



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 10:52 PM UTC

PDB ID : 2XZQ / pdb_00002xzq
Title : CRYSTAL STRUCTURE ANALYSIS OF THE ANTI-(4-HYDROXY-3-NITROPHENYL)- ACETYL MURINE GERMLINE MONOCLONAL ANTIBODY BBE6.12H3 FAB FRAGMENT IN COMPLEX WITH A PHAGE DISPLAY DERIVED DODECAPEPTIDE YQLRPNAETLRF
Authors : Khan, T.; Salunke, D.M.
Deposited on : 2010-12-01
Resolution : 2.40 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

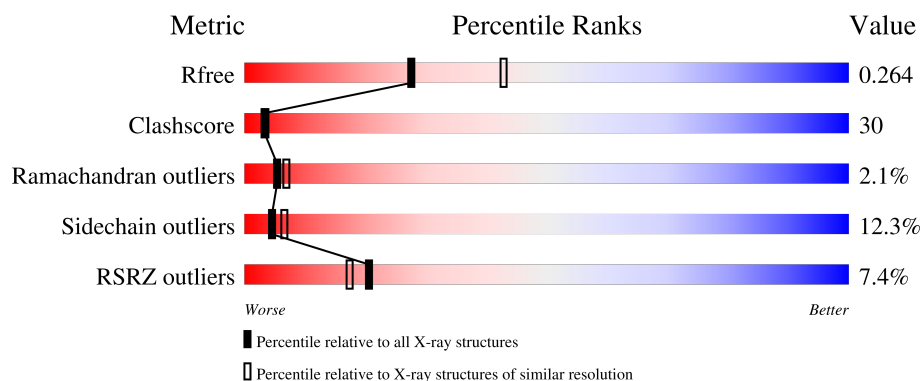
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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 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	H	220	 6% 52% 36% 8% ..
2	L	211	 4% 56% 38% 5%
3	P	12	 75% 8% 25% 8% 33% 25%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 3434 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 ANTI-NP MURINE GERMLINE MONOCLONAL ANTI-BODY BBE6.12H3, HEAVY CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	215	Total	C	N	O	S	0	0	0
			1635	1043	268	317	7			

- Molecule 2 is a protein called ANTI-NP MURINE GERMLINE MONOCLONAL ANTI-BODY BBE6.12H3, LIGHT CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	211	Total	C	N	O	S	0	0	0
			1598	996	270	326	6			

- Molecule 3 is a protein called PHAGE DISPLAY DERIVED ANTIGEN.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	P	9	Total	C	N	O	0	0	0
			73	44	14	15			

- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	H	1	Total	Zn	0	0
			1	1		
4	P	1	Total	Zn	0	0
			1	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	H	61	Total	O	0	0
			61	61		
5	L	63	Total	O	0	0
			63	63		

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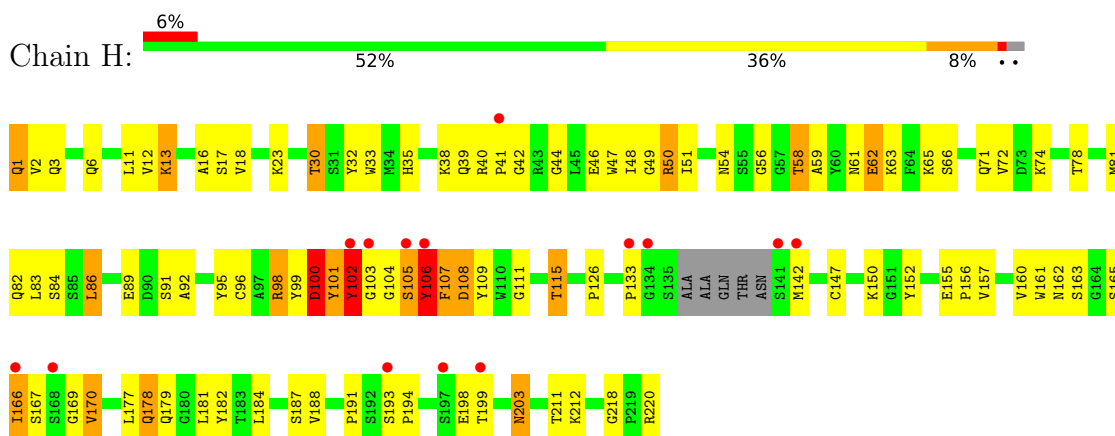
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	P	2	Total	O	0	0
			2	2		

