



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:31 PM UTC

PDB ID : 8G8D / pdb_00008g8d
Title : Crystal structure of DH1346 Fab in complex with HIV proximal MPER peptide
Authors : Niyongabo, A.; Janus, B.M.; Ofek, G.
Deposited on : 2023-02-17
Resolution : 2.02 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
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

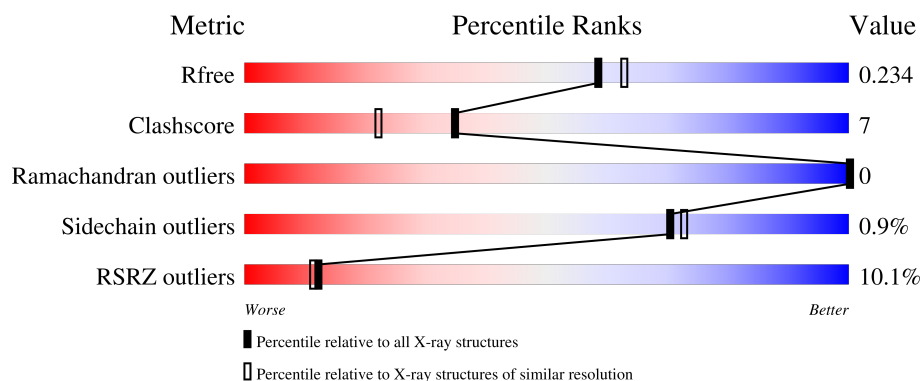
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.02 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	13299 (2.04-2.00)
Clashscore	190562	1022 (2.02-2.02)
Ramachandran outliers	187476	1014 (2.02-2.02)
Sidechain outliers	187428	1014 (2.02-2.02)
RSRZ outliers	180081	13314 (2.04-2.00)

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	230	3% 92% 7%
1	H	230	4% 88% 10% .
2	B	215	6% 82% 16% .
2	L	215	27% 80% 17% .
3	C	34	12% 74% 12% 15%

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Mol	Chain	Length	Quality of chain
3	P	34	15% 74% 9% • 15%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 14370 atoms, of which 6798 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 DH1346 heavy chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	H	228	Total	C	H	N	O	S	0	0	0
			3386	1092	1672	284	331	7			
1	A	230	Total	C	H	N	O	S	0	0	0
			3401	1098	1677	286	333	7			

- Molecule 2 is a protein called DH1346 light chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	L	208	Total	C	H	N	O	S	0	0	0
			3017	973	1464	261	315	4			
2	B	211	Total	C	H	N	O	S	0	0	0
			3093	995	1503	267	324	4			

- Molecule 3 is a protein called gp41 MPER peptide.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	P	29	Total	C	H	N	O	0	0	0
			507	180	241	40	46			
3	C	29	Total	C	H	N	O	0	0	0
			507	180	241	40	46			

- Molecule 4 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	H	3	Total	Na	0	0
			3	3		
4	L	3	Total	Na	0	0
			3	3		
4	A	2	Total	Na	0	0
			2	2		
4	B	2	Total	Na	0	0
			2	2		

- Molecule 5 is FLUORIDE ION (CCD ID: F) (formula: F).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	L	1	Total F 1 1	0	0

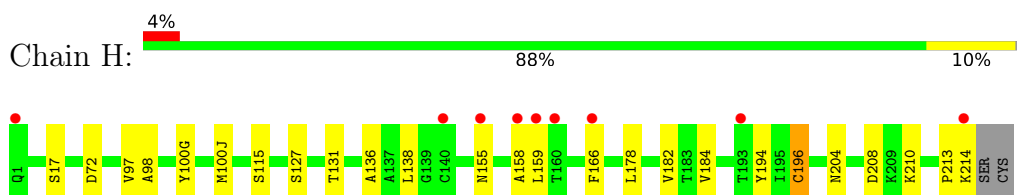
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	H	132	Total O 132 132	0	0
6	L	74	Total O 74 74	0	0
6	P	3	Total O 3 3	0	0
6	A	144	Total O 144 144	0	0
6	B	82	Total O 82 82	0	0
6	C	13	Total O 13 13	0	0

