



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:20 AM UTC

PDB ID : 4O51 / pdb_00004o51
Title : Crystal structure of the QAA variant of anti-hinge rabbit antibody 2095-2 in complex with IDES hinge peptide
Authors : Malia, T.J.; Luo, J.; Teplyakov, A.; Gilliland, G.L.
Deposited on : 2013-12-19
Resolution : 2.20 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	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

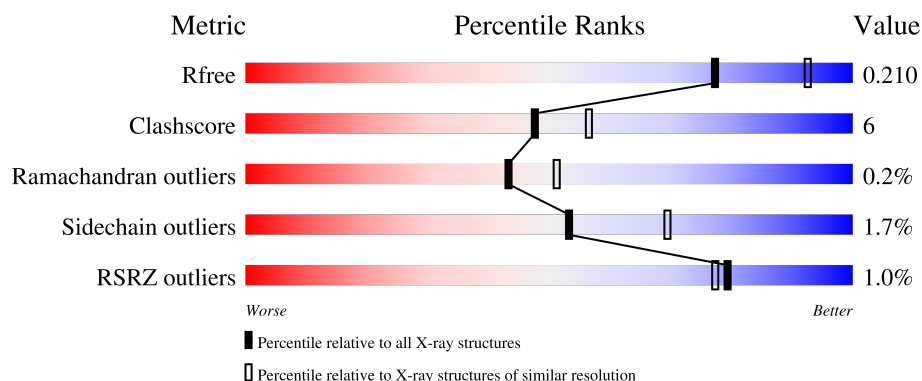
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	216	
1	C	216	
1	E	216	
1	L	216	
2	B	223	

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Mol	Chain	Length	Quality of chain
2	D	223	% 80% 15% • •
2	F	223	% 81% 15% •
2	H	223	2% 80% 15% •
3	M	14	36% 7% 57%
3	N	14	7% 29% 14% 57%
3	O	14	7% 21% 21% 57%
3	P	14	29% 14% 57%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 13803 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 QAA-2095-2 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	215	Total	C	N	O	S	0	0	0
			1624	1014	274	332	4			
1	A	215	Total	C	N	O	S	0	0	0
			1624	1014	274	332	4			
1	C	215	Total	C	N	O	S	0	0	0
			1624	1014	274	332	4			
1	E	215	Total	C	N	O	S	0	0	0
			1624	1014	274	332	4			

- Molecule 2 is a protein called QAA-2095-2 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	214	Total	C	N	O	S	0	0	0
			1578	1001	259	313	5			
2	B	214	Total	C	N	O	S	0	0	0
			1578	1001	259	313	5			
2	D	214	Total	C	N	O	S	0	0	0
			1578	1001	259	313	5			
2	F	214	Total	C	N	O	S	0	0	0
			1578	1001	259	313	5			

- Molecule 3 is a protein called IDES hinge peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	M	6	Total	C	N	O	0	0	0
			42	27	6	9			
3	N	6	Total	C	N	O	0	0	0
			42	27	6	9			
3	O	6	Total	C	N	O	0	0	0
			42	27	6	9			
3	P	6	Total	C	N	O	0	0	0
			42	27	6	9			

- Molecule 4 is water.

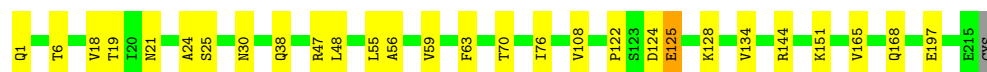
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	L	117	Total 117	O 117	0	0
4	H	116	Total 116	O 116	0	0
4	M	4	Total 4	O 4	0	0
4	A	82	Total 82	O 82	0	0
4	B	94	Total 94	O 94	0	0
4	N	1	Total 1	O 1	0	0
4	C	99	Total 99	O 99	0	0
4	D	97	Total 97	O 97	0	0
4	O	1	Total 1	O 1	0	0
4	E	108	Total 108	O 108	0	0
4	F	106	Total 106	O 106	0	0
4	P	2	Total 2	O 2	0	0

3 Residue-property plots

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.

