



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 25, 2017 – 11:22 PM EDT

PDB ID : 4D9R
Title : Inhibiting Alternative Pathway Complement Activation by Targeting the Exosite on Factor D
Authors : Murray, J.M.; Wiesmann, C.
Deposited on : unknown
Resolution : 2.42 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
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20029824
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-20029824

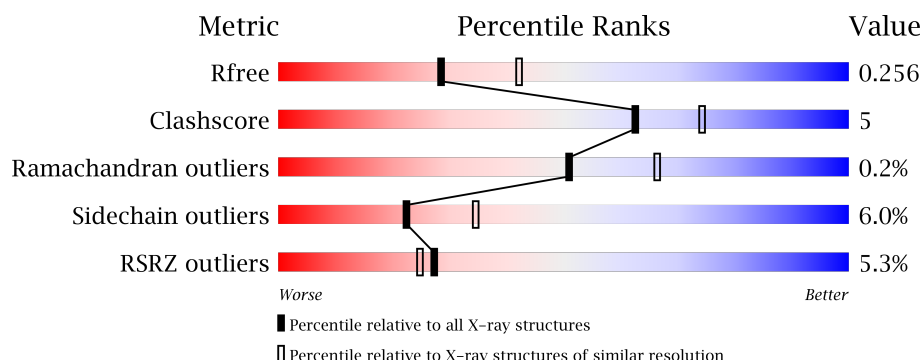
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.42 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	3709 (2.44-2.40)
Clashscore	112137	4241 (2.44-2.40)
Ramachandran outliers	110173	4178 (2.44-2.40)
Sidechain outliers	110143	4179 (2.44-2.40)
RSRZ outliers	101464	3740 (2.44-2.40)

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	228	<div> <div>2%</div> <div>89%</div> <div>10%</div> <div>.</div> </div>
1	B	228	<div> <div>21%</div> <div>86%</div> <div>13%</div> <div>.</div> </div>
2	D	214	<div> <div>%</div> <div>83%</div> <div>15%</div> <div>.</div> </div>
2	L	214	<div> <div>2%</div> <div>83%</div> <div>15%</div> <div>.</div> </div>
3	E	218	<div> <div>%</div> <div>87%</div> <div>11%</div> <div>..</div> </div>

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Mol	Chain	Length	Quality of chain
3	H	218	<div><div></div><div>3%</div><div>83%</div><div>15%</div><div></div></div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10296 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 Complement factor D.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	228	Total	C	N	O	S	0	3	0
			1721	1063	327	321	10			
1	B	228	Total	C	N	O	S	0	3	0
			1721	1063	327	321	10			

- Molecule 2 is a protein called Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	214	Total	C	N	O	S	0	0	0
			1630	1014	271	339	6			
2	D	213	Total	C	N	O	S	0	0	0
			1624	1011	270	338	5			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
L	105	GLU	-	LINKER	UNP P01834
L	106	ILE	-	LINKER	UNP P01834
L	107	LYS	-	LINKER	UNP P01834
L	108	ARG	-	LINKER	UNP P01834
D	105	GLU	-	LINKER	UNP P01834
D	106	ILE	-	LINKER	UNP P01834
D	107	LYS	-	LINKER	UNP P01834
D	108	ARG	-	LINKER	UNP P01834

- Molecule 3 is a protein called Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	218	Total	C	N	O	S	19	0	0
			1631	1030	267	327	7			
3	E	216	Total	C	N	O	S	0	0	0
			1618	1024	265	323	6			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
H	117	VAL	-	LINKER	UNP P01857
H	118	THR	-	LINKER	UNP P01857
H	119	VAL	-	LINKER	UNP P01857
H	120	SER	-	LINKER	UNP P01857
H	121	SER	-	LINKER	UNP P01857
E	117	VAL	-	LINKER	UNP P01857
E	118	THR	-	LINKER	UNP P01857
E	119	VAL	-	LINKER	UNP P01857
E	120	SER	-	LINKER	UNP P01857
E	121	SER	-	LINKER	UNP P01857

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	D	2	Total Cl 2 2	0	0

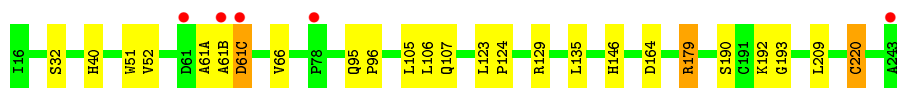
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	75	Total O 75 75	0	0
5	L	63	Total O 63 63	0	0
5	H	75	Total O 75 75	0	0
5	B	16	Total O 16 16	0	0
5	E	60	Total O 60 60	0	0
5	D	60	Total O 60 60	0	0

