Full wwPDB X-ray Structure Validation Report (i)

Sep 5, 2023 – 05:19 PM EDT

PDB ID	:	4E42
Title	:	Structural basis for the recognition of mutant self by a tumor-specific, MHC class II-restricted T cell receptor G4
Authors	:	Deng, L.; Langley, R.J.; Wang, Q.; Topalian, S.L.; Mariuzza, R.A.
Deposited on	:	2012-03-11
Resolution	:	2.70 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 (i) 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 (i)) were used in the production of this report:

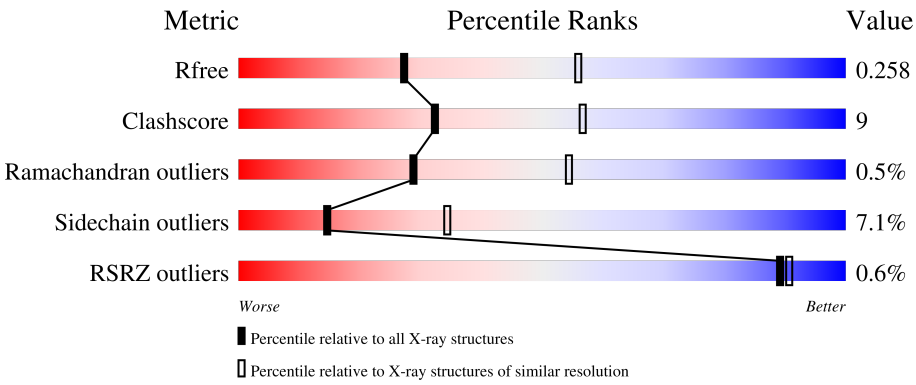
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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 <=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	203	71% 18% • 8%
1	C	203	% 72% 14% • • 10%
2	B	239	77% 17% • •
2	D	239	% 77% 15% • •

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
3	CL	B	302	-	-	X	-
4	NO3	B	301	-	-	X	-

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 6845 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 G4 alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	187	1480	932	246	295	7	0	0	0
1	C	182	1445	913	240	285	7	0	0	0

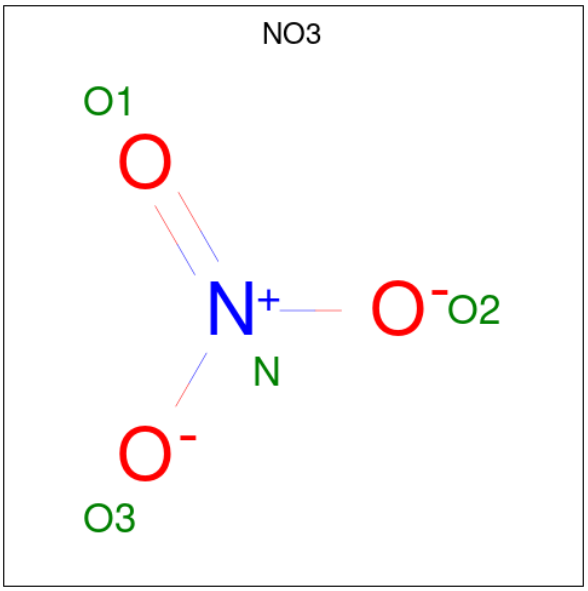
- Molecule 2 is a protein called T cell receptor G4 beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	231	1854	1175	324	350	5	0	0	0
2	D	229	1842	1167	322	348	5	0	0	0

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Cl		
3	A	1	1	1	0	0
3	B	4	4	4	0	0
3	D	4	4	4	0	0

- Molecule 4 is NITRATE ION (three-letter code: NO3) (formula: NO_3).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total N O 4 1 3	0	0
4	D	1	Total N O 4 1 3	0	0

- Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total Na 1 1	0	0
5	D	1	Total Na 1 1	0	0

- Molecule 6 is water.