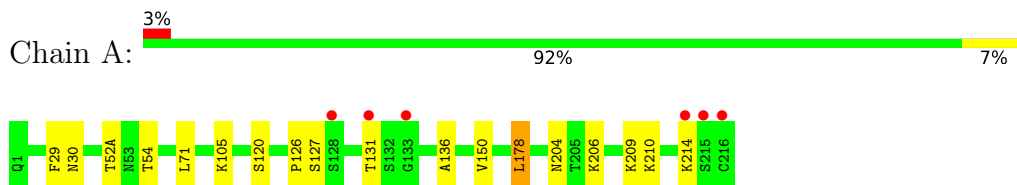
3 Residue-property plots [i](#)

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.

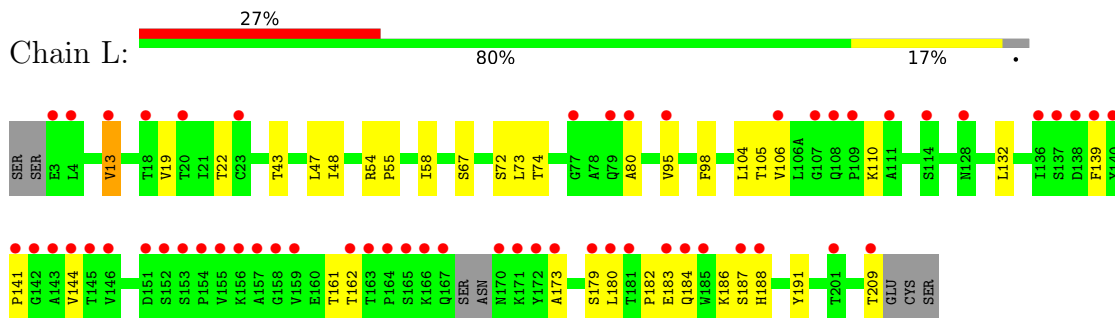
- Molecule 1: DH1346 heavy chain



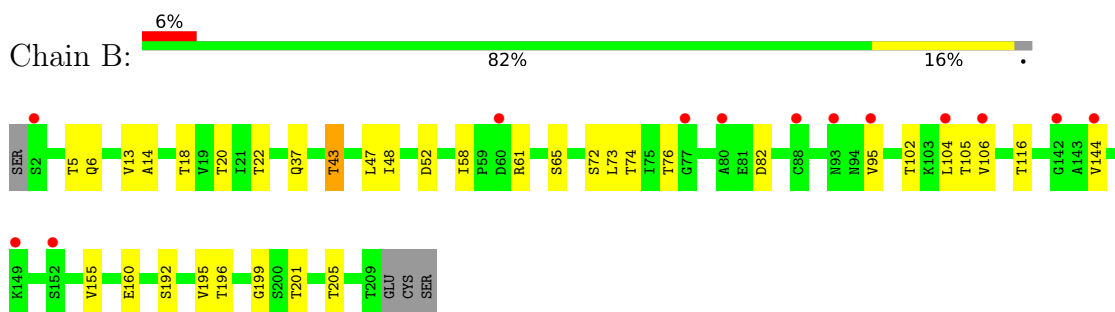
- Molecule 1: DH1346 heavy chain



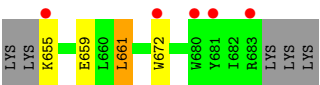
- Molecule 2: DH1346 light chain



- Molecule 2: DH1346 light chain



- Molecule 3: gp41 MPER peptide



● Molecule 3: gp41 MPER peptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	41.07Å 115.51Å 231.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.98 – 2.02 37.98 – 2.02	Depositor EDS
% Data completeness (in resolution range)	91.7 (37.98-2.02) 91.7 (37.98-2.02)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.49 (at 2.03Å)	Xtriage
Refinement program	PHENIX 1.20	Depositor
R, R_{free}	0.198 , 0.234 0.198 , 0.234	Depositor DCC
R_{free} test set	5097 reflections (6.97%)	wwPDB-VP
Wilson B-factor (Å ²)	30.0	Xtriage
Anisotropy	0.241	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 56.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	14370	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.23% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NA, F