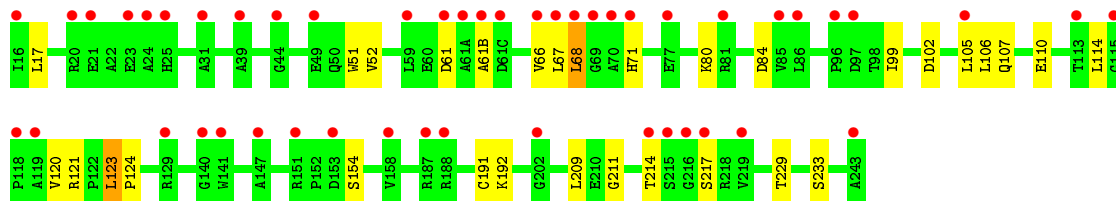
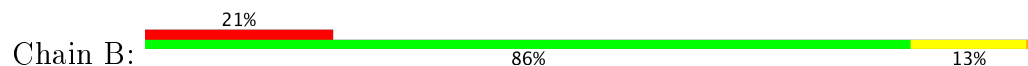
3 Residue-property plots [i](#)

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.

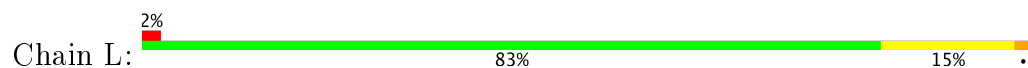
• Molecule 1: Complement factor D



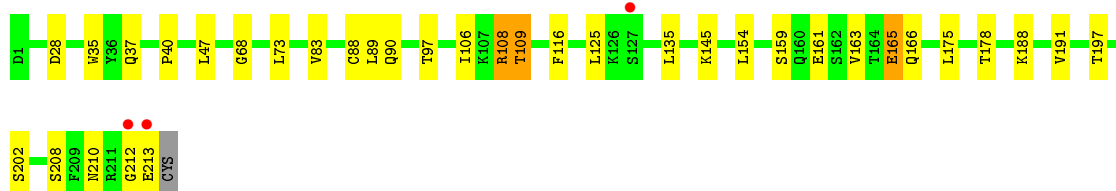
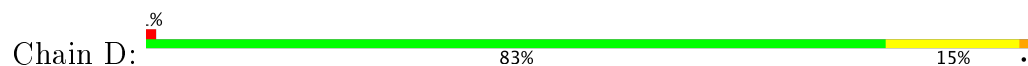
• Molecule 1: Complement factor D