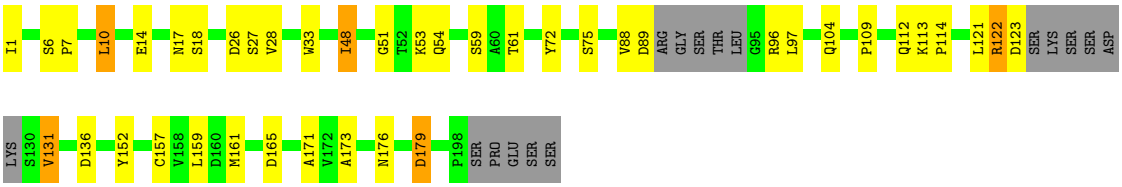
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	38	Total O 38 38	0	0
6	B	79	Total O 79 79	0	0
6	C	31	Total O 31 31	0	0
6	D	57	Total O 57 57	0	0

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.

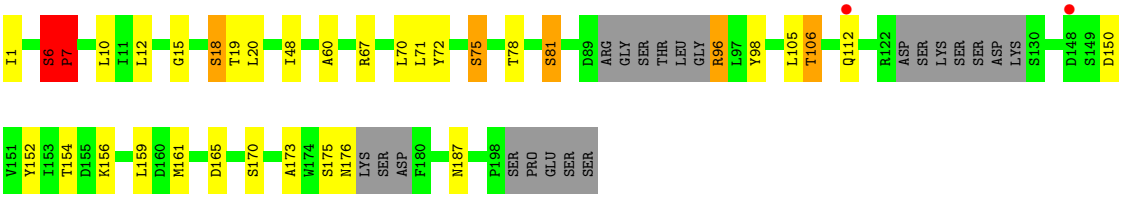
• Molecule 1: T cell receptor G4 alpha chain

Chain A:



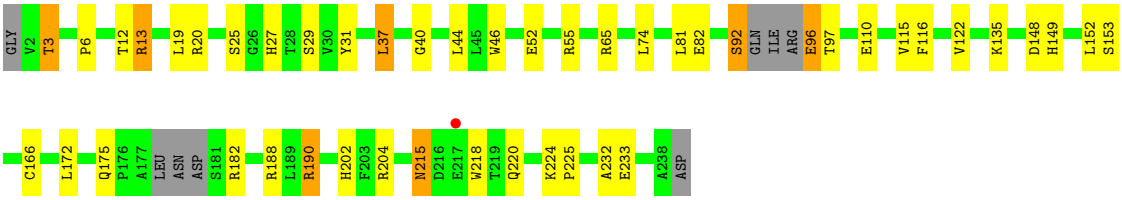
• Molecule 1: T cell receptor G4 alpha chain

Chain C:



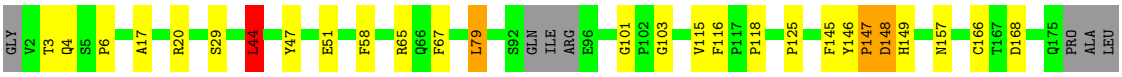
• Molecule 2: T cell receptor G4 beta chain

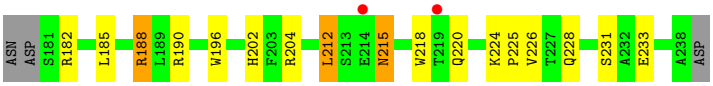
Chain B:



• Molecule 2: T cell receptor G4 beta chain

Chain D:




ASN
ASP
S181
R182
L185
R188
L189
R190
W196
H202
F203
R204
L212
S213
E214
N215
W218
T219
Q220
K224
P225
V226
T227
Q228
S231
A232
E233
A238
ASP

4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	168.07Å 75.36Å 130.86Å 90.00° 123.48° 90.00°	Depositor
Resolution (Å)	44.63 – 2.70 44.65 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.9 (44.63-2.70) 99.6 (44.65-2.30)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.92 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.200 , 0.263 0.199 , 0.258	Depositor DCC
R_{free} test set	3039 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å^2)	42.9	Xtriage
Anisotropy	0.205	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 42.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	6845	wwPDB-VP
Average B, all atoms (Å^2)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.43% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: CL, NA, NO3