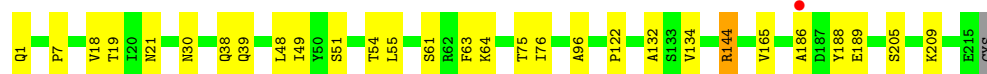
- Molecule 1: QAA-2095-2 light chain

Chain L:  87% 12%



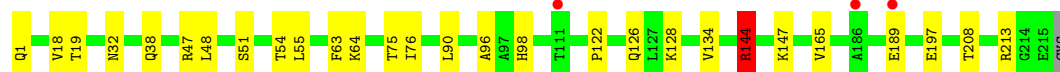
- Molecule 1: QAA-2095-2 light chain

Chain A:  86% 13%



- Molecule 1: QAA-2095-2 light chain

Chain C:  87% 12%



- Molecule 1: QAA-2095-2 light chain

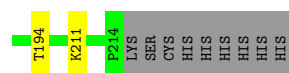
Chain E:  85% 14%



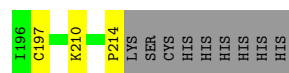
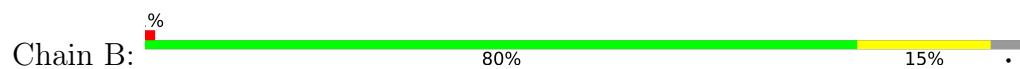
- Molecule 2: QAA-2095-2 heavy chain

Chain H:  80% 15%

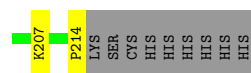
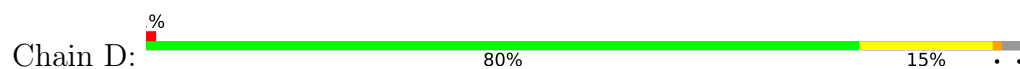




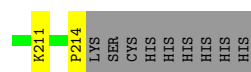
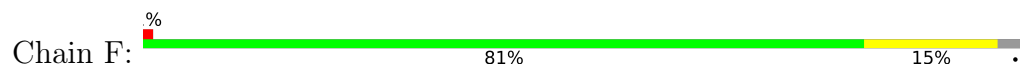
- Molecule 2: QAA-2095-2 heavy chain



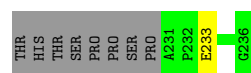
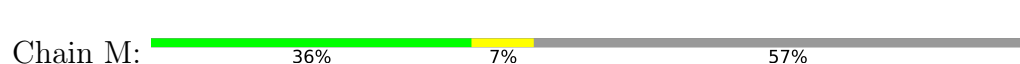
- Molecule 2: QAA-2095-2 heavy chain



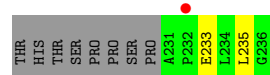
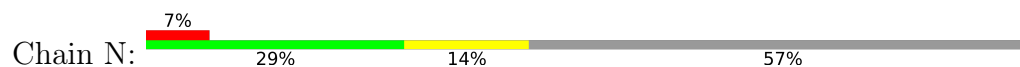
- Molecule 2: QAA-2095-2 heavy chain



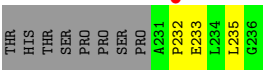
- Molecule 3: IDES hinge peptide



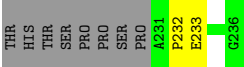
- Molecule 3: IDES hinge peptide



- Molecule 3: IDES hinge peptide



● Molecule 3: IDES hinge peptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	128.51Å 93.87Å 74.73Å 90.00° 90.04° 90.00°	Depositor
Resolution (Å)	48.74 – 2.20 48.74 – 2.20	Depositor EDS
% Data completeness (in resolution range)	89.0 (48.74-2.20) 88.9 (48.74-2.20)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.36 (at 2.20Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_1232)	Depositor
R, R_{free}	0.182 , 0.213 0.184 , 0.210	Depositor DCC
R_{free} test set	2000 reflections (2.33%)	wwPDB-VP
Wilson B-factor (Å ²)	39.0	Xtriage
Anisotropy	0.351	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 32.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.40$, $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	0.436 for -h,-k,l	Xtriage
Reported twinning fraction	0.450 for -h,-k,l	Depositor
Outliers	0 of 79799 reflections	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	13803	wwPDB-VP
Average B, all atoms (Å ²)	41.0	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 99.21 % 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 2.2972e-12. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: PCA

