:GLU:HG2	1.88	0.53
2:B:14:LYS:HB3	2:B:17:GLN:NE2	2.25	0.52
2:B:13:LEU:HD11	2:B:19:MET:HG2	1.91	0.52
2:E:95:GLU:HG3	4:E:325:HOH:O	2.11	0.51
2:B:16:GLY:HA2	2:B:80:SER:HB2	1.93	0.50
2:E:25:GLN:O	2:E:72:ARG:HG3	2.10	0.50
1:A:153:TYR:O	1:A:174:ALA:HA	2.12	0.50
2:E:98:TRP:CD1	2:E:99:GLN:HG3	2.47	0.49
2:E:11:GLN:HG2	2:E:19:MET:SD	2.52	0.49
1:D:159:VAL:HG22	1:D:170:ASN:OD1	2.13	0.49
1:A:19:VAL:CG1	1:A:75:ILE:HB	2.44	0.48
2:B:46:ILE:HG22	2:B:47:TYR:HD1	1.77	0.48
2:E:174:ASP:HB2	2:E:191:LEU:HD12	1.96	0.48
2:E:14:LYS:HA	2:E:115:THR:O	2.14	0.48
2:E:13:LEU:HD11	2:E:19:MET:HG2	1.95	0.48
2:B:112:LEU:HD23	2:B:113:THR:N	2.28	0.47
1:A:43:GLU:OE2	1:A:45:LYS:NZ	2.46	0.47
2:B:123:PRO:HD3	2:B:231:PRO:HB3	1.96	0.47
1:D:19:VAL:HG12	1:D:75:ILE:HB	1.96	0.47
2:E:155:HIS:HD2	4:E:342:HOH:O	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:187:PHE:HB2	1:A:192:ILE:HD11	1.96	0.47
1:D:168:LYS:NZ	4:D:311:HOH:O	2.37	0.46
2:E:194:ARG:N	2:E:194:ARG:HD2	2.30	0.46
2:E:38:ASP:OD2	2:E:44:ARG:NH2	2.48	0.46
1:D:19:VAL:CG1	1:D:75:ILE:HB	2.45	0.46
1:A:38:GLN:NE2	4:A:316:HOH:O	2.48	0.46
1:A:160:LEU:HB3	2:B:172:CYS:HB2	1.98	0.46
1:A:45:LYS:HE2	4:A:401:HOH:O	2.16	0.45
2:E:115:THR:HG21	2:E:152:TYR:CE1	2.48	0.45
1:A:155:THR:HG22	1:A:156:ASP:O	2.15	0.45
1:D:184:ALA:HA	1:D:198:PHE:CE1	2.52	0.45
1:A:121:GLN:HB2	1:A:183:CYS:SG	2.56	0.45
1:A:159:VAL:HG22	1:A:170:ASN:OD1	2.16	0.45
1:A:32:SER:HB2	1:A:91:GLN:HB3	1.99	0.45
2:E:119:LYS:O	2:E:228:ARG:NH2	2.50	0.45
1:D:25:TYR:CZ	1:D:69:GLN:HA	2.51	0.45
1:D:10:PRO:HB3	1:D:105:LYS:HB3	1.98	0.45
2:B:179:LYS:HD3	2:B:182:PRO:HA	1.99	0.45
1:D:87:LEU:HD23	1:D:103:GLY:HA3	1.99	0.45
1:D:37:ARG:NH2	1:D:81:SER:O	2.50	0.45
1:D:71:ILE:HD11	1:D:88:CYS:SG	2.57	0.44
1:D:97:LYS:HE3	2:E:58:GLY:HA2	1.98	0.44
1:D:155:THR:CG2	2:E:194:ARG:HH12	2.29	0.44
2:B:115:THR:HG23	2:B:152:TYR:OH	2.18	0.44
1:D:32:SER:HB2	1:D:91:GLN:HB3	2.00	0.44
2:E:178:LEU:C	2:E:178:LEU:HD12	2.38	0.44
1:D:147:SER:HB2	1:D:154:ILE:HG13	2.00	0.44
1:A:134:LEU:HD11	1:A:171:SER:HB2	1.99	0.44
2:E:131:PRO:HD3	2:E:144:LEU:HG	2.01	0.43
2:B:235:ILE:HD12	2:E:241:TRP:CE2	2.53	0.43
1:D:152:VAL:HG22	1:D:176:SER:HB2	2.00	0.43
1:D:165:MET:CE	2:E:198:SER:HB3	2.48	0.43
2:E:218:LEU:HD12	2:E:231:PRO:HD2	2.00	0.43
2:E:13:LEU:HD11	2:E:19:MET:CG	2.49	0.43
1:D:6:GLN:HB2	4:D:323:HOH:O	2.18	0.42
2:B:13:LEU:HD11	2:B:19:MET:CG	2.49	0.42
1:D:155:THR:HG21	2:E:194:ARG:NH1	2.33	0.42
2:E:154:ASP:HB3	2:E:189:TYR:CD1	2.55	0.42
2:B:235:ILE:HD12	2:E:241:TRP:CD2	2.55	0.42
1:A:109:ILE:CG2	1:A:140:SER:HB3	2.50	0.42
1:A:35:TRP:CZ2	1:A:71:ILE:HG13	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:71:ILE:HD11	1:A:88:CYS:SG	2.60	0.42
2:E:174:ASP:HB2	2:E:191:LEU:CD1	2.50	0.42
2:E:171:VAL:HA	2:E:194:ARG:O	2.20	0.41
2:B:37:GLN:HB2	2:B:43:LEU:HD23	2.02	0.41
1:A:30:SER:O	1:A:66:ARG:NH1	2.42	0.41
2:B:164:GLY:HA2	3:B:301:SO4:O2	2.19	0.41
1:D:37:ARG:NH1	1:D:39:ASP:OD1	2.53	0.41
2:B:89:TYR:O	2:B:109:GLY:HA2	2.21	0.41
3:B:301:SO4:O4	2:E:210:ARG:NH2	2.53	0.41
2:E:112:LEU:HD23	2:E:113:THR:N	2.36	0.41
2:E:115:THR:CG2	2:E:152:TYR:HE1	2.33	0.41
1:D:37:ARG:CZ	1:D:39:ASP:OD1	2.69	0.41
2:E:46:ILE:HG22	2:E:47:TYR:HD1	1.86	0.41
1:A:160:LEU:C	1:A:160:LEU:HD12	2.42	0.40
1:A:176:SER:HB3	1:A:181:PHE:CG	2.56	0.40
1:A:45:LYS:CE	4:A:401:HOH:O	2.70	0.40
2:B:156:VAL:HA	2:B:214:GLN:O	2.21	0.40
2:E:148:ALA:O	2:E:190:ALA:HA	2.22	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	184/205 (90%)	176 (96%)	8 (4%)	0	100	100
1	D	189/205 (92%)	186 (98%)	3 (2%)	0	100	100
2	B	236/246 (96%)	231 (98%)	5 (2%)	0	100	100
2	E	238/246 (97%)	233 (98%)	5 (2%)	0	100	100
All	All	847/902 (94%)	826 (98%)	21 (2%)	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	161/185 (87%)	158 (98%)	3 (2%)	57	68
1	D	162/185 (88%)	158 (98%)	4 (2%)	47	58
2	B	199/213 (93%)	194 (98%)	5 (2%)	47	58
2	E	202/213 (95%)	202 (100%)	0	100	100
All	All	724/796 (91%)	712 (98%)	12 (2%)	60	72