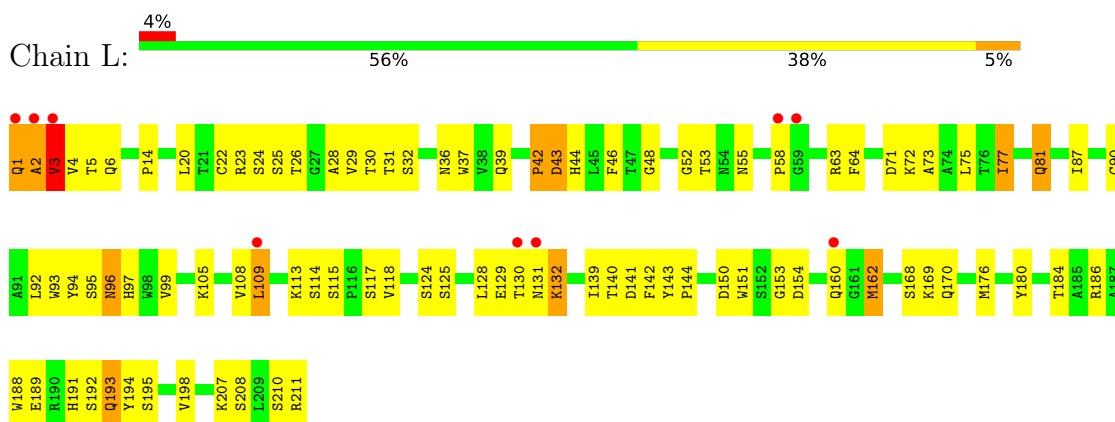
3 Residue-property plots

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.

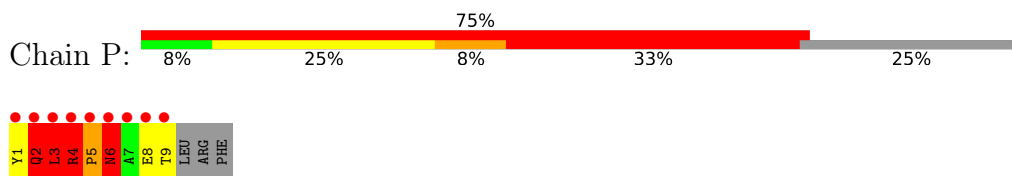
- Molecule 1: ANTI-NP MURINE GERMLINE MONOCLONAL ANTIBODY BBE6.12H3, HEAVY CHAIN



- Molecule 2: ANTI-NP MURINE GERMLINE MONOCLONAL ANTIBODY BBE6.12H3, LIGHT CHAIN



- Molecule 3: PHAGE DISPLAY DERIVED ANTIGEN



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	55.92Å 65.28Å 113.56Å 90.00° 90.19° 90.00°	Depositor
Resolution (Å)	50.00 – 2.40 50.00 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.4 (50.00-2.40) 99.5 (50.00-2.40)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.50 (at 2.39Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.241 , 0.251 0.239 , 0.264	Depositor DCC
R_{free} test set	1626 reflections (10.07%)	wwPDB-VP
Wilson B-factor (Å ²)	26.9	Xtriage
Anisotropy	0.928	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 37.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.023 for -h,-k,l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	3434	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.22% 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: ZN