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	0.63	0/1650	0.78	2/2246 (0.1%)
1	C	0.62	1/1650 (0.1%)	0.79	1/2246 (0.0%)
1	E	0.63	1/1650 (0.1%)	0.79	1/2246 (0.0%)
1	L	0.60	0/1650	0.78	0/2246
2	B	0.60	0/1613	0.81	0/2212
2	D	0.57	0/1613	0.82	0/2212
2	F	0.57	0/1613	0.81	1/2212 (0.0%)
2	H	0.56	0/1613	0.82	1/2212 (0.0%)
3	M	0.48	0/42	0.92	0/55
3	N	0.66	0/42	0.75	0/55
3	O	0.82	0/42	0.81	0/55
3	P	0.47	0/42	0.90	0/55
All	All	0.60	2/13220 (0.0%)	0.80	6/18052 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	147	LYS	C-O	-5.28	1.18	1.24
1	E	171	LYS	C-O	-5.00	1.17	1.23

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	73	SER	N-CA-C	-10.49	100.22	113.23
2	F	73	SER	N-CA-C	-6.16	105.59	113.23
1	A	188	TYR	N-CA-C	-5.82	105.08	111.82
1	E	144	ARG	N-CA-C	5.49	117.26	111.28
1	A	144	ARG	N-CA-C	5.32	117.08	111.28
1	C	144	ARG	N-CA-C	5.26	117.02	111.28

There are no chirality outliers.

There are no planarity outliers.

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	1624	0	1572	15	0
1	C	1624	0	1572	17	0
1	E	1624	0	1572	20	0
1	L	1624	0	1572	16	0
2	B	1578	0	1552	19	0
2	D	1578	0	1552	24	0
2	F	1578	0	1552	21	0
2	H	1578	0	1552	18	0
3	M	42	0	42	1	0
3	N	42	0	42	2	0
3	O	42	0	42	4	0
3	P	42	0	42	2	0
4	A	82	0	0	0	0
4	B	94	0	0	1	0
4	C	99	0	0	2	0
4	D	97	0	0	2	0
4	E	108	0	0	2	0
4	F	106	0	0	2	0
4	H	116	0	0	1	0
4	L	117	0	0	1	0
4	M	4	0	0	0	0
4	N	1	0	0	0	0
4	O	1	0	0	0	0
4	P	2	0	0	0	0
All	All	13803	0	12664	143	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 6.