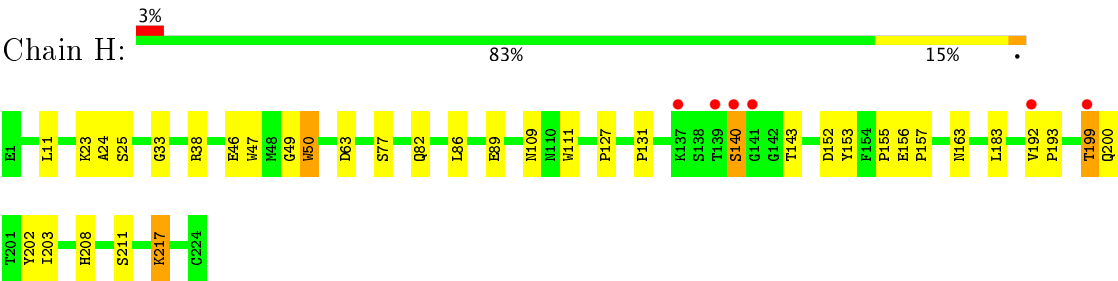
• Molecule 2: Fab light chain



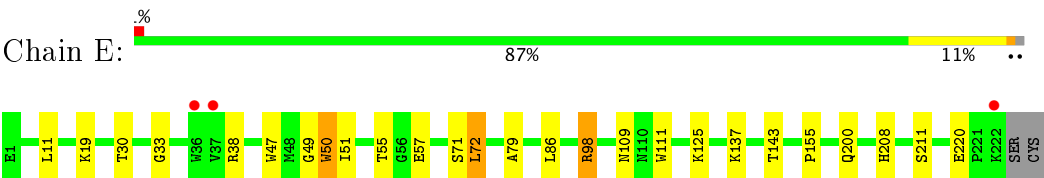
• Molecule 2: Fab light chain



● Molecule 3: Fab heavy chain



● Molecule 3: Fab heavy chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	132.05Å 132.05Å 180.29Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.84 – 2.42 19.84 – 2.42	Depositor EDS
% Data completeness (in resolution range)	98.7 (19.84-2.42) 98.7 (19.84-2.42)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.28 (at 2.41Å)	Xtriage
Refinement program	BUSTER-TNT BUSTER 2.9.6, BUSTER 2.11.2	Depositor
R, R_{free}	0.213 , 0.251 0.215 , 0.256	Depositor DCC
R_{free} test set	3064 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	46.2	Xtriage
Anisotropy	0.021	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 46.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10296	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.92% 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:
CL