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	0.18	0/1770	0.39	0/2415
1	H	0.24	0/1760	0.40	0/2401
2	B	0.21	0/1631	0.41	0/2236
2	L	0.24	0/1592	0.45	2/2182 (0.1%)
3	C	0.17	0/277	0.36	0/380
3	P	0.45	0/277	0.52	0/380
All	All	0.23	0/7307	0.41	2/9994 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	141	PRO	N-CA-C	-6.07	101.69	111.03
2	L	141	PRO	CA-C-O	-5.58	115.26	121.23

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	1724	1677	1677	19	0
1	H	1714	1672	1672	23	0
2	B	1590	1503	1508	30	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	L	1553	1464	1463	34	0
3	C	266	241	241	2	2
3	P	266	241	241	5	2
4	A	2	0	0	0	0
4	B	2	0	0	0	0
4	H	3	0	0	0	0
4	L	3	0	0	0	0
5	L	1	0	0	0	0
6	A	144	0	0	9	1
6	B	82	0	0	9	0
6	C	13	0	0	0	0
6	H	132	0	0	4	1
6	L	74	0	0	13	0
6	P	3	0	0	0	0
All	All	7572	6798	6802	104	3

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 7.

All (104) 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:H:97:VAL:HG11	3:P:661:LEU:HD22	1.53	0.89
2:L:67:SER:OG	6:L:401:HOH:O	1.91	0.88
2:B:160:GLU:OE1	6:B:402:HOH:O	1.90	0.87
2:B:199:GLY:O	6:B:401:HOH:O	1.90	0.86
2:L:54:ARG:NH1	6:L:403:HOH:O	2.12	0.83
2:L:95:VAL:HG21	3:P:672:TRP:CE3	2.15	0.81
1:A:209:LYS:NZ	6:A:402:HOH:O	2.13	0.80
1:A:120:SER:OG	6:A:401:HOH:O	2.02	0.78
2:L:209:THR:OG1	6:L:402:HOH:O	2.01	0.78
1:A:131:THR:HG22	1:A:136:ALA:HB2	1.67	0.76
1:H:100(J):MET:HE1	2:L:98:PHE:CZ	2.26	0.70
2:B:5:THR:HG23	6:B:413:HOH:O	1.90	0.70
1:H:115:SER:O	6:H:401:HOH:O	2.10	0.70
1:H:166:PHE:O	6:H:402:HOH:O	2.11	0.68
2:B:52:ASP:OD1	6:B:404:HOH:O	2.11	0.68
2:B:155:VAL:O	6:B:403:HOH:O	2.11	0.67
1:H:72:ASP:OD2	6:H:403:HOH:O	2.13	0.67
1:H:127:SER:O	1:H:131:THR:HG23	1.95	0.66
1:A:126:PRO:O	6:A:403:HOH:O	2.13	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:THR:HG22	1:A:136:ALA:CB	2.28	0.64
2:L:55:PRO:HB2	2:L:58:ILE:HD13	1.81	0.62
2:B:37:GLN:CB	2:B:47:LEU:HD11	2.29	0.62
2:L:74:THR:HG23	6:L:416:HOH:O	2.00	0.62
2:B:116:THR:HG22	6:B:415:HOH:O	2.00	0.60
2:L:47:LEU:C	2:L:48:ILE:HD12	2.28	0.59
2:B:37:GLN:HB2	2:B:47:LEU:HD11	1.84	0.59
1:A:150:VAL:CG2	1:A:178:LEU:HD21	2.32	0.59
2:B:201:THR:HG23	6:B:401:HOH:O	2.03	0.58
2:B:47:LEU:C	2:B:48:ILE:HD12	2.28	0.58
1:H:159:LEU:CD2	1:H:182:VAL:HG21	2.35	0.56
1:H:213:PRO:O	1:H:214:LYS:HB3	2.06	0.55
2:B:104:LEU:HD23	2:B:105:THR:N	2.20	0.55
2:B:192:SER:OG	2:B:205:THR:HG22	2.07	0.55
1:H:131:THR:HG22	1:H:136:ALA:HB2	1.89	0.55
1:A:52(A):THR:HA	1:A:71:LEU:HD11	1.88	0.54
2:B:13:VAL:HG12	2:B:104:LEU:HD21	1.89	0.54
2:L:182:PRO:HD2	6:L:470:HOH:O	2.08	0.54
2:L:80:ALA:HA	2:L:106:VAL:HG21	1.91	0.53
2:L:161:THR:C	6:L:406:HOH:O	2.52	0.53
2:B:20:THR:HG22	2:B:74:THR:HG22	1.91	0.53
2:L:132:LEU:HD13	2:L:180:LEU:HD12	1.89	0.53
2:B:106:VAL:O	2:B:106:VAL:HG23	2.08	0.53
1:H:131:THR:HG22	1:H:136:ALA:CB	2.39	0.53