All (143) 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:H:72:SER:OG	2:H:75:THR:N	2.21	0.74
1:E:25:SER:HG	1:E:93:TYR:HH	1.35	0.74
1:L:151:LYS:NZ	1:L:197:GLU:OE1	2.25	0.69
2:B:50:ILE:HG12	2:B:56:ILE:HG12	1.74	0.69
1:E:151:LYS:NZ	1:E:197:GLU:OE1	2.25	0.69
2:F:118:LYS:NZ	4:F:353:HOH:O	2.25	0.69
2:F:3:VAL:HG12	2:F:23:VAL:HG12	1.74	0.68
2:H:3:VAL:HG12	2:H:23:VAL:HG12	1.78	0.66
1:C:144:ARG:HD3	1:C:165:VAL:HG11	1.77	0.66
2:D:207:LYS:NZ	4:D:309:HOH:O	2.28	0.66
2:H:50:ILE:HG12	2:H:56:ILE:HG12	1.79	0.65
1:A:21:ASN:ND2	1:E:201:GLN:OE1	2.29	0.65
2:B:77:ASP:OD2	4:B:382:HOH:O	2.15	0.64
1:C:128:LYS:HE3	4:C:337:HOH:O	1.95	0.64
2:D:50:ILE:HG12	2:D:56:ILE:HG12	1.79	0.64
1:A:122:PRO:HD3	1:A:134:VAL:HG22	1.81	0.62
1:L:25:SER:OG	2:D:4:GLU:OE2	2.14	0.62
1:E:78:ASP:OD1	4:E:307:HOH:O	2.16	0.61
1:C:122:PRO:HD3	1:C:134:VAL:HG22	1.83	0.61
2:B:127:PRO:HG2	2:B:214:PRO:HG3	1.84	0.60
2:F:132:THR:HG23	2:F:136:THR:H	1.67	0.60
1:A:51:SER:HB2	1:A:54:THR:HG22	1.83	0.59
2:F:193:GLN:NE2	4:F:367:HOH:O	2.35	0.59
2:F:54:GLY:HA2	2:F:70:ARG:HH12	1.67	0.58
1:E:108:VAL:O	1:E:168:GLN:NE2	2.36	0.57
2:B:95:ARG:HB3	2:B:103:VAL:HG22	1.87	0.57
2:F:132:THR:CG2	2:F:136:THR:H	2.16	0.57
2:D:95:ARG:HB3	2:D:103:VAL:HG22	1.86	0.56
1:E:19:THR:HG23	1:E:73:THR:CG2	2.35	0.56
2:H:169:ALA:HB2	2:H:179:LEU:HD23	1.87	0.56
2:F:50:ILE:HD11	2:F:68:ILE:HG22	1.88	0.56
2:F:132:THR:HG23	2:F:136:THR:C	2.31	0.56
2:D:127:PRO:HG2	2:D:214:PRO:HG3	1.88	0.55
1:L:125:GLU:HA	1:L:128:LYS:HE2	1.88	0.55
2:F:50:ILE:HG12	2:F:56:ILE:HG12	1.88	0.55
2:F:132:THR:HG23	2:F:136:THR:O	2.07	0.54
2:H:11:VAL:HG11	2:H:17:LEU:HD22	1.88	0.53
1:L:21:ASN:ND2	4:L:362:HOH:O	2.18	0.52
1:L:122:PRO:HD3	1:L:134:VAL:HG22	1.89	0.52
2:B:160:LEU:HD21	2:B:183:VAL:HG21	1.90	0.52
1:C:51:SER:HB2	1:C:54:THR:HG22	1.90	0.52
2:D:67:ILE:HG22	2:D:79:LYS:HB3	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:186:ALA:O	1:A:189:GLU:HG2	2.10	0.52
1:E:144:ARG:NE	4:E:316:HOH:O	2.30	0.52
2:H:50:ILE:HD11	2:H:68:ILE:HG22	1.91	0.51
1:A:96:ALA:HA	3:N:235:LEU:HB2	1.92	0.51
2:D:57:TRP:CH2	3:O:233:GLU:HG2	2.45	0.51
1:C:38:GLN:HB2	1:C:48:LEU:HD11	1.92	0.51
1:E:19:THR:HG23	1:E:73:THR:HG23	1.92	0.51
2:B:50:ILE:HD11	2:B:68:ILE:HG22	1.93	0.51
1:A:55:LEU:HD21	1:A:63:PHE:O	2.11	0.50
1:C:96:ALA:HA	3:O:235:LEU:HB2	1.93	0.50
2:D:160:LEU:HD21	2:D:183:VAL:HG21	1.92	0.50
2:B:128:SER:HB3	2:B:131:SER:HB2	1.93	0.50
1:C:126:GLN:NE2	4:C:396:HOH:O	2.45	0.50
2:D:50:ILE:HD11	2:D:68:ILE:HG22	1.94	0.50
1:C:144:ARG:O	1:C:144:ARG:HD2	2.12	0.50
1:E:122:PRO:HD3	1:E:134:VAL:HG22	1.93	0.50
1:L:47:ARG:NH2	2:H:100:ASP:HA	2.28	0.49
1:L:124:ASP:O	1:L:128:LYS:HG3	2.13	0.49
1:E:56:ALA:O	1:E:59:VAL:HG12	2.12	0.49
1:C:64:LYS:HB3	1:C:75:THR:HG23	1.93	0.49
2:H:10:LEU:HD21	2:H:115:ALA:O	2.13	0.49
2:D:120:PRO:HB3	2:D:146:TYR:HB3	1.95	0.49
2:F:9:ARG:CZ	2:F:11:VAL:HG12	2.43	0.49
1:L:55:LEU:HD21	1:L:63:PHE:O	2.13	0.48
1:L:108:VAL:O	1:L:168:GLN:NE2	2.45	0.48
2:F:10:LEU:HD21	2:F:115:ALA:O	2.12	0.48
2:D:133:SER:O	2:D:136:THR:HG22	2.14	0.48
1:L:56:ALA:O	1:L:59:VAL:HG12	2.14	0.48
2:B:57:TRP:CH2	3:N:233:GLU:HG2	2.48	0.48
1:A:18:VAL:HG22	1:A:76:ILE:HB	1.96	0.48
2:B:67:ILE:HG22	2:B:79:LYS:HB3	1.95	0.47
2:D:95:ARG:O	2:D:102:THR:HA	2.15	0.47
1:E:38:GLN:HB2	1:E:48:LEU:HD11	1.96	0.47
