



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:08 PM UTC

PDB ID : 5UY3 / pdb_00005uy3
Title : Crystal structure of human Fab PGT144, a broadly reactive and potent HIV-1 neutralizing antibody
Authors : Julien, J.-P.; Lee, J.H.; Wilson, I.A.
Deposited on : 2017-02-23
Resolution : 2.90 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

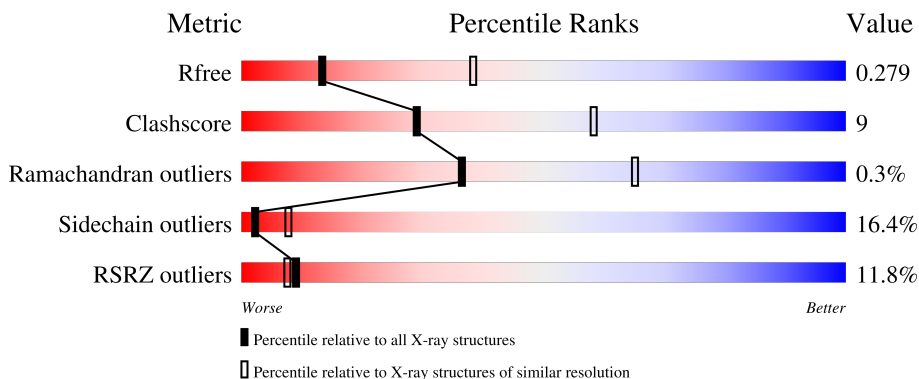
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	B	216	12% 73% 24% .
1	L	216	8% 70% 25% 5%
2	A	238	11% 71% 20% 5% .
2	H	238	14% 61% 30% 5% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6856 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 Antibody PGT144 Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	216	Total	C	N	O	S	0	0	0
			1655	1031	292	326	6			
1	B	216	Total	C	N	O	S	0	0	0
			1655	1031	292	326	6			

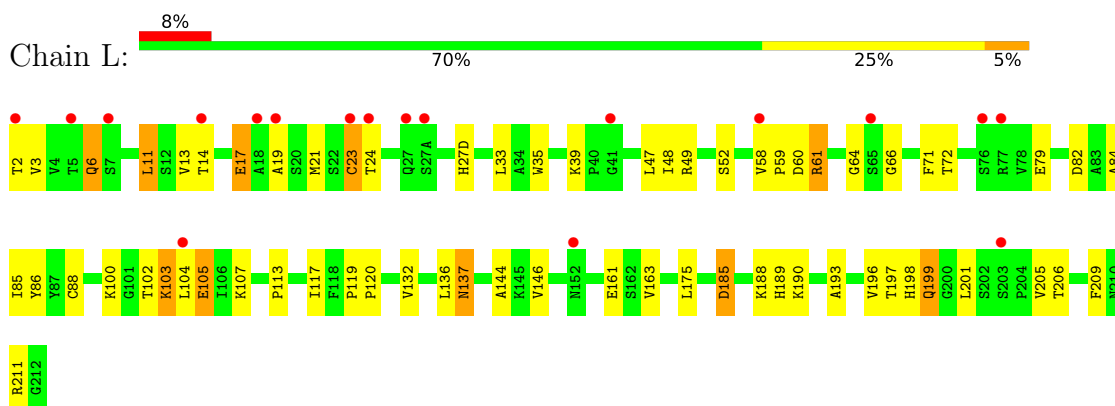
- Molecule 2 is a protein called Antibody PGT144 Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	230	Total	C	N	O	S	0	0	0
			1773	1116	302	349	6			
2	A	230	Total	C	N	O	S	0	0	0
			1773	1116	302	349	6			