2:L:54:ARG:HD2	6:L:403:HOH:O	2.09	0.52
2:B:48:ILE:HD13	2:B:73:LEU:CD1	2.40	0.52
1:H:100(J):MET:HE1	2:L:98:PHE:HZ	1.74	0.52
2:B:144:VAL:HG13	2:B:195:VAL:HG13	1.92	0.51
2:B:196:THR:OG1	2:B:201:THR:HG22	2.10	0.51
2:B:144:VAL:CG1	2:B:195:VAL:HG13	2.39	0.51
3:P:655:LYS:HZ3	3:P:659:GLU:HB3	1.74	0.51
2:B:18:THR:HG22	2:B:76:THR:HA	1.92	0.51
2:L:104:LEU:HD23	2:L:105:THR:N	2.26	0.50
2:L:186:LYS:O	6:L:405:HOH:O	2.20	0.50
1:H:204:ASN:HB3	1:A:210:LYS:HB3	1.93	0.50
2:L:139:PHE:CE1	2:L:173:ALA:HA	2.46	0.50
2:B:61:ARG:NH1	2:B:82:ASP:OD2	2.45	0.49
1:H:178:LEU:C	1:H:178:LEU:HD12	2.37	0.49
2:L:95:VAL:HG21	3:P:672:TRP:CD2	2.47	0.49
1:H:159:LEU:HD21	1:H:182:VAL:HG21	1.93	0.49
1:H:210:LYS:HB3	1:A:204:ASN:HB3	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:48:ILE:HD13	2:L:73:LEU:CD1	2.43	0.48
2:L:106:VAL:HG23	2:L:106:VAL:O	2.14	0.48
1:A:178:LEU:C	1:A:178:LEU:HD12	2.39	0.48
2:B:65:SER:OG	6:B:405:HOH:O	2.20	0.48
2:B:13:VAL:HG22	2:B:14:ALA:N	2.29	0.47
2:L:58:ILE:N	2:L:58:ILE:HD12	2.29	0.47
3:C:672:TRP:O	3:C:676:THR:HG23	2.15	0.47
1:H:155:ASN:HB2	1:H:158:ALA:HB3	1.97	0.47
2:B:6:GLN:HG2	2:B:102:THR:OG1	2.15	0.47
1:A:127:SER:O	1:A:131:THR:HG23	2.15	0.46
2:L:183:GLU:N	6:L:404:HOH:O	2.49	0.45
2:L:188:HIS:HB2	2:L:191:TYR:CE1	2.51	0.45
2:L:179:SER:O	2:L:180:LEU:HD23	2.16	0.45
1:H:138:LEU:C	1:H:138:LEU:HD12	2.41	0.45
6:A:483:HOH:O	3:C:662:GLU:HG2	2.17	0.45
1:A:206:LYS:NZ	6:A:404:HOH:O	2.20	0.45
2:L:13:VAL:HG21	2:L:19:VAL:HG11	1.98	0.45
1:H:97:VAL:CG1	3:P:661:LEU:HD22	2.38	0.45
2:B:58:ILE:HD12	2:B:58:ILE:N	2.32	0.44
2:L:144:VAL:HA	6:L:410:HOH:O	2.17	0.43
1:H:155:ASN:CB	1:H:158:ALA:HB3	2.49	0.43
2:L:67:SER:CB	6:L:401:HOH:O	2.61	0.43
1:A:30:ASN:ND2	6:A:405:HOH:O	2.21	0.43
1:A:206:LYS:HD2	6:A:404:HOH:O	2.19	0.42
2:L:184:GLN:HA	2:L:187:SER:HB3	2.00	0.42
1:H:17:SER:N	6:H:405:HOH:O	2.29	0.42
1:A:150:VAL:HG22	1:A:178:LEU:HD21	2.01	0.42
2:L:22:THR:HG22	2:L:72:SER:OG	2.19	0.42
1:H:184:VAL:HG11	1:H:194:TYR:CE1	2.55	0.42
2:B:20:THR:CG2	2:B:74:THR:HG22	2.49	0.41
2:B:22:THR:HG22	2:B:72:SER:OG	2.20	0.41
2:B:95:VAL:HG12	6:B:443:HOH:O	2.20	0.41
1:A:105:LYS:HD3	2:B:43:THR:HG21	2.02	0.41
1:H:196:CYS:O	1:H:208:ASP:HA	2.20	0.41
2:B:13:VAL:CG1	2:B:104:LEU:HD21	2.51	0.41
1:A:29:PHE:CZ	1:A:71:LEU:HD22	2.56	0.41
2:L:188:HIS:ND1	2:L:191:TYR:OH	2.48	0.41
2:L:132:LEU:HD12	2:L:132:LEU:N	2.37	0.40
2:L:184:GLN:HG3	6:L:404:HOH:O	2.20	0.40
1:A:54:THR:HA	6:A:522:HOH:O	2.21	0.40
1:H:98:ALA:HB3	1:H:100(G):TYR:CZ	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:214:LYS:NZ	6:A:418:HOH:O	2.47	0.40
2:L:104:LEU:HD23	2:L:104:LEU:C	2.47	0.40
2:L:162:THR:C	6:L:406:HOH:O	2.63	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:P:655:LYS:NZ	3:C:681:TYR:OH[2_455]	2.08	0.12
3:P:655:LYS:HZ3	3:C:681:TYR:OH[2_455]	1.56	0.04
6:H:471:HOH:O	6:A:525:HOH:O[3_554]	2.18	0.02