2:H:88:THR:HG23	2:H:111:THR:HA	1.96	0.47
1:C:47:ARG:NH2	2:D:100:ASP:HA	2.30	0.47
2:D:189:SER:HB2	2:D:193:GLN:CG	2.44	0.47
1:E:185:LYS:O	1:E:189:GLU:HG2	2.15	0.46
1:C:197:GLU:HG2	1:C:208:THR:OG1	2.16	0.46
1:C:55:LEU:HD21	1:C:63:PHE:O	2.16	0.46
2:B:120:PRO:HB3	2:B:146:TYR:HB3	1.97	0.46
2:F:169:ALA:HB2	2:F:179:LEU:HD23	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:58:TYR:HB2	2:D:63:LYS:HG2	1.98	0.45
1:L:144:ARG:CZ	1:L:165:VAL:HG21	2.46	0.45
2:D:88:THR:HG23	2:D:111:THR:HA	1.99	0.45
1:C:18:VAL:HG22	1:C:76:ILE:HB	1.98	0.45
1:L:24:ALA:O	1:L:70:THR:HG23	2.17	0.45
1:E:144:ARG:CZ	1:E:165:VAL:HG21	2.46	0.45
1:A:122:PRO:HG3	1:A:132:ALA:HB1	1.99	0.45
2:D:57:TRP:HZ2	3:O:232:PRO:HA	1.82	0.45
1:E:47:ARG:NH2	2:F:100:ASP:HA	2.31	0.45
2:H:132:THR:CG2	2:H:136:THR:H	2.30	0.44
2:F:127:PRO:HG2	2:F:214:PRO:HG3	2.00	0.44
2:H:37:ARG:HB3	2:H:47:ILE:HD11	1.99	0.44
2:B:197:CYS:SG	2:B:210:LYS:HB3	2.58	0.44
2:F:57:TRP:CH2	3:P:233:GLU:HG2	2.52	0.44
2:F:160:LEU:HD21	2:F:183:VAL:HG21	2.00	0.44
2:D:61:TRP:HE3	2:D:65:ARG:HH21	1.66	0.44
2:H:57:TRP:CH2	3:M:233:GLU:HG2	2.53	0.43
1:C:189:GLU:OE1	1:C:213:ARG:NH1	2.51	0.43
2:D:189:SER:HB2	2:D:193:GLN:HG2	1.99	0.43
1:E:48:LEU:O	1:E:59:VAL:HG11	2.18	0.43
2:F:28:LEU:HD21	2:F:76:VAL:HG23	1.99	0.43
2:F:194:THR:HG23	2:F:211:LYS:HE3	1.99	0.43
2:D:128:SER:HB3	2:D:131:SER:HB2	2.00	0.43
2:F:37:ARG:HB3	2:F:47:ILE:HD11	2.00	0.43
1:E:124:ASP:O	1:E:128:LYS:HG3	2.16	0.43
1:A:38:GLN:HB2	1:A:48:LEU:HD11	1.99	0.43
2:B:3:VAL:HG12	2:B:23:VAL:HG12	2.00	0.43
1:C:90:LEU:HD11	1:C:98:HIS:HB3	2.00	0.43
1:C:32:ASN:O	1:C:51:SER:HA	2.19	0.43
2:D:189:SER:O	2:D:193:GLN:HB2	2.18	0.43
1:L:38:GLN:HB2	1:L:48:LEU:HD11	2.01	0.43
1:L:55:LEU:HD22	1:L:59:VAL:HG13	2.01	0.43
2:H:28:LEU:HD21	2:H:76:VAL:HG23	2.00	0.43
1:L:48:LEU:O	1:L:59:VAL:HG11	2.18	0.42
2:B:127:PRO:HG3	2:B:139:LEU:HB3	2.01	0.42
2:D:202:LYS:NZ	4:D:329:HOH:O	2.52	0.42
2:H:79:LYS:HE3	2:H:81:THR:HG22	2.01	0.42
1:A:64:LYS:HB3	1:A:75:THR:HG23	2.00	0.42
2:B:95:ARG:O	2:B:102:THR:HA	2.19	0.42
2:H:71:ALA:HB2	2:H:77:ASP:OD2	2.19	0.42
2:H:189:SER:HB3	4:H:364:HOH:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:71:ALA:HB2	2:B:77:ASP:CG	2.45	0.42
2:D:127:PRO:HG3	2:D:139:LEU:HB3	2.01	0.42
1:C:64:LYS:HB3	1:C:75:THR:CG2	2.49	0.42
2:B:88:THR:HG23	2:B:111:THR:HA	2.02	0.42
2:F:57:TRP:HZ2	3:P:232:PRO:HA	1.85	0.42
2:H:132:THR:HG23	2:H:136:THR:C	2.45	0.41
1:A:7:PRO:HD3	1:E:201:GLN:HG2	2.01	0.41
2:D:57:TRP:HH2	3:O:233:GLU:HG2	1.84	0.41
1:A:49:ILE:HG12	1:A:55:LEU:HD23	2.02	0.41
1:L:18:VAL:HG22	1:L:76:ILE:HB	2.01	0.41
1:A:144:ARG:CZ	1:A:165:VAL:HG21	2.50	0.41
2:B:185:VAL:HG11	2:B:195:TYR:CE1	2.55	0.41
1:E:41:PRO:HG2	1:E:167:GLU:HG3	2.03	0.41
2:H:194:THR:HG23	2:H:211:LYS:HE3	2.02	0.41
1:E:3:LEU:HD21	1:E:91:GLY:HA3	2.03	0.41
1:E:55:LEU:HD21	1:E:63:PHE:O	2.21	0.41
1:A:39:GLN:NE2	2:B:38:GLN:OE1	2.52	0.40
1:A:209:LYS:HA	1:A:209:LYS:HD3	1.90	0.40
2:B:71:ALA:HB3	2:B:75:THR:OG1	2.22	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	213/216 (99%)	204 (96%)	9 (4%)	0	100	100
1	C	213/216 (99%)	205 (96%)	8 (4%)	0	100	100
1	E	213/216 (99%)	202 (95%)	11 (5%)	0	100	100
1	L	213/216 (99%)	202 (95%)	11 (5%)	0	100	100
2	B	212/223 (95%)	204 (96%)	7 (3%)	1 (0%)	24	27