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	0.51	0/1770	0.75	0/2408
1	B	0.49	0/1770	0.71	0/2408
2	D	0.51	0/1656	0.75	3/2252 (0.1%)
2	L	0.49	0/1662	0.76	0/2260
3	E	0.51	0/1658	0.77	0/2261
3	H	0.51	0/1671	0.75	0/2277
All	All	0.50	0/10187	0.75	3/13866 (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
2	L	0	1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	108	ARG	NE-CZ-NH1	6.12	123.36	120.30
2	D	108	ARG	NE-CZ-NH2	-6.09	117.25	120.30
2	D	212	GLY	N-CA-C	5.04	125.70	113.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	L	138	ASN	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	1721	0	1704	10	0
1	B	1721	0	1704	15	0
2	D	1624	0	1582	17	0
2	L	1630	0	1586	22	0
3	E	1618	0	1579	14	0
3	H	1631	0	1588	22	0
4	D	2	0	0	0	0
5	A	75	0	0	0	0
5	B	16	0	0	0	0
5	D	60	0	0	1	0
5	E	60	0	0	2	0
5	H	75	0	0	2	0
5	L	63	0	0	1	0
All	All	10296	0	9743	98	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 (98) 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:108:ARG:HH21	2:D:109:THR:HG22	1.07	1.16
3:H:156:GLU:HG3	3:H:157:PRO:HA	1.27	1.13
3:H:156:GLU:HG3	3:H:157:PRO:CA	1.98	0.94
2:D:108:ARG:NH2	2:D:109:THR:HG22	1.81	0.93
1:B:71:HIS:HE2	1:B:154:SER:HB2	1.39	0.85
1:B:71:HIS:NE2	1:B:154:SER:HB2	1.92	0.84
2:L:39:LYS:HB2	2:L:42:LYS:HG3	1.61	0.80
3:H:156:GLU:CG	3:H:157:PRO:HA	2.12	0.77
3:H:192:VAL:HG22	3:H:193:PRO:HD2	1.68	0.74
2:L:198:HIS:CD2	2:L:200:GLY:H	2.07	0.72
3:E:137:LYS:HE2	2:D:208:SER:O	1.94	0.68
3:E:109:ASN:HD21	3:E:111:TRP:HE1	1.41	0.66
2:L:29:ILE:HG13	2:L:92:ASP:HB2	1.77	0.66
2:D:161:GLU:HG3	5:D:405:HOH:O	1.95	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:71:HIS:CD2	1:B:154:SER:HB2	2.32	0.65
1:B:102:ASP:OD2	1:B:214:THR:HG23	1.95	0.65
3:H:109:ASN:HD21	3:H:111:TRP:HE1	1.43	0.64
1:A:52:VAL:HG21	1:A:66:VAL:HG11	1.80	0.64
2:L:198:HIS:HD2	2:L:200:GLY:H	1.46	0.63
1:A:40:HIS:HE1	1:A:193:GLY:O	1.83	0.62
3:E:55:THR:HG22	3:E:57:GLU:H	1.63	0.62
3:H:38:ARG:NH1	3:H:46:GLU:OE1	2.32	0.62
3:H:192:VAL:HG21	3:H:202:TYR:OH	2.01	0.61
1:A:123:LEU:HD12	1:A:124:PRO:HD2	1.85	0.58
1:B:102:ASP:HB3	1:B:214:THR:HG23	1.85	0.58
3:E:51:ILE:HD13	3:E:72:LEU:HD13	1.85	0.58
1:B:123:LEU:HD13	1:B:209:LEU:HB2	1.86	0.57
1:A:146:HIS:HE1	1:A:220:CYS:O	1.88	0.56
3:H:38:ARG:NH2	3:H:89:GLU:O	2.39	0.56
3:H:155:PRO:O	3:H:208:HIS:HE1	1.90	0.55
3:H:11:LEU:HB2	3:H:155:PRO:HG3	1.89	0.54
2:D:145:LYS:HB3	2:D:197:THR:HB	1.90	0.53
1:B:67:LEU:HD11	1:B:80:LYS:HG3	1.91	0.53
3:E:11:LEU:HB2	3:E:155:PRO:HG3	1.91	0.53
2:L:41:GLY:H	2:L:42:LYS:HE3	1.74	0.52
3:E:208:HIS:HD2	3:E:211:SER:OG	1.92	0.52
3:E:155:PRO:O	3:E:208:HIS:HE1	1.92	0.52
1:A:61(A):ALA:O	1:A:61(C):ASP:N	2.43	0.51
3:H:208:HIS:HD2	3:H:211:SER:OG	1.93	0.51
2:L:145:LYS:HB3	2:L:197:THR:HB	1.93	0.51
2:L:141:PRO:O	2:L:198:HIS:HE1	1.94	0.50
2:L:2:ILE:HD13	2:L:29:ILE:HD11	1.92	0.50
1:B:102:ASP:HB3	1:B:214:THR:CG2	2.42	0.50
3:E:72:LEU:HD12	3:E:79:ALA:HA	1.93	0.50
3:H:33:GLY:HA3	3:H:50:TRP:CZ3	2.47	0.49
3:E:33:GLY:HA3	3:E:50:TRP:CZ3	2.48	0.48
1:B:123:LEU:HD23	1:B:124:PRO:HD2	1.94	0.48
2:D:163:VAL:HG22	2:D:175:LEU:HD12	1.94	0.48
2:L:31:ASP:OD1	2:L:51:GLY:HA2	2.13	0.48
2:L:90:GLN:HE21	2:L:92:ASP:HB3	1.78	0.48
2:L:29:ILE:HG13	2:L:92:ASP:CB	2.43	0.48
2:D:108:ARG:HH21	2:D:109:THR:CG2	2.00	0.48
2:D:40:PRO:HG3	2:D:165:GLU:HG3	1.97	0.47
2:L:35:TRP:CZ3	2:L:88:CYS:HB3	2.49	0.47
1:A:32:SER:OG	1:A:40:HIS:HD2	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:140:SER:N	3:H:143:THR:O	2.38	0.46
1:B:51:TRP:CH2	1:B:107:GLN:HB2	2.51	0.45
2:D:35:TRP:CZ3	2:D:88:CYS:HB3	2.51	0.45
3:H:47:TRP:CZ3	3:H:49:GLY:HA2	2.52	0.45
2:L:159:SER:HA	2:L:178:THR:O	2.17	0.45
3:E:137:LYS:CE	2:D:208:SER:O	2.63	0.44
3:E:98:ARG:HD2	5:E:301:HOH:O	2.17	0.44
2:D:37:GLN:HB2	2:D:47:LEU:HD11	2.00	0.44
1:B:102:ASP:CB	1:B:214:THR:HG23	2.47	0.44
2:D:191:VAL:HG22	2:D:210:ASN:ND2	2.33	0.44
2:D:28:ASP:OD1	2:D:68:GLY:HA2	2.17	0.44
1:A:51:TRP:CH2	1:A:107:GLN:HB2	2.53	0.43
1:B:52:VAL:HG21	1:B:66:VAL:HG11	2.00	0.43
2:D:83:VAL:HG21	2:D:166:GLN:HB3	1.99	0.43
2:L:123:GLU:HB2	5:L:309:HOH:O	2.18	0.43
3:H:192:VAL:HG21	3:H:202:TYR:CZ	2.54	0.43
1:B:68:LEU:HD22	1:B:120:VAL:HG13	2.00	0.43
3:E:47:TRP:CZ3	3:E:49:GLY:HA2	2.53	0.43
2:L:39:LYS:CB	2:L:42:LYS:HG3	2.39	0.43
2:D:159:SER:HA	2:D:178:THR:O	2.19	0.43
3:H:152:ASP:HB3	3:H:183:LEU:HD13	2.01	0.43
2:L:191:VAL:HG22	2:L:210:ASN:ND2	2.34	0.42
1:B:71:HIS:CD2	1:B:154:SER:CB	3.02	0.42
1:B:211:GLY:HA2	1:B:229:THR:O	2.18	0.42
2:L:29:ILE:CG1	2:L:92:ASP:HB2	2.48	0.42
1:A:95:GLN:HB2	1:A:96:PRO:CD	2.50	0.42
1:A:123:LEU:CD2	1:A:209:LEU:HB2	2.49	0.42
3:E:51:ILE:HG21	5:E:306:HOH:O	2.18	0.42
3:H:127:PRO:HB3	3:H:153:TYR:HB3	2.02	0.42
3:H:163:ASN:ND2	3:H:203:ILE:H	2.18	0.42
3:H:199:THR:HB	5:H:317:HOH:O	2.19	0.41
2:L:37:GLN:HB2	2:L:47:LEU:HD11	2.02	0.41
2:L:27:THR:HG22	2:L:28:ASP:N	2.35	0.41
2:L:2:ILE:HG21	2:L:29:ILE:HD11	2.02	0.41
3:H:131:PRO:HD3	3:H:217:LYS:HE2	2.02	0.41
3:H:24:ALA:O	3:H:77:SER:OG	2.38	0.41
1:A:179:ARG:H	1:A:179:ARG:HG2	1.82	0.41
2:D:35:TRP:CD2	2:D:73:LEU:HB2	2.55	0.41
2:L:69:THR:HG22	2:L:70:ASP:OD1	2.21	0.40
2:D:116:PHE:HD2	2:D:135:LEU:HD23	1.86	0.40
3:E:71:SER:O	3:E:72:LEU:CD1	2.69	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:124:GLN:HE21	2:L:129:THR:HG23	1.86	0.40
3:H:23:LYS:HD3	5:H:301:HOH:O	2.20	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	229/228 (100%)	219 (96%)	9 (4%)	1 (0%)	38	52
1	B	229/228 (100%)	222 (97%)	6 (3%)	1 (0%)	38	52
2	D	211/214 (99%)	203 (96%)	8 (4%)	0	100	100
2	L	212/214 (99%)	205 (97%)	6 (3%)	1 (0%)	32	45
3	E	214/218 (98%)	210 (98%)	4 (2%)	0	100	100
3	H	216/218 (99%)	209 (97%)	7 (3%)	0	100	100
All	All	1311/1320 (99%)	1268 (97%)	40 (3%)	3 (0%)	51	66