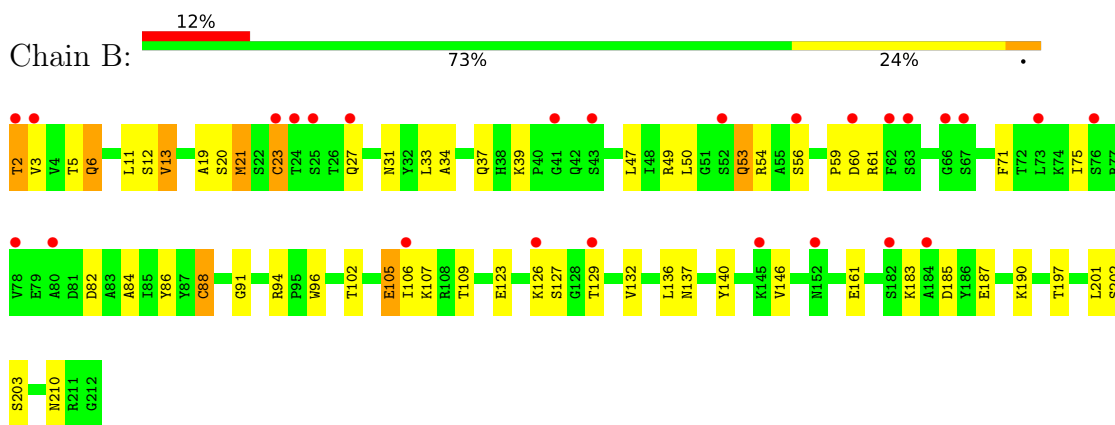
3 Residue-property plots [i](#)

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.

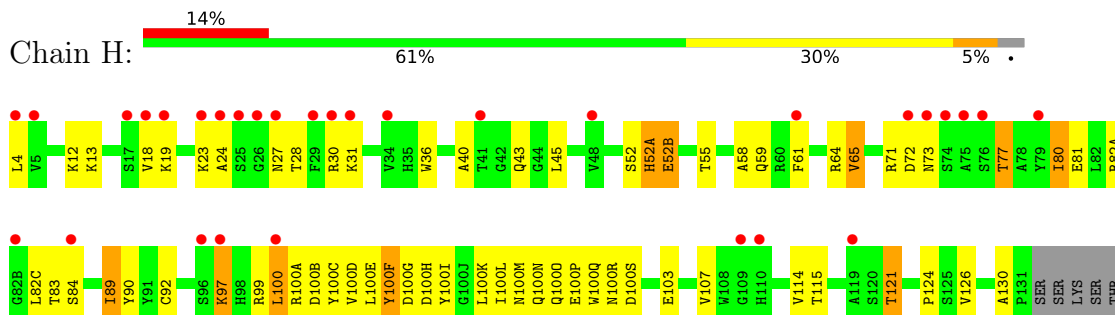
- Molecule 1: Antibody PGT144 Fab light chain