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	D	212/223 (95%)	205 (97%)	6 (3%)	1 (0%)	24	27
2	F	212/223 (95%)	205 (97%)	6 (3%)	1 (0%)	24	27
2	H	212/223 (95%)	206 (97%)	5 (2%)	1 (0%)	24	27
3	M	4/14 (29%)	4 (100%)	0	0	100	100
3	N	4/14 (29%)	4 (100%)	0	0	100	100
3	O	4/14 (29%)	4 (100%)	0	0	100	100
3	P	4/14 (29%)	4 (100%)	0	0	100	100
All	All	1716/1812 (95%)	1649 (96%)	63 (4%)	4 (0%)	43	51

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	97	LEU
2	B	97	LEU
2	D	97	LEU
2	F	97	LEU

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	181/182 (100%)	177 (98%)	4 (2%)	45	61
1	C	181/182 (100%)	179 (99%)	2 (1%)	65	79
1	E	181/182 (100%)	179 (99%)	2 (1%)	65	79
1	L	181/182 (100%)	177 (98%)	4 (2%)	45	61
2	B	179/188 (95%)	175 (98%)	4 (2%)	45	61
2	D	179/188 (95%)	176 (98%)	3 (2%)	53	69
2	F	179/188 (95%)	176 (98%)	3 (2%)	53	69
2	H	179/188 (95%)	176 (98%)	3 (2%)	53	69
3	M	4/12 (33%)	4 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	N	4/12 (33%)	4 (100%)	0	100	100
3	O	4/12 (33%)	4 (100%)	0	100	100
3	P	4/12 (33%)	4 (100%)	0	100	100
All	All	1456/1528 (95%)	1431 (98%)	25 (2%)	53	69