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	H	0.49	0/1684	1.17	19/2300 (0.8%)
2	L	0.51	1/1636 (0.1%)	1.13	11/2234 (0.5%)
3	P	0.98	0/74	1.75	2/100 (2.0%)
All	All	0.52	1/3394 (0.0%)	1.17	32/4634 (0.7%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	L	3	VAL	C-O	-5.36	1.17	1.24

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	130	THR	N-CA-C	-14.54	88.96	110.48
2	L	2	ALA	N-CA-C	-14.39	95.01	113.72
2	L	3	VAL	N-CA-C	11.27	132.78	109.34
1	H	142	MET	N-CA-C	10.35	126.95	113.18
1	H	103	GLY	N-CA-C	-10.10	99.19	115.66
1	H	199	THR	N-CA-C	9.30	124.04	110.28
3	P	3	LEU	N-CA-C	9.29	130.60	110.80
1	H	106	TYR	N-CA-C	8.92	123.56	109.65
1	H	169	GLY	N-CA-C	-7.20	105.85	115.21
2	L	2	ALA	CA-C-N	6.94	134.46	121.97
2	L	2	ALA	C-N-CA	6.94	134.46	121.97
1	H	101	TYR	N-CA-C	6.89	123.69	114.12
1	H	111	GLY	N-CA-C	-6.75	101.72	112.66
3	P	4	ARG	N-CA-C	6.67	124.54	109.81
1	H	155	GLU	N-CA-C	-6.59	100.51	110.39
1	H	218	GLY	CA-C-N	6.53	126.92	119.93
1	H	218	GLY	C-N-CA	6.53	126.92	119.93

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	96	CYS	N-CA-C	-5.94	99.22	108.90
1	H	30	THR	N-CA-C	5.91	120.10	113.01
2	L	153	GLY	N-CA-C	-5.90	99.19	113.18
1	H	44	GLY	N-CA-C	5.63	124.33	112.17
1	H	150	LYS	N-CA-C	5.58	118.90	109.76
2	L	130	THR	CA-C-N	-5.54	113.85	122.16
2	L	130	THR	C-N-CA	-5.54	113.85	122.16
1	H	61	ASN	N-CA-C	-5.46	101.77	109.96
2	L	184	THR	N-CA-C	-5.41	101.72	110.32
1	H	56	GLY	N-CA-C	-5.23	108.50	115.40
1	H	100	ASP	CB-CA-C	5.14	118.11	109.48
1	H	106	TYR	CB-CA-C	-5.13	101.83	112.44
2	L	24	SER	N-CA-C	-5.07	101.42	109.59
1	H	107	PHE	N-CA-C	-5.05	99.06	107.80
2	L	109	LEU	N-CA-C	-5.04	102.74	110.30

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	H	1635	0	1588	98	1
2	L	1598	0	1531	95	1
3	P	73	0	65	32	0
4	H	1	0	0	0	0
4	P	1	0	0	0	0
5	H	61	0	0	7	0
5	L	63	0	0	4	0
5	P	2	0	0	0	0
All	All	3434	0	3184	197	1

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 30.