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	0.61	1/1514 (0.1%)	0.68	0/2060
1	C	0.58	1/1478 (0.1%)	0.66	1/2011 (0.0%)
2	B	0.67	1/1906 (0.1%)	0.68	0/2596
2	D	0.59	1/1893 (0.1%)	0.65	3/2577 (0.1%)
All	All	0.61	4/6791 (0.1%)	0.67	4/9244 (0.0%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	7	PRO	N-CD	5.59	1.55	1.47
1	A	7	PRO	N-CD	5.56	1.55	1.47
2	D	6	PRO	N-CD	5.46	1.55	1.47
2	B	6	PRO	N-CD	5.28	1.55	1.47

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	D	44	LEU	CA-CB-CG	5.94	128.97	115.30
2	D	146	TYR	C-N-CD	5.40	139.74	128.40
2	D	147	PRO	CA-N-CD	-5.33	104.04	111.50
1	C	6	SER	C-N-CD	5.04	138.99	128.40

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	1480	0	1394	33	0
1	C	1445	0	1364	22	0
2	B	1854	0	1759	45	0
2	D	1842	0	1747	31	0
3	A	1	0	0	0	0
3	B	4	0	0	2	0
3	D	4	0	0	1	0
4	B	4	0	0	2	0
4	D	4	0	0	1	0
5	B	1	0	0	0	0
5	D	1	0	0	0	0
6	A	38	0	0	2	0
6	B	79	0	0	5	0
6	C	31	0	0	0	0
6	D	57	0	0	0	0
All	All	6845	0	6264	114	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 (114) 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:D:148:ASP:CG	2:D:149:HIS:H	1.41	1.24
2:B:148:ASP:CG	2:B:149:HIS:H	1.46	1.16
1:C:96:ARG:HH11	1:C:96:ARG:HG2	1.22	1.03
2:D:148:ASP:CG	2:D:149:HIS:N	2.13	0.97
2:B:13:ARG:HG3	2:B:13:ARG:HH11	1.28	0.96
2:B:148:ASP:CG	2:B:149:HIS:N	2.21	0.92
2:B:82:GLU:HG2	6:B:423:HOH:O	1.73	0.88
1:C:161:MET:HE2	2:D:190:ARG:HD3	1.56	0.87
2:D:3:THR:HG22	3:D:302:CL:CL	2.11	0.87
1:A:122:ARG:CG	1:A:122:ARG:HH11	1.90	0.85
2:B:116:PHE:CD1	2:B:182:ARG:HD3	2.13	0.84
1:C:161:MET:CE	2:D:190:ARG:HD3	2.07	0.83
2:B:202:HIS:HE1	2:B:233:GLU:OE1	1.59	0.83
1:A:122:ARG:HH11	1:A:122:ARG:HG3	1.46	0.80
1:C:96:ARG:HG2	1:C:96:ARG:NH1	2.01	0.73
1:C:170:SER:OG	2:D:188:ARG:HD3	1.88	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:161:MET:CE	2:B:190:ARG:HB3	2.20	0.72
2:B:13:ARG:HH11	2:B:13:ARG:CG	2.02	0.70
2:B:13:ARG:HG3	2:B:13:ARG:NH1	2.03	0.69
2:D:148:ASP:OD1	2:D:149:HIS:N	2.19	0.68
2:D:116:PHE:CD1	2:D:182:ARG:HD3	2.29	0.68
2:D:168:ASP:OD1	2:D:188:ARG:NH2	2.28	0.66
2:B:3:THR:HG22	3:B:302:CL:CL	2.34	0.65