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	228/230 (99%)	226 (99%)	2 (1%)	0	100	100
1	H	226/230 (98%)	224 (99%)	2 (1%)	0	100	100
2	B	209/215 (97%)	202 (97%)	7 (3%)	0	100	100
2	L	204/215 (95%)	191 (94%)	13 (6%)	0	100	100
3	C	27/34 (79%)	27 (100%)	0	0	100	100
3	P	27/34 (79%)	27 (100%)	0	0	100	100
All	All	921/958 (96%)	897 (97%)	24 (3%)	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	189/193 (98%)	188 (100%)	1 (0%)	81	81
1	H	189/193 (98%)	188 (100%)	1 (0%)	81	81
2	B	177/185 (96%)	176 (99%)	1 (1%)	78	79
2	L	170/185 (92%)	167 (98%)	3 (2%)	51	52
3	C	27/33 (82%)	27 (100%)	0	100	100
3	P	27/33 (82%)	26 (96%)	1 (4%)	30	24
All	All	779/822 (95%)	772 (99%)	7 (1%)	70	73

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	196	CYS
2	L	13	VAL
2	L	43	THR
2	L	110	LYS
3	P	661	LEU
1	A	178	LEU
2	B	43	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	H	56	ASN
2	L	79	GLN
2	L	93	ASN
2	L	197	HIS
2	B	95(C)	ASN

5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 11 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 [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	230/230 (100%)	-0.04	6 (2%) 57 57	21, 34, 59, 163	0
1	H	228/230 (99%)	0.06	9 (3%) 43 43	22, 38, 63, 79	0
2	B	211/215 (98%)	0.64	13 (6%) 26 25	25, 52, 81, 97	0
2	L	208/215 (96%)	1.32	57 (27%) 1 1	26, 64, 102, 114	0
3	C	29/34 (85%)	0.64	4 (13%) 6 5	29, 46, 81, 102	0
3	P	29/34 (85%)	0.70	5 (17%) 4 3	30, 46, 94, 105	0
All	All	935/958 (97%)	0.49	94 (10%) 12 11	21, 43, 88, 163	0