All (197) 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:62:GLU:HA	1:H:65:LYS:HE3	1.26	1.13
1:H:102:TYR:CD1	1:H:102:TYR:N	2.28	1.01
2:L:14:PRO:HD3	2:L:109:LEU:O	1.67	0.95
2:L:94:TYR:HB3	3:P:2:GLN:HB2	1.49	0.93
2:L:1:GLN:HG3	2:L:99:VAL:HG13	1.50	0.93
3:P:4:ARG:HB2	3:P:5:PRO:CD	1.99	0.92
2:L:32:SER:HB2	3:P:4:ARG:HA	1.50	0.91
1:H:106:TYR:OH	2:L:55:ASN:HB3	1.71	0.91
2:L:131:ASN:O	2:L:132:LYS:HD3	1.74	0.88
2:L:170:GLN:HE21	2:L:176:MET:HB3	1.42	0.84
1:H:59:ALA:HB1	3:P:8:GLU:OE2	1.78	0.84
1:H:193:SER:OG	1:H:194:PRO:HD3	1.77	0.82
1:H:62:GLU:HA	1:H:65:LYS:CE	2.09	0.81
3:P:5:PRO:O	3:P:6:ASN:HB2	1.78	0.81
1:H:58:THR:HG22	5:H:2014:HOH:O	1.80	0.80
1:H:102:TYR:C	1:H:104:GLY:H	1.88	0.80
2:L:94:TYR:HB3	3:P:2:GLN:CB	2.12	0.78
1:H:102:TYR:N	1:H:102:TYR:HD1	1.80	0.78
1:H:81:MET:HE3	1:H:83:LEU:HD21	1.64	0.78
2:L:141:ASP:H	2:L:170:GLN:HE22	1.30	0.77
1:H:162:ASN:HB2	1:H:166:ILE:HD13	1.67	0.77
1:H:38:LYS:HE2	1:H:40:ARG:HD2	1.65	0.76
1:H:178:GLN:O	1:H:179:GLN:HG2	1.85	0.76
1:H:13:LYS:H	1:H:13:LYS:HE3	1.50	0.76
1:H:163:SER:H	1:H:203:ASN:HD21	1.33	0.75
1:H:32:TYR:CD2	1:H:98:ARG:HD3	2.20	0.75
1:H:13:LYS:HE3	1:H:13:LYS:N	2.02	0.75
2:L:94:TYR:HB3	3:P:2:GLN:CG	2.18	0.74
1:H:17:SER:HA	1:H:86:LEU:HD22	1.69	0.73
5:H:2041:HOH:O	2:L:44:HIS:HB2	1.89	0.72
1:H:107:PHE:O	2:L:48:GLY:HA3	1.90	0.71
2:L:26:THR:HG21	3:P:1:TYR:N	2.06	0.71
2:L:118:VAL:HG22	2:L:139:ILE:HD12	1.72	0.71
1:H:71:GLN:HG3	1:H:72:VAL:H	1.57	0.70
3:P:4:ARG:HB2	3:P:5:PRO:HD3	1.73	0.70
2:L:94:TYR:HA	3:P:3:LEU:H	1.57	0.70
2:L:1:GLN:CG	2:L:99:VAL:HG13	2.21	0.69
1:H:6:GLN:HE22	1:H:95:TYR:HA	1.55	0.69
1:H:170:VAL:HA	1:H:187:SER:O	1.93	0.69
2:L:125:SER:O	2:L:129:GLU:HG3	1.93	0.69
1:H:50:ARG:HG2	1:H:51:ILE:N	2.08	0.69
2:L:93:TRP:HD1	3:P:4:ARG:O	1.76	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:166:ILE:HD12	1:H:166:ILE:N	2.07	0.68
2:L:208:SER:HB2	5:L:2058:HOH:O	1.93	0.68
2:L:2:ALA:HB3	2:L:99:VAL:HG21	1.75	0.68
2:L:5:THR:HG23	2:L:23:ARG:HB2	1.75	0.67
2:L:14:PRO:CD	2:L:109:LEU:O	2.44	0.65