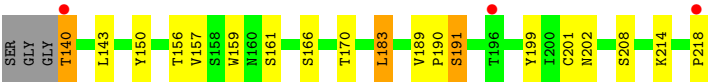


- Molecule 1: Antibody PGT144 Fab light chain

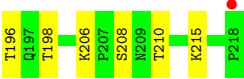
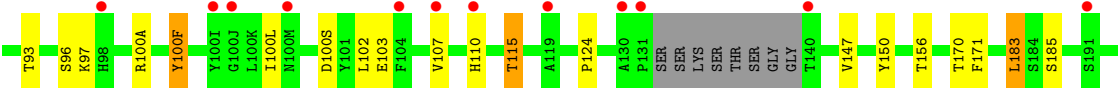
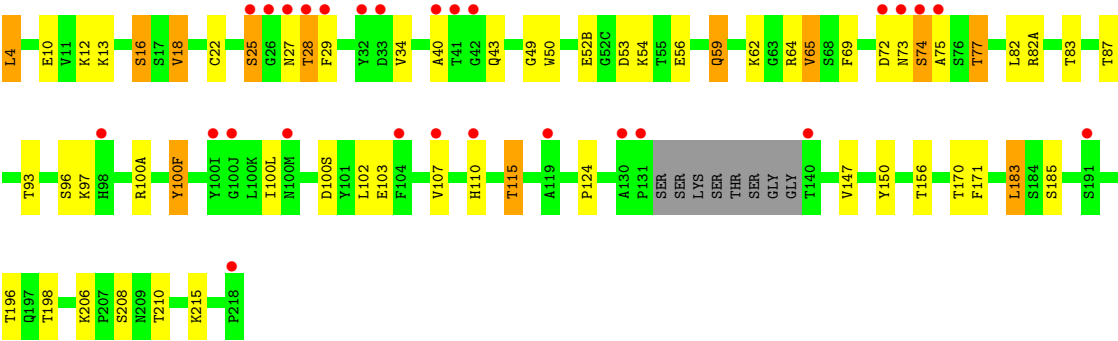


- Molecule 2: Antibody PGT144 Fab heavy chain





● Molecule 2: Antibody PGT144 Fab heavy chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	57.60Å 67.55Å 69.55Å 82.04° 74.89° 81.26°	Depositor
Resolution (Å)	36.29 – 2.90 36.29 – 2.90	Depositor EDS
% Data completeness (in resolution range)	96.6 (36.29-2.90) 92.6 (36.29-2.90)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.69 (at 2.90Å)	Xtriage
Refinement program	PHENIX (1.10_2155: ???)	Depositor
R, R_{free}	0.241 , 0.282 0.236 , 0.279	Depositor DCC
R_{free} test set	1065 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	34.7	Xtriage
Anisotropy	0.457	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 52.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	6856	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.69% 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: TYS

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	B	0.12	0/1691	0.33	0/2293
1	L	0.13	0/1691	0.33	0/2293
2	A	0.13	0/1799	0.35	0/2446
2	H	0.12	0/1799	0.36	0/2446
All	All	0.12	0/6980	0.34	0/9478

There are no bond length outliers.

There are no bond angle outliers.

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	B	1655	0	1621	20	0
1	L	1655	0	1621	38	0
2	A	1773	0	1706	29	0
2	H	1773	0	1707	34	0
All	All	6856	0	6655	118	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 9.

All (118) close contacts within the same asymmetric unit are listed below, sorted by their clash

magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:61:ARG:NH2	1:L:82:ASP:OD2	2.14	0.81
2:A:27:ASN:ND2	2:A:74:SER:O	2.14	0.80
1:B:190:LYS:HE3	1:B:210:ASN:HB3	1.68	0.75
1:L:49:ARG:NH1	2:H:103:GLU:OE2	2.20	0.75
1:B:49:ARG:NH1	2:A:103:GLU:OE2	2.25	0.70
2:H:52(A):HIS:HA	2:H:71:ARG:HH22	1.55	0.69
2:H:124:PRO:HB3	2:H:150:TYR:HB3	1.75	0.68
2:H:140:THR:HA	2:H:190:PRO:HA	1.76	0.67
2:A:52(B):GLU:OE1	2:A:97:LYS:NZ	2.25	0.67
2:A:147:VAL:HB	2:A:183:LEU:HD23	1.77	0.67
2:H:64:ARG:NH1	2:H:82(A):ARG:O	2.29	0.66
1:L:35:TRP:HB2	1:L:48:ILE:HB	1.77	0.65
1:B:39:LYS:HG2	1:B:84:ALA:HB2	1.77	0.65
2:A:124:PRO:HB3	2:A:150:TYR:HB3	1.80	0.63
2:A:59:GLN:HA	2:A:62:LYS:HG3	1.79	0.63
2:H:27:ASN:HD21	2:H:30:ARG:HG3	1.66	0.61
2:H:18:VAL:HG23	2:H:82(C):LEU:HD11	1.83	0.60
1:L:136:LEU:HD11	1:L:196:VAL:HG11	1.83	0.60
1:B:183:LYS:NZ	1:B:187:GLU:OE2	2.31	0.60
2:A:12:LYS:HG3	2:A:18:VAL:HG22	1.83	0.59
2:H:4:LEU:HD22	2:H:24:ALA:HA	1.84	0.58
1:B:61:ARG:NH2	1:B:82:ASP:OD2	2.36	0.58
2:H:24:ALA:HB1	2:H:28:THR:HB	1.86	0.58
1:L:21:MET:HE3	1:L:102:THR:HB	1.84	0.58
2:H:12:LYS:HG3	2:H:18:VAL:HG22	1.86	0.58
2:H:40:ALA:HB3	2:H:43:GLN:HB2	1.84	0.58
2:A:171:PHE:O	2:A:183:LEU:HD12	2.05	0.57
2:H:97:LYS:HE3	2:H:99:ARG:HB2	1.86	0.56
2:H:130:ALA:HB1	2:H:218:PRO:HA	1.88	0.55
1:L:47:LEU:HA	1:L:58:VAL:HG21	1.90	0.54
1:L:189:HIS:O	1:L:211:ARG:HD3	2.08	0.53
2:H:100(F):TYS:O2	2:H:100(F):TYS:HE1	2.08	0.53
2:H:30:ARG:HA	2:H:52(A):HIS:HB2	1.91	0.53
2:A:72:ASP:HB3	2:A:75:ALA:H	1.73	0.53
1:B:105:GLU:OE2	1:B:140:TYR:OH	2.13	0.53
1:L:21:MET:HE1	1:L:86:TYR:HB2	1.89	0.52
1:L:189:HIS:O	1:L:211:ARG:NH1	2.38	0.52
1:B:49:ARG:HG2	1:B:53:GLN:HB2	1.91	0.52
2:A:65:VAL:HG13	2:A:82:LEU:HD13	1.92	0.51
2:H:23:LYS:HG2	2:H:77:THR:HG23	1.92	0.51
2:A:72:ASP:HB2	2:A:77:THR:H	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:65:VAL:HA	2:H:81:GLU:O	2.12	0.50
1:L:11:LEU:O	1:L:105:GLU:HG3	2.10	0.50
1:L:39:LYS:HG2	1:L:84:ALA:HB2	1.93	0.49
2:A:100(F):TYS:HE1	2:A:100(L):ILE:HG12	1.93	0.49
1:L:144:ALA:HB2	1:L:198:HIS:HD2	1.77	0.49
2:A:87:THR:HG23	2:A:115:THR:HA	1.94	0.49
1:B:123:GLU:OE1	1:B:123:GLU:N	2.42	0.49
2:A:72:ASP:HB3	2:A:75:ALA:N	2.28	0.49
2:H:58:ALA:HB3	2:H:61:PHE:HD2	1.77	0.48
1:L:14:THR:O	1:L:17:GLU:HB2	2.13	0.48
1:L:185:ASP:HA	1:L:188:LYS:HG3	1.95	0.48
2:A:170:THR:HG23	2:A:185:SER:HB2	1.95	0.48
1:L:60:ASP:OD1	1:L:60:ASP:N	2.45	0.47
1:L:27(D):HIS:NE2	2:H:100:LEU:HD23	2.28	0.47
2:A:49:GLY:HA3	2:A:69:PHE:CZ	2.49	0.47
1:B:2:THR:HB	1:B:27:GLN:NE2	2.30	0.47
2:A:13:LYS:H	2:A:13:LYS:HD2	1.79	0.47
2:H:18:VAL:CG2	2:H:82(C):LEU:HD11	2.45	0.47
1:B:37:GLN:HB2	1:B:47:LEU:HD11	1.96	0.46
2:A:25:SER:O	2:A:28:THR:OG1	2.33	0.46
2:A:50:TRP:HE1	2:A:56:GLU:HB2	1.80	0.46