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	6	THR
1	L	19	THR
1	L	30	ASN
1	L	125	GLU
2	H	11	VAL
2	H	80	VAL
2	H	103	VAL
1	A	19	THR
1	A	30	ASN
1	A	61	SER
1	A	205	SER
2	B	60	SER
2	B	67	ILE
2	B	132	THR
2	B	192	THR
1	C	19	THR
1	C	144	ARG
2	D	67	ILE
2	D	136	THR
2	D	192	THR
1	E	19	THR
1	E	30	ASN
2	F	42	LYS
2	F	103	VAL
2	F	132	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	39	GLN
1	L	71	GLN
2	H	38	GLN

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Mol	Chain	Res	Type
2	H	200	ASN
1	A	21	ASN
1	A	23	GLN
2	B	55	ASN
2	B	193	GLN
1	C	38	GLN
1	E	162	GLN
1	E	201	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

8 non-standard protein/DNA/RNA residues 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$
2	PCA	H	1	2	7,8,9	1.92	1 (14%)	9,10,12	2.35	5 (55%)
2	PCA	D	1	2	7,8,9	1.91	1 (14%)	9,10,12	2.25	5 (55%)
1	PCA	A	1	1	7,8,9	1.89	1 (14%)	9,10,12	2.14	5 (55%)
1	PCA	E	1	1	7,8,9	1.93	1 (14%)	9,10,12	2.12	5 (55%)
2	PCA	B	1	2	7,8,9	1.92	1 (14%)	9,10,12	2.28	5 (55%)
2	PCA	F	1	2	7,8,9	1.92	1 (14%)	9,10,12	2.40	6 (66%)
1	PCA	L	1	1	7,8,9	1.93	1 (14%)	9,10,12	2.10	4 (44%)
1	PCA	C	1	1	7,8,9	1.90	1 (14%)	9,10,12	2.13	5 (55%)

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
2	PCA	H	1	2	-	0/0/11/13	0/1/1/1
2	PCA	D	1	2	-	0/0/11/13	0/1/1/1
1	PCA	A	1	1	-	0/0/11/13	0/1/1/1
1	PCA	E	1	1	-	0/0/11/13	0/1/1/1
2	PCA	B	1	2	-	0/0/11/13	0/1/1/1
2	PCA	F	1	2	-	0/0/11/13	0/1/1/1
1	PCA	L	1	1	-	0/0/11/13	0/1/1/1
1	PCA	C	1	1	-	0/0/11/13	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	L	1	PCA	CD-N	4.98	1.46	1.34
2	B	1	PCA	CD-N	4.97	1.46	1.34
2	F	1	PCA	CD-N	4.97	1.46	1.34
2	D	1	PCA	CD-N	4.95	1.46	1.34
1	A	1	PCA	CD-N	4.94	1.46	1.34
2	H	1	PCA	CD-N	4.93	1.46	1.34
1	E	1	PCA	CD-N	4.92	1.46	1.34
1	C	1	PCA	CD-N	4.92	1.46	1.34