1:A:161:MET:HE1	2:B:190:ARG:HD2	1.79	0.64
1:A:171:ALA:HA	2:B:188:ARG:HH21	1.62	0.64
1:C:96:ARG:HH11	1:C:96:ARG:CG	2.04	0.63
1:A:171:ALA:HA	2:B:188:ARG:NH2	2.13	0.63
2:D:215:ASN:ND2	2:D:215:ASN:H	1.97	0.63
1:C:19:THR:HG22	1:C:72:TYR:CD2	2.33	0.62
1:C:81:SER:HB3	1:C:106:THR:HA	1.81	0.61
2:D:157:ASN:ND2	2:D:202:HIS:H	2.01	0.59
1:A:161:MET:HE1	2:B:190:ARG:CD	2.32	0.59
2:B:202:HIS:HD2	4:B:301:NO3:O1	1.86	0.59
2:D:65:ARG:HD2	2:D:67:PHE:CE1	2.38	0.58
2:D:226:VAL:O	2:D:228:GLN:HG2	2.03	0.58
1:A:179:ASP:OD1	1:A:179:ASP:N	2.29	0.57
1:C:12:LEU:HD11	1:C:18:SER:HB2	1.85	0.57
1:A:161:MET:HE2	2:B:190:ARG:HB3	1.87	0.56
1:A:122:ARG:CG	1:A:122:ARG:NH1	2.61	0.56
1:C:159:LEU:HB3	2:D:166:CYS:HB2	1.87	0.56
2:B:202:HIS:CE1	2:B:233:GLU:OE1	2.50	0.56
2:D:157:ASN:HD21	2:D:202:HIS:H	1.52	0.56
1:A:89:ASP:HB2	1:A:97:LEU:HD23	1.88	0.55
2:D:218:TRP:HB2	2:D:224:LYS:HG3	1.88	0.55
1:A:112:GLN:O	1:A:114:PRO:HD2	2.05	0.55
2:B:218:TRP:HB2	2:B:224:LYS:HG3	1.88	0.55
1:A:161:MET:HE1	2:B:190:ARG:HB3	1.89	0.55
1:C:10:LEU:HD23	1:C:105:LEU:HD13	1.88	0.55
2:D:212:LEU:HD22	2:D:225:PRO:HG2	1.88	0.54
2:B:19:LEU:HD22	2:B:74:LEU:HD23	1.89	0.54
1:A:121:LEU:HB2	1:A:131:VAL:HG12	1.88	0.54
2:B:52:GLU:OE2	2:B:55:ARG:HG3	2.08	0.53
1:C:161:MET:HE1	2:D:190:ARG:HD3	1.87	0.53
1:A:1:ILE:HD12	1:A:26:ASP:HB3	1.91	0.52
1:A:54:GLN:HG3	1:A:59:SER:OG	2.10	0.52
2:B:116:PHE:CE1	2:B:182:ARG:HD3	2.44	0.52
2:B:13:ARG:CZ	2:B:110:GLU:HG2	2.40	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:152:LEU:C	2:B:152:LEU:HD23	2.31	0.51
1:C:20:LEU:HD22	1:C:71:LEU:HD23	1.93	0.51
1:A:17:ASN:ND2	1:A:72:TYR:OH	2.41	0.51
1:A:96:ARG:HD2	6:A:421:HOH:O	2.10	0.51
2:B:37:LEU:HD12	2:B:37:LEU:H	1.76	0.50
2:D:125:PRO:HD2	2:D:196:TRP:CZ2	2.46	0.50
2:B:122:VAL:HG23	2:B:232:ALA:HB3	1.93	0.50
2:D:17:ALA:HB2	2:D:79:LEU:HD21	1.94	0.49
2:B:82:GLU:HB2	6:B:479:HOH:O	2.12	0.49
2:D:118:PRO:HB3	2:D:145:PHE:HB3	1.94	0.49
1:A:28:VAL:HG11	1:A:88:VAL:HB	1.94	0.48
1:A:161:MET:HE2	2:B:190:ARG:CB	2.43	0.48
2:B:27:HIS:CG	2:B:92:SER:HB2	2.47	0.48
2:B:148:ASP:OD1	2:B:149:HIS:N	2.35	0.48
1:C:48:ILE:HG23	1:C:48:ILE:O	2.13	0.48