1:B:13:VAL:HG21	1:B:19:ALA:HB2	1.97	0.46
1:L:113:PRO:HB2	1:L:136:LEU:HD12	1.97	0.46
2:H:157:VAL:HA	2:H:202:ASN:O	2.16	0.46
1:L:66:GLY:HA3	1:L:71:PHE:HA	1.98	0.45
1:L:85:ILE:HG12	1:L:103:LYS:HD2	1.96	0.45
1:B:6:GLN:HG2	1:B:23:CYS:SG	2.56	0.45
1:L:47:LEU:HD12	1:L:48:ILE:HG12	1.98	0.45
1:L:137:ASN:HD22	1:L:137:ASN:HA	1.60	0.45
2:H:58:ALA:HB3	2:H:61:PHE:CD2	2.51	0.45
2:A:13:LYS:O	2:A:16:SER:OG	2.34	0.45
1:L:48:ILE:HG21	1:L:64:GLY:HA3	1.99	0.44
2:A:53:ASP:OD1	2:A:53:ASP:N	2.49	0.44
1:B:21:MET:HE1	1:B:86:TYR:HB2	1.99	0.44
2:A:54:LYS:HD2	2:A:100(S):ASP:OD2	2.17	0.44
2:H:52(A):HIS:NE2	2:H:52(B):GLU:OE2	2.51	0.44
2:H:90:TYR:HE2	2:H:114:VAL:HB	1.83	0.44
2:H:121:THR:HG22	2:H:208:SER:HB3	2.00	0.44
2:H:140:THR:N	2:H:191:SER:OG	2.51	0.44
1:B:31:ASN:HB3	1:B:71:PHE:CZ	2.53	0.43
1:L:61:ARG:H	1:L:61:ARG:HG2	1.44	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:19:LYS:HA	2:H:80:ILE:O	2.19	0.43
2:H:18:VAL:O	2:H:81:GLU:HA	2.19	0.43
2:A:208:SER:OG	2:A:210:THR:OG1	2.28	0.43
1:L:33:LEU:HD13	1:L:71:PHE:CD1	2.53	0.43
2:A:64:ARG:NH1	2:A:82(A):ARG:O	2.46	0.43
1:L:193:ALA:HB1	1:L:206:THR:HG22	2.00	0.42
1:L:199:GLN:HE21	1:L:199:GLN:HB2	1.67	0.42
2:H:189:VAL:HG11	2:H:199:TYR:CZ	2.54	0.42
1:L:120:PRO:HD3	1:L:132:VAL:HG22	2.00	0.42
1:L:21:MET:O	1:L:72:THR:HG23	2.20	0.42
1:L:119:PRO:HB3	1:L:209:PHE:CE1	2.53	0.42
1:L:163:VAL:HG22	1:L:175:LEU:HD12	2.02	0.42
2:H:89:ILE:H	2:H:89:ILE:HG13	1.55	0.42
1:L:201:LEU:HD13	1:L:205:VAL:HG23	2.00	0.42
2:A:40:ALA:HB3	2:A:43:GLN:HB2	2.00	0.42
1:B:50:LEU:HD12	1:B:53:GLN:HG3	2.01	0.42
1:L:6:GLN:HE21	1:L:6:GLN:HB3	1.73	0.42
1:L:11:LEU:HD13	1:L:11:LEU:HA	1.85	0.42
1:B:34:ALA:O	1:B:88:CYS:HA	2.20	0.42
1:L:59:PRO:HB2	1:L:61:ARG:HG2	2.02	0.41
2:A:198:THR:HG23	2:A:215:LYS:HE3	2.02	0.41
1:L:6:GLN:HG2	1:L:23:CYS:SG	2.60	0.41
1:B:59:PRO:HB2	1:B:61:ARG:HG2	2.02	0.41
1:L:13:VAL:HG21	1:L:19:ALA:HB2	2.02	0.41
2:H:214:LYS:HB2	2:H:214:LYS:HE2	1.83	0.41
1:B:21:MET:HE3	1:B:102:THR:HB	2.03	0.41
2:A:100(A):ARG:NH2	2:A:100(F):TYS:O3	2.45	0.41
2:H:36:TRP:CZ3	2:H:92:CYS:HB3	2.56	0.41
1:L:136:LEU:HD13	1:L:136:LEU:HA	1.71	0.41
2:H:170:THR:HG23	2:H:183:LEU:HD11	2.03	0.41
1:B:91:GLY:HA2	1:B:96:TRP:CD1	2.56	0.41
2:A:4:LEU:HD13	2:A:22:CYS:SG	2.61	0.41
2:H:159:TRP:CZ2	2:H:201:CYS:HB3	2.57	0.40
2:A:64:ARG:O	2:A:82(A):ARG:HG2	2.21	0.40
1:L:82:ASP:O	1:L:86:TYR:OH	2.31	0.40
1:B:33:LEU:HD11	1:B:88:CYS:SG	2.61	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	B	214/216 (99%)	206 (96%)	8 (4%)	0	100	100
1	L	214/216 (99%)	203 (95%)	11 (5%)	0	100	100
2	A	225/238 (94%)	216 (96%)	7 (3%)	2 (1%)	14	41
2	H	225/238 (94%)	216 (96%)	8 (4%)	1 (0%)	30	59
All	All	878/908 (97%)	841 (96%)	34 (4%)	3 (0%)	36	65