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	61(B)	ALA
1	B	61(B)	ALA
2	L	40	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	185/182 (102%)	175 (95%)	10 (5%)	26	40
1	B	185/182 (102%)	169 (91%)	16 (9%)	12	18
2	D	190/191 (100%)	179 (94%)	11 (6%)	23	36
2	L	191/191 (100%)	180 (94%)	11 (6%)	23	36
3	E	182/184 (99%)	171 (94%)	11 (6%)	22	34
3	H	184/184 (100%)	175 (95%)	9 (5%)	29	45
All	All	1117/1114 (100%)	1049 (94%)	68 (6%)	22	33

All (68) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	61(C)	ASP
1	A	105	LEU
1	A	106	LEU
1	A	129	ARG
1	A	135	LEU
1	A	164	ASP
1	A	179	ARG
1	A	190	SER
1	A	192	LYS
1	A	220	CYS
2	L	29	ILE
2	L	31	ASP
2	L	42	LYS
2	L	69	THR
2	L	89	LEU
2	L	90	GLN
2	L	97	THR
2	L	106	ILE
2	L	107	LYS
2	L	125	LEU
2	L	214	CYS
3	H	25	SER
3	H	50	TRP
3	H	63	ASP
3	H	82	GLN
3	H	86	LEU
3	H	140	SER
3	H	199	THR

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Mol	Chain	Res	Type
3	H	200	GLN
3	H	217	LYS
1	B	17	LEU
1	B	61	ASP
1	B	68	LEU
1	B	84	ASP
1	B	99	ILE
1	B	105	LEU
1	B	106	LEU
1	B	110	GLU
1	B	114	LEU
1	B	121	ARG
1	B	123	LEU
1	B	191	CYS
1	B	192	LYS
1	B	217	SER
1	B	233[A]	SER
1	B	233[B]	SER
3	E	19	LYS
3	E	30	THR
3	E	38	ARG
3	E	50	TRP
3	E	72	LEU
3	E	86	LEU
3	E	98	ARG
3	E	125	LYS
3	E	143	THR
3	E	200	GLN
3	E	220	GLU
2	D	89	LEU
2	D	90	GLN
2	D	97	THR
2	D	106	ILE
2	D	109	THR
2	D	125	LEU
2	D	154	LEU
2	D	165	GLU
2	D	188	LYS
2	D	202	SER
2	D	213	GLU