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	1	PCA	CB-CA-C	-3.72	107.55	112.66
2	H	1	PCA	CB-CA-C	-3.69	107.59	112.66
2	B	1	PCA	CB-CA-C	-3.08	108.43	112.66
2	B	1	PCA	OE-CD-CG	-3.08	121.22	126.72
1	A	1	PCA	OE-CD-CG	-3.00	121.36	126.72
2	D	1	PCA	CB-CA-C	-3.00	108.54	112.66
1	L	1	PCA	CB-CA-N	2.99	111.46	103.24
1	E	1	PCA	CB-CA-N	2.98	111.44	103.24
2	F	1	PCA	CA-N-CD	-2.97	103.40	113.58
1	L	1	PCA	OE-CD-CG	-2.97	121.42	126.72
2	D	1	PCA	OE-CD-CG	-2.96	121.43	126.72
2	H	1	PCA	OE-CD-CG	-2.94	121.47	126.72
2	F	1	PCA	OE-CD-CG	-2.93	121.49	126.72
1	C	1	PCA	OE-CD-CG	-2.92	121.51	126.72
1	E	1	PCA	OE-CD-CG	-2.92	121.51	126.72
1	C	1	PCA	CA-N-CD	-2.91	103.61	113.58
1	A	1	PCA	CA-N-CD	-2.90	103.64	113.58
1	C	1	PCA	CB-CA-N	2.90	111.22	103.24
2	D	1	PCA	CA-N-CD	-2.90	103.66	113.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1	PCA	CB-CA-N	2.89	111.19	103.24
2	H	1	PCA	CA-N-CD	-2.89	103.70	113.58
1	L	1	PCA	CA-N-CD	-2.87	103.74	113.58
2	B	1	PCA	CA-N-CD	-2.86	103.80	113.58
1	E	1	PCA	CA-N-CD	-2.83	103.88	113.58
2	B	1	PCA	CB-CA-N	2.80	110.94	103.24
2	D	1	PCA	CB-CA-N	2.77	110.85	103.24
1	A	1	PCA	CG-CD-N	2.65	114.86	108.39
2	B	1	PCA	CG-CD-N	2.63	114.83	108.39
1	C	1	PCA	CG-CD-N	2.63	114.81	108.39
1	L	1	PCA	CG-CD-N	2.62	114.80	108.39
2	H	1	PCA	CB-CA-N	2.62	110.45	103.24
2	F	1	PCA	CB-CA-N	2.61	110.41	103.24
2	D	1	PCA	CG-CD-N	2.59	114.74	108.39
1	E	1	PCA	CG-CD-N	2.59	114.72	108.39
2	H	1	PCA	CG-CD-N	2.52	114.55	108.39
2	F	1	PCA	CG-CD-N	2.51	114.52	108.39
1	A	1	PCA	CB-CA-C	-2.37	109.40	112.66
2	F	1	PCA	O-C-CA	-2.33	118.77	124.77
1	C	1	PCA	CB-CA-C	-2.17	109.68	112.66
1	E	1	PCA	CB-CA-C	-2.15	109.71	112.66

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

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	214/216 (99%)	-0.72	1 (0%) 87 85	21, 43, 70, 126	0
1	C	214/216 (99%)	-0.73	3 (1%) 73 71	21, 43, 69, 120	0
1	E	214/216 (99%)	-1.27	0 100 100	20, 36, 56, 69	0
1	L	214/216 (99%)	-1.24	0 100 100	23, 36, 56, 72	0
2	B	213/223 (95%)	-0.98	3 (1%) 73 71	20, 36, 69, 138	0
2	D	213/223 (95%)	-1.00	3 (1%) 73 71	21, 36, 73, 160	0
2	F	213/223 (95%)	-1.07	2 (0%) 81 79	20, 36, 67, 182	0
2	H	213/223 (95%)	-1.04	4 (1%) 66 63	21, 36, 72, 161	0
3	M	6/14 (42%)	0.18	0 100 100	42, 52, 69, 106	0
3	N	6/14 (42%)	0.84	1 (16%) 4 3	53, 62, 78, 89	0
3	O	6/14 (42%)	0.48	1 (16%) 4 3	49, 58, 68, 87	0
3	P	6/14 (42%)	0.14	0 100 100	38, 50, 55, 97	0
All	All	1732/1812 (95%)	-0.99	18 (1%) 79 77	20, 38, 69, 182	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	134	GLY	4.1
2	B	133	SER	4.0
1	A	186	ALA	3.9
2	D	134	GLY	3.8
2	B	132	THR	3.7
2	F	132	THR	3.5
2	F	133	SER	3.4
2	H	135	GLY	3.2
2	H	132	THR	3.0
2	D	132	THR	2.8
2	H	131	SER	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	186	ALA	2.5
2	D	133	SER	2.4
1	C	189	GLU	2.3
3	N	232	PRO	2.3
1	C	111	THR	2.1
3	O	232	PRO	2.1
2	H	133	SER	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
2	PCA	D	1	8/9	0.98	0.07	39,52,54,55	0
2	PCA	H	1	8/9	0.99	0.03	41,49,52,58	0
1	PCA	A	1	8/9	0.99	0.03	27,34,43,53	0
2	PCA	B	1	8/9	0.99	0.05	41,46,47,48	0
1	PCA	C	1	8/9	0.99	0.03	30,36,45,52	0
1	PCA	L	1	8/9	0.99	0.03	30,32,34,37	0
1	PCA	E	1	8/9	0.99	0.02	29,31,35,36	0
2	PCA	F	1	8/9	0.99	0.04	39,47,52,56	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.