1:C:152:TYR:O	1:C:173:ALA:HA	2.13	0.48
2:D:204:ARG:NH1	4:D:301:NO3:O3	2.47	0.48
1:A:159:LEU:HB3	2:B:166:CYS:HB2	1.95	0.47
2:D:115:VAL:HG12	2:D:225:PRO:HB2	1.95	0.47
2:D:218:TRP:CB	2:D:224:LYS:HG3	2.45	0.47
1:A:33:TRP:HD1	1:A:48:ILE:HD11	1.80	0.46
2:B:27:HIS:ND1	2:B:92:SER:HB2	2.31	0.46
1:A:161:MET:CE	2:B:190:ARG:NE	2.79	0.46
1:A:14:GLU:HG2	1:A:109:PRO:HA	1.98	0.45
2:D:47:TYR:HA	2:D:51:GLU:O	2.16	0.45
6:B:469:HOH:O	1:C:98:TYR:HE2	1.98	0.45
1:C:156:LYS:HA	1:C:170:SER:O	2.16	0.45
1:A:152:TYR:O	1:A:173:ALA:HA	2.17	0.45
2:B:116:PHE:CD1	2:B:182:ARG:CD	2.93	0.45
2:B:115:VAL:HG12	2:B:225:PRO:HB2	2.00	0.44
2:B:204:ARG:NH1	4:B:301:NO3:O2	2.50	0.44
2:D:202:HIS:HE1	2:D:233:GLU:OE1	2.00	0.44
1:A:122:ARG:O	1:A:123:ASP:HB3	2.18	0.44
1:A:51:GLY:O	1:A:61:THR:HA	2.18	0.44
2:D:168:ASP:HB2	2:D:185:LEU:HD12	1.98	0.44
1:A:122:ARG:HG3	1:A:122:ARG:NH1	2.23	0.44
2:B:40:GLY:HA2	6:B:431:HOH:O	2.17	0.43
2:D:4:GLN:NE2	2:D:103:GLY:H	2.15	0.43
1:A:136:ASP:OD1	2:B:190:ARG:NH2	2.50	0.43
1:A:122:ARG:HH11	1:A:122:ARG:HG2	1.77	0.43
2:B:135:LYS:CE	6:B:415:HOH:O	2.67	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:31:TYR:CE1	2:B:46:TRP:CD1	3.07	0.43
2:B:13:ARG:HD3	2:B:81:LEU:HG	2.02	0.42
1:A:48:ILE:HG23	1:A:53:LYS:HG3	2.01	0.42
2:B:96:GLU:HB3	2:B:97:THR:H	1.61	0.41
2:D:44:LEU:HD13	2:D:58:PHE:CG	2.55	0.41
1:A:10:LEU:HD11	1:A:18:SER:OG	2.21	0.41
1:C:6:SER:HA	1:C:7:PRO:HA	1.86	0.41
1:C:60:ALA:HA	1:C:70:LEU:O	2.20	0.41
1:C:161:MET:HE1	2:D:190:ARG:CD	2.49	0.41
1:A:157:CYS:HB2	6:A:426:HOH:O	2.21	0.41
2:B:152:LEU:HD23	2:B:153:SER:N	2.35	0.41
1:C:78:THR:O	1:C:81:SER:OG	2.39	0.41
2:B:3:THR:CG2	3:B:302:CL:CL	3.06	0.41
1:C:15:GLY:O	1:C:75:SER:HA	2.21	0.41
2:B:115:VAL:O	2:B:225:PRO:HG3	2.22	0.40
2:D:4:GLN:HE21	2:D:101:GLY:HA3	1.86	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	181/203 (89%)	176 (97%)	4 (2%)	1 (1%)	25	50
1	C	174/203 (86%)	167 (96%)	7 (4%)	0	100	100
2	B	225/239 (94%)	215 (96%)	8 (4%)	2 (1%)	17	40
2	D	223/239 (93%)	210 (94%)	12 (5%)	1 (0%)	34	60
All	All	803/884 (91%)	768 (96%)	31 (4%)	4 (0%)	29	54