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	107	VAL
2	A	25	SER
2	A	107	VAL

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	B	186/186 (100%)	152 (82%)	34 (18%)	2	6
1	L	186/186 (100%)	162 (87%)	24 (13%)	4	13
2	A	195/201 (97%)	173 (89%)	22 (11%)	5	19
2	H	195/201 (97%)	150 (77%)	45 (23%)	1	3
All	All	762/774 (98%)	637 (84%)	125 (16%)	2	8

All (125) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	2	THR
1	L	3	VAL
1	L	6	GLN
1	L	11	LEU
1	L	17	GLU
1	L	23	CYS
1	L	24	THR
1	L	52	SER
1	L	61	ARG
1	L	79	GLU
1	L	88	CYS
1	L	100	LYS
1	L	103	LYS
1	L	104	LEU
1	L	105	GLU
1	L	107	LYS
1	L	117	ILE
1	L	137	ASN
1	L	146	VAL
1	L	161	GLU
1	L	185	ASP
1	L	190	LYS
1	L	197	THR
1	L	199	GLN
2	H	13	LYS
2	H	31	LYS
2	H	45	LEU
2	H	52	SER
2	H	52(A)	HIS
2	H	52(B)	GLU
2	H	55	THR
2	H	59	GLN
2	H	65	VAL
2	H	72	ASP
2	H	73	ASN
2	H	77	THR
2	H	80	ILE
2	H	83	THR
2	H	84	SER
2	H	89	ILE
2	H	97	LYS
2	H	100	LEU
2	H	100(A)	ARG

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Mol	Chain	Res	Type
2	H	100(B)	ASP
2	H	100(C)	TYR
2	H	100(D)	VAL
2	H	100(E)	LEU
2	H	100(G)	ASP
2	H	100(H)	ASP
2	H	100(I)	TYR
2	H	100(K)	LEU
2	H	100(L)	ILE
2	H	100(M)	ASN
2	H	100(N)	GLN
2	H	100(O)	GLN
2	H	100(P)	GLU
2	H	100(Q)	TRP
2	H	100(R)	ASN
2	H	100(S)	ASP
2	H	115	THR
2	H	121	THR
2	H	126	VAL
2	H	140	THR
2	H	143	LEU
2	H	156	THR
2	H	161	SER
2	H	166	SER
2	H	183	LEU
2	H	191	SER
1	B	2	THR
1	B	3	VAL
1	B	5	THR
1	B	6	GLN
1	B	11	LEU
1	B	12	SER
1	B	13	VAL
1	B	20	SER
1	B	21	MET
1	B	23	CYS
1	B	53	GLN
1	B	54	ARG
1	B	56	SER
1	B	60	ASP
1	B	75	ILE
1	B	88	CYS