2:L:140:THR:HB	2:L:170:GLN:HE22	1.61	0.65
2:L:3:VAL:HG23	2:L:25:SER:OG	1.97	0.64
2:L:26:THR:HG21	3:P:1:TYR:H2	1.61	0.64
2:L:141:ASP:H	2:L:170:GLN:NE2	1.95	0.64
2:L:5:THR:CG2	2:L:23:ARG:HB2	2.27	0.64
1:H:102:TYR:C	1:H:104:GLY:N	2.53	0.64
2:L:26:THR:CG2	3:P:1:TYR:H1	2.11	0.64
2:L:139:ILE:HD12	2:L:198:VAL:HG21	1.80	0.64
2:L:140:THR:HB	2:L:170:GLN:NE2	2.14	0.63
1:H:106:TYR:HH	2:L:55:ASN:HB3	1.64	0.63
2:L:2:ALA:CB	2:L:99:VAL:HG21	2.29	0.63
2:L:94:TYR:CA	3:P:3:LEU:H	2.11	0.63
2:L:151:TRP:CD1	2:L:162:MET:HG3	2.34	0.63
2:L:93:TRP:CD1	3:P:4:ARG:O	2.52	0.62
1:H:105:SER:HB3	5:H:2039:HOH:O	1.98	0.62
1:H:71:GLN:HG3	1:H:72:VAL:N	2.14	0.62
1:H:40:ARG:O	1:H:41:PRO:C	2.39	0.62
1:H:12:VAL:HG21	1:H:18:VAL:CG1	2.30	0.62
1:H:100:ASP:OD1	1:H:106:TYR:CD1	2.52	0.62
2:L:96:ASN:C	2:L:96:ASN:HD22	2.06	0.61
2:L:1:GLN:H2	2:L:3:VAL:HG12	1.65	0.61
2:L:6:GLN:NE2	2:L:90:CYS:H	2.00	0.60
1:H:66:SER:HB2	5:H:2021:HOH:O	2.01	0.60
1:H:191:PRO:HB2	1:H:194:PRO:CD	2.32	0.59
2:L:192:SER:O	2:L:210:SER:HA	2.01	0.59
1:H:98:ARG:HD2	1:H:108:ASP:OD1	2.02	0.59
2:L:3:VAL:HG23	2:L:25:SER:CB	2.33	0.59
2:L:195:SER:HB2	5:L:2058:HOH:O	2.02	0.59
2:L:81:GLN:O	2:L:108:VAL:HG11	2.03	0.59
1:H:133:PRO:O	1:H:220:ARG:OXT	2.20	0.58
1:H:191:PRO:HB2	1:H:194:PRO:HD2	1.85	0.58
1:H:194:PRO:O	1:H:198:GLU:N	2.32	0.58
1:H:38:LYS:CB	1:H:48:ILE:HD11	2.33	0.57
2:L:42:PRO:HG2	2:L:169:LYS:HD2	1.86	0.57
1:H:17:SER:HB3	1:H:84:SER:HA	1.86	0.57
2:L:26:THR:CG2	3:P:1:TYR:N	2.68	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:104:GLY:O	1:H:105:SER:CB	2.51	0.57
1:H:191:PRO:O	1:H:194:PRO:HD2	2.03	0.57
2:L:1:GLN:H2	2:L:3:VAL:CG1	2.18	0.56
1:H:99:TYR:CD2	1:H:106:TYR:O	2.59	0.56
2:L:94:TYR:HA	3:P:3:LEU:HA	1.86	0.56
1:H:12:VAL:CG1	1:H:16:ALA:HB3	2.36	0.55
2:L:193:GLN:C	2:L:193:GLN:HE21	2.14	0.55
1:H:160:VAL:HG12	1:H:161:TRP:N	2.22	0.55
1:H:193:SER:HG	1:H:194:PRO:HD3	1.72	0.55
1:H:178:GLN:O	1:H:179:GLN:CG	2.54	0.55
1:H:13:LYS:H	1:H:13:LYS:CE	2.19	0.55
1:H:47:TRP:CZ2	1:H:49:GLY:HA2	2.42	0.55
5:L:2029:HOH:O	3:P:2:GLN:HB3	2.07	0.55