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	53	SER
1	A	155	THR
1	A	191	ILE
2	B	9	LYS
2	B	33	TYR
2	B	194	ARG
2	B	221	ASN
2	B	232	VAL
1	D	53	SER
1	D	144	VAL
1	D	155	THR
1	D	164	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	188	ASN
2	B	17	GLN
2	B	120	ASN
2	E	155	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand 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	SO4	B	301	-	4,4,4	0.33	0	6,6,6	0.35	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.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	301	SO4	2	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	192/205 (93%)	0.05	1 (0%) 91 95	24, 38, 67, 78	0
1	D	193/205 (94%)	0.16	3 (1%) 72 80	25, 41, 67, 84	0
2	B	240/246 (97%)	0.02	2 (0%) 86 91	23, 35, 55, 75	0
2	E	242/246 (98%)	0.06	6 (2%) 57 67	22, 38, 63, 84	0
All	All	867/902 (96%)	0.07	12 (1%) 75 83	22, 37, 66, 84	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	E	227	ASP	3.5
2	E	221	ASN	3.5
1	D	53	SER	3.0
2	E	225	THR	3.0
1	D	180	ASP	2.8
2	E	223	GLU	2.8
2	B	227	ASP	2.6
2	E	137	SER	2.5
2	E	222	ASP	2.4
1	A	192	ILE	2.3
2	B	225	THR	2.1
1	D	192	ILE	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 monosaccharides 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(\AA^2)	Q<0.9
3	SO4	B	301	5/5	0.99	0.10	33,34,36,38	0

6.5 Other polymers [i](#)

There are no such residues in this entry.