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

Mol	Chain	Res	Type
1	A	40	HIS
1	A	50	GLN
1	A	146	HIS
1	A	157	HIS
2	L	90	GLN
2	L	147	GLN
2	L	198	HIS
2	L	210	ASN
3	H	109	ASN
3	H	163	ASN
3	H	179	GLN
3	H	208	HIS
1	B	157	HIS
3	E	82	GLN
3	E	109	ASN
3	E	179	GLN
3	E	208	HIS
2	D	37	GLN
2	D	210	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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

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 ⓘ

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	228/228 (100%)	-0.23	5 (2%) 62 59	29, 42, 69, 92	0
1	B	228/228 (100%)	0.94	48 (21%) 1 1	35, 80, 126, 164	0
2	D	213/214 (99%)	-0.23	3 (1%) 75 73	25, 46, 68, 119	0
2	L	214/214 (100%)	-0.27	4 (1%) 67 64	29, 43, 65, 84	0
3	E	216/218 (99%)	-0.17	3 (1%) 75 73	32, 46, 70, 99	0
3	H	215/218 (98%)	-0.25	6 (2%) 53 50	31, 41, 76, 109	0
All	All	1314/1320 (99%)	-0.03	69 (5%) 27 25	25, 46, 97, 164	0

All (69) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	H	139	THR	8.1
3	H	141	GLY	6.7
1	B	216	GLY	6.6
1	B	217	SER	6.5
2	D	212	GLY	6.2
1	B	153	ASP	5.3
3	H	140	SER	5.0
1	B	113	THR	4.2
1	B	69	GLY	4.2
3	E	222	LYS	4.1
1	B	61(A)	ALA	3.9
1	B	39	ALA	3.8
1	B	202	GLY	3.8
2	L	51	GLY	3.7
1	B	61(C)	ASP	3.5
1	A	61(C)	ASP	3.5
1	B	187	ARG	3.5
2	L	57	GLY	3.4
1	B	16	ILE	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	214	THR	3.1
1	A	61(B)	ALA	3.1
1	B	140	GLY	3.1
1	B	61(B)	ALA	3.0
1	B	49	GLU	3.0
2	D	213	GLU	2.9
2	L	88	CYS	2.9
1	B	66	VAL	2.8
1	B	25	HIS	2.8
1	A	243	ALA	2.8
1	B	31	ALA	2.8
2	D	127	SER	2.7
1	B	21	GLU	2.6
1	B	215	SER	2.6
1	B	20	ARG	2.6
1	B	115	GLY	2.5
1	B	86	LEU	2.5
1	B	219	VAL	2.5
1	B	71	HIS	2.5
1	B	77	GLU	2.4
3	H	137	LYS	2.4
1	B	67	LEU	2.4
3	E	37	VAL	2.3
1	B	243	ALA	2.3
1	B	23	GLU	2.3
1	B	129	ARG	2.3
2	L	41	GLY	2.3
1	B	97	ASP	2.3
1	B	59	LEU	2.3
1	B	158	VAL	2.2
3	H	199	THR	2.2
1	B	70	ALA	2.2
1	B	96	PRO	2.2
1	B	119	ALA	2.2
1	B	118	PRO	2.2
1	B	44	GLY	2.2
1	B	61	ASP	2.1
1	B	151	ARG	2.1
1	B	141	TRP	2.1
3	E	36	TRP	2.1
1	B	68	LEU	2.1
1	B	188	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	85	VAL	2.1
3	H	192	VAL	2.1
1	B	147	ALA	2.1
1	B	105	LEU	2.1
1	A	61	ASP	2.0
1	A	78	PRO	2.0
1	B	24	ALA	2.0
1	B	81	ARG	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
4	CL	D	302	1/1	0.97	0.16	0.27	63,63,63,63	0
4	CL	D	301	1/1	0.99	0.03	-3.79	35,35,35,35	0

6.5 Other polymers [i](#)

There are no such residues in this entry.