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	220	GLN
2	B	215	ASN
1	A	113	LYS
2	D	147	PRO

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	169/184 (92%)	158 (94%)	11 (6%)	17 38
1	C	165/184 (90%)	149 (90%)	16 (10%)	8 19
2	B	201/208 (97%)	186 (92%)	15 (8%)	13 31
2	D	200/208 (96%)	190 (95%)	10 (5%)	24 51
All	All	735/784 (94%)	683 (93%)	52 (7%)	14 34

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	SER
1	A	10	LEU
1	A	27	SER
1	A	48	ILE
1	A	75	SER
1	A	104	GLN
1	A	122	ARG
1	A	131	VAL
1	A	165	ASP
1	A	176	ASN
1	A	179	ASP
2	B	3	THR
2	B	12	THR
2	B	13	ARG
2	B	20	ARG
2	B	25	SER
2	B	29	SER
2	B	37	LEU

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Mol	Chain	Res	Type
2	B	44	LEU
2	B	65	ARG
2	B	92	SER
2	B	96	GLU
2	B	172	LEU
2	B	175	GLN
2	B	190	ARG
2	B	215	ASN
1	C	1	ILE
1	C	6	SER
1	C	7	PRO
1	C	18	SER
1	C	67	ARG
1	C	75	SER
1	C	81	SER
1	C	96	ARG
1	C	106	THR
1	C	112	GLN
1	C	150	ASP
1	C	154	THR
1	C	165	ASP
1	C	175	SER
1	C	176	ASN
1	C	187	ASN
2	D	20	ARG
2	D	29	SER
2	D	44	LEU
2	D	79	LEU
2	D	148	ASP
2	D	188	ARG
2	D	212	LEU
2	D	215	ASN
2	D	220	GLN
2	D	231	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	188	ASN
2	B	75	ASN
2	B	132	HIS
2	B	202	HIS

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Mol	Chain	Res	Type
2	B	215	ASN
2	B	220	GLN
1	C	5	GLN
1	C	13	GLN
1	C	77	GLN
1	C	184	ASN
1	C	187	ASN
2	D	4	GLN
2	D	114	ASN
2	D	157	ASN
2	D	202	HIS
2	D	215	ASN
2	D	220	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)

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)

Of 13 ligands modelled in this entry, 11 are monoatomic - leaving 2 for Mogul analysis.

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	NO3	D	301	-	1,3,3	3.47	1 (100%)	0,3,3	-	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
4	NO3	B	301	-	1,3,3	3.61	1 (100%)	0,3,3	-	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	301	NO3	O1-N	3.61	1.40	1.24
4	D	301	NO3	O1-N	3.47	1.40	1.24

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	301	NO3	1	0
4	B	301	NO3	2	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 (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	187/203 (92%)	-0.36	0	100	100	24, 38, 60, 66	0
1	C	182/203 (89%)	-0.29	2 (1%)	80	82	30, 47, 63, 72	0
2	B	231/239 (96%)	-0.46	1 (0%)	92	93	23, 33, 55, 62	0
2	D	229/239 (95%)	-0.40	2 (0%)	84	85	29, 41, 60, 72	0
All	All	829/884 (93%)	-0.38	5 (0%)	89	91	23, 40, 60, 72	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	148	ASP	2.9
1	C	112	GLN	2.4
2	D	219	THR	2.2
2	D	214	GLU	2.1
2	B	217	GLU	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(Å^2)	Q<0.9
3	CL	D	302	1/1	0.88	0.15	55,55,55,55	0
5	NA	B	306	1/1	0.88	0.26	32,32,32,32	0
5	NA	D	306	1/1	0.90	0.24	33,33,33,33	0
3	CL	B	304	1/1	0.91	0.15	66,66,66,66	0
3	CL	B	302	1/1	0.91	0.22	44,44,44,44	0
3	CL	D	305	1/1	0.92	0.13	52,52,52,52	0
3	CL	D	304	1/1	0.95	0.13	54,54,54,54	0
3	CL	D	303	1/1	0.95	0.06	54,54,54,54	0
3	CL	A	301	1/1	0.96	0.08	55,55,55,55	0
3	CL	B	305	1/1	0.96	0.20	46,46,46,46	0
4	NO3	D	301	4/4	0.98	0.16	39,39,39,40	0
3	CL	B	303	1/1	0.98	0.10	54,54,54,54	0
4	NO3	B	301	4/4	0.98	0.14	32,32,32,33	0

6.5 Other polymers (i)

There are no such residues in this entry.