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Mol	Chain	Res	Type
1	B	94	ARG
1	B	105	GLU
1	B	106	ILE
1	B	107	LYS
1	B	109	THR
1	B	126	LYS
1	B	127	SER
1	B	129	THR
1	B	132	VAL
1	B	136	LEU
1	B	137	ASN
1	B	146	VAL
1	B	161	GLU
1	B	185	ASP
1	B	197	THR
1	B	201	LEU
1	B	202	SER
1	B	203	SER
2	A	4	LEU
2	A	10	GLU
2	A	16	SER
2	A	18	VAL
2	A	28	THR
2	A	29	PHE
2	A	34	VAL
2	A	59	GLN
2	A	65	VAL
2	A	73	ASN
2	A	74	SER
2	A	77	THR
2	A	83	THR
2	A	93	THR
2	A	96	SER
2	A	102	LEU
2	A	110	HIS
2	A	115	THR
2	A	156	THR
2	A	183	LEU
2	A	196	THR
2	A	206	LYS

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	27	GLN
1	L	70	HIS
1	L	199	GLN
2	H	27	ASN
2	H	169	HIS
1	B	27	GLN
1	B	37	GLN
1	B	53	GLN
1	B	189	HIS
2	A	73	ASN
2	A	209	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 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	TYS	A	100(F)	2	15,16,17	1.25	3 (20%)	15,22,24	0.99	1 (6%)
2	TYS	H	100(F)	2	15,16,17	1.21	3 (20%)	15,22,24	1.14	2 (13%)

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	TYS	A	100(F)	2	-	2/10/11/13	0/1/1/1
2	TYS	H	100(F)	2	-	5/10/11/13	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	100(F)	TYS	OH-CZ	-3.06	1.37	1.42
2	H	100(F)	TYS	OH-CZ	-2.93	1.37	1.42
2	A	100(F)	TYS	OH-S	-2.36	1.53	1.58
2	H	100(F)	TYS	O3-S	2.20	1.64	1.50
2	A	100(F)	TYS	O3-S	2.16	1.63	1.50
2	H	100(F)	TYS	OH-S	-2.03	1.54	1.58

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	100(F)	TYS	OH-CZ-CE1	2.81	124.21	118.70
2	A	100(F)	TYS	OH-CZ-CE2	2.30	123.21	118.70
2	H	100(F)	TYS	OH-S-O2	2.09	113.94	107.56

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	H	100(F)	TYS	CZ-OH-S-O3
2	A	100(F)	TYS	CE1-CZ-OH-S
2	A	100(F)	TYS	CE2-CZ-OH-S
2	H	100(F)	TYS	CZ-OH-S-O1
2	H	100(F)	TYS	CZ-OH-S-O2
2	H	100(F)	TYS	CE1-CZ-OH-S
2	H	100(F)	TYS	CE2-CZ-OH-S

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	100(F)	TYS	2	0
2	H	100(F)	TYS	1	0

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	B	216/216 (100%)	0.89	26 (12%) 9 7	18, 43, 61, 69	0
1	L	216/216 (100%)	0.80	18 (8%) 17 14	18, 39, 61, 70	0
2	A	229/238 (96%)	0.86	27 (11%) 9 7	16, 38, 67, 87	0
2	H	229/238 (96%)	1.10	34 (14%) 5 4	14, 46, 80, 96	0
All	All	890/908 (98%)	0.92	105 (11%) 9 7	14, 41, 66, 96	0