All (94) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	216	CYS	8.6
1	A	215	SER	6.4
2	L	3	GLU	6.2
2	L	159	VAL	5.2
2	L	155	VAL	5.1
2	L	164	PRO	5.0
2	L	139	PHE	5.0
2	L	140	TYR	4.8
2	L	142	GLY	4.6
2	L	128	ASN	4.6
2	L	172	TYR	4.3
2	L	187	SER	4.2
2	L	141	PRO	4.1
2	L	109	PRO	4.0
2	L	144	VAL	3.9
2	L	156	LYS	3.8
2	L	137	SER	3.8
2	L	180	LEU	3.8
3	P	683	ARG	3.7

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Mol	Chain	Res	Type	RSRZ
2	B	95	VAL	3.7
2	B	2	SER	3.7
3	P	680	TRP	3.7
2	L	163	THR	3.7
2	L	165	SER	3.6
2	L	108	GLN	3.5
3	C	680	TRP	3.4
3	P	681	TYR	3.4
2	B	93	ASN	3.3
1	A	131	THR	3.3
2	L	179	SER	3.3
2	L	170	ASN	3.3
2	L	106	VAL	3.2
2	L	157	ALA	3.2
2	L	183	GLU	3.2
2	B	106	VAL	3.2
2	L	167	GLN	3.2
2	L	184	GLN	3.2
1	H	140	CYS	3.1
1	H	158	ALA	3.1
2	B	77	GLY	3.1
3	P	655	LYS	3.0
2	L	4	LEU	3.0
2	L	151	ASP	3.0
3	C	666	TRP	3.0
2	L	162	THR	3.0
2	L	188	HIS	3.0
2	L	13	VAL	2.9
2	L	171	LYS	2.8
2	L	153	SER	2.8
2	L	107	GLY	2.8
2	L	95	VAL	2.8
1	H	214	LYS	2.7
2	L	143	ALA	2.7
2	L	146	VAL	2.7
2	L	209	THR	2.7
2	L	158	GLY	2.6
2	L	181	THR	2.6
1	H	159	LEU	2.6
2	L	138	ASP	2.5
2	L	145	THR	2.5
2	L	173	ALA	2.5

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Mol	Chain	Res	Type	RSRZ
1	H	1	GLN	2.4
2	B	80	ALA	2.4
2	B	149	LYS	2.4
2	B	142	GLY	2.4
1	A	133	GLY	2.4
2	L	77	GLY	2.4
2	L	166	LYS	2.4
2	B	152	SER	2.4
2	L	201	THR	2.3
1	H	160	THR	2.3
2	L	18	THR	2.3
3	C	683	ARG	2.3
1	A	128	SER	2.3
1	H	193	THR	2.3
2	L	23	CYS	2.3
2	B	104	LEU	2.3
3	C	681	TYR	2.3
3	P	672	TRP	2.3
2	L	114	SER	2.2
2	L	136	ILE	2.2
2	L	80	ALA	2.2
2	L	79	GLN	2.2
2	L	152	SER	2.1
2	B	60	ASP	2.1
1	A	214	LYS	2.1
2	L	154	PRO	2.1
2	B	144	VAL	2.1
2	L	20	THR	2.1
2	L	111	ALA	2.1
1	H	155	ASN	2.0
1	H	166	PHE	2.0
2	L	185	TRP	2.0
2	B	88	CYS	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 oligosaccharides 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
4	NA	L	303	1/1	0.31	0.32	104,104,104,104	0
4	NA	H	303	1/1	0.79	0.20	60,60,60,60	0
4	NA	B	301	1/1	0.80	0.17	67,67,67,67	0
4	NA	L	301	1/1	0.84	0.11	56,56,56,56	0
4	NA	A	301	1/1	0.91	0.10	44,44,44,44	0
4	NA	H	302	1/1	0.93	0.12	40,40,40,40	0
4	NA	A	302	1/1	0.94	0.10	45,45,45,45	0
4	NA	H	301	1/1	0.94	0.15	35,35,35,35	0
4	NA	L	302	1/1	0.95	0.14	64,64,64,64	0
4	NA	B	302	1/1	0.97	0.05	49,49,49,49	0
5	F	L	304	1/1	0.97	0.06	31,31,31,31	0

6.5 Other polymers [i](#)

There are no such residues in this entry.