All (105) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	A	28	THR	4.7
2	H	84	SER	4.6
2	H	29	PHE	4.5
2	A	27	ASN	4.3
2	A	131	PRO	4.3
2	H	75	ALA	4.3
2	H	76	SER	4.2
2	H	17	SER	4.0
1	B	27	GLN	3.8
1	B	63	SER	3.7
2	H	74	SER	3.7
2	A	29	PHE	3.7
2	A	110	HIS	3.4
1	L	14	THR	3.4
1	B	24	THR	3.4
1	L	23	CYS	3.3
2	A	41	THR	3.3
1	B	62	PHE	3.3
2	H	48	VAL	3.2
2	A	74	SER	3.1
1	B	41	GLY	3.1

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Mol	Chain	Res	Type	RSRZ
1	L	19	ALA	3.1
1	B	3	VAL	3.1
2	H	23	LYS	3.1
2	A	72	ASP	3.0
2	H	18	VAL	3.0
1	L	7	SER	3.0
1	L	76	SER	3.0
2	H	31	LYS	2.9
1	L	58	VAL	2.9
1	B	126	LYS	2.8
2	H	72	ASP	2.8
2	A	191	SER	2.8
2	A	100(J)	GLY	2.8
2	H	119	ALA	2.8
2	A	119	ALA	2.8
1	B	106	ILE	2.8
1	B	43	SER	2.8
1	L	152	ASN	2.7
2	H	82(B)	GLY	2.7
2	A	25	SER	2.7
2	A	130	ALA	2.7
1	L	5	THR	2.7
2	A	42	GLY	2.7
2	H	100	LEU	2.7
1	L	65	SER	2.7
2	A	73	ASN	2.6
2	H	4	LEU	2.6
2	H	79	TYR	2.6
1	B	182	SER	2.5
1	L	41	GLY	2.5
2	H	61	PHE	2.5
1	B	2	THR	2.5
2	H	24	ALA	2.5
2	A	40	ALA	2.5
2	A	75	ALA	2.5
2	H	73	ASN	2.5
1	B	184	ALA	2.4
1	L	2	THR	2.4
2	H	140	THR	2.4
2	H	34	VAL	2.4
2	A	218	PRO	2.4
2	A	26	GLY	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	76	SER	2.4
2	H	96	SER	2.4
1	B	145	LYS	2.3
2	H	5	VAL	2.3
2	A	104	PHE	2.3
1	B	56	SER	2.3
1	B	66	GLY	2.3
1	B	152	ASN	2.3
2	A	98	HIS	2.3
1	L	77	ARG	2.2
2	A	107	VAL	2.2
2	A	100(I)	TYR	2.2
1	B	73	LEU	2.2
2	A	33	ASP	2.2
1	L	24	THR	2.2
2	H	109	GLY	2.2
2	H	97	LYS	2.2
1	L	18	ALA	2.2
1	L	27	GLN	2.2
2	H	27	ASN	2.2
2	H	41	THR	2.2
1	B	25	SER	2.2
1	B	67	SER	2.2
1	B	60	ASP	2.1
2	A	32	TYR	2.1
1	L	27(A)	SER	2.1
1	B	23	CYS	2.1
2	H	30	ARG	2.1
2	H	110	HIS	2.1
2	H	218	PRO	2.1
2	H	26	GLY	2.1
1	B	78	VAL	2.1
1	B	52	SER	2.1
2	A	100(M)	ASN	2.1
2	A	140	THR	2.1
2	H	196	THR	2.0
1	B	80	ALA	2.0
1	L	203	SER	2.0
2	H	25	SER	2.0
1	L	104	LEU	2.0
1	B	129	THR	2.0
2	H	19	LYS	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($\text{\AA}^2$)	Q<0.9
2	TYS	H	100(F)	16/17	0.83	0.15	39,50,64,64	0
2	TYS	A	100(F)	16/17	0.88	0.13	47,52,58,58	4

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.