2:L:81:GLN:HE21	2:L:81:GLN:HA	1.72	0.54
1:H:194:PRO:O	1:H:198:GLU:HG2	2.07	0.54
2:L:64:PHE:CE1	2:L:77:ILE:HD11	2.42	0.54
2:L:94:TYR:CB	3:P:2:GLN:HB2	2.31	0.54
1:H:220:ARG:HH12	2:L:124:SER:HA	1.72	0.54
2:L:77:ILE:HG22	2:L:77:ILE:O	2.08	0.54
2:L:94:TYR:HA	3:P:3:LEU:N	2.22	0.54
1:H:40:ARG:O	1:H:42:GLY:N	2.40	0.54
1:H:39:GLN:O	1:H:92:ALA:HB1	2.07	0.54
1:H:162:ASN:CB	1:H:166:ILE:HD13	2.36	0.53
1:H:178:GLN:O	1:H:178:GLN:HG3	2.08	0.53
2:L:139:ILE:CD1	2:L:198:VAL:HG21	2.39	0.53
2:L:42:PRO:O	2:L:43:ASP:HB2	2.09	0.53
1:H:162:ASN:HB2	1:H:166:ILE:CD1	2.37	0.53
2:L:94:TYR:HB3	3:P:2:GLN:HG3	1.90	0.53
1:H:106:TYR:CD1	1:H:106:TYR:N	2.76	0.52
1:H:166:ILE:N	1:H:166:ILE:CD1	2.73	0.52
3:P:4:ARG:CB	3:P:5:PRO:HD3	2.40	0.52
1:H:62:GLU:CA	1:H:65:LYS:HE3	2.18	0.52
2:L:3:VAL:HG23	2:L:25:SER:HB3	1.92	0.52
2:L:4:VAL:HA	2:L:23:ARG:O	2.10	0.52
1:H:166:ILE:HG21	1:H:188:VAL:HG21	1.92	0.52
1:H:115:THR:HG23	5:H:2042:HOH:O	2.10	0.51
2:L:154:ASP:H	2:L:193:GLN:HE21	1.58	0.51
1:H:38:LYS:HB3	1:H:48:ILE:HD11	1.93	0.51
2:L:154:ASP:N	2:L:193:GLN:NE2	2.59	0.50
1:H:91:SER:O	1:H:92:ALA:HB2	2.11	0.50
2:L:94:TYR:C	3:P:2:GLN:HB2	2.36	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:162:ASN:ND2	1:H:166:ILE:HD13	2.27	0.50
2:L:96:ASN:C	2:L:96:ASN:ND2	2.70	0.49
1:H:104:GLY:O	1:H:105:SER:HB3	2.12	0.49
1:H:160:VAL:CG1	1:H:161:TRP:N	2.75	0.49
2:L:64:PHE:HE1	2:L:77:ILE:HD11	1.76	0.49
2:L:162:MET:HA	2:L:180:TYR:O	2.12	0.49
2:L:37:TRP:CD2	2:L:75:LEU:HB2	2.48	0.49
2:L:207:LYS:HD3	2:L:207:LYS:HA	1.66	0.48
2:L:63:ARG:CZ	2:L:81:GLN:HG3	2.44	0.48
1:H:18:VAL:O	1:H:82:GLN:HA	2.14	0.48
2:L:28:ALA:HA	2:L:71:ASP:HB2	1.96	0.48
2:L:6:GLN:HE22	2:L:90:CYS:H	1.60	0.47
1:H:2:VAL:HG11	1:H:109:TYR:CD2	2.49	0.47
1:H:12:VAL:HG21	1:H:18:VAL:HG13	1.95	0.47
1:H:33:TRP:CE3	1:H:50:ARG:HD2	2.49	0.47
1:H:71:GLN:CG	1:H:72:VAL:H	2.24	0.47
1:H:38:LYS:HB2	1:H:48:ILE:HD11	1.95	0.47
2:L:39:GLN:O	2:L:46:PHE:HA	2.15	0.47
3:P:4:ARG:HB2	3:P:5:PRO:HD2	1.93	0.47
1:H:106:TYR:N	1:H:106:TYR:HD1	2.13	0.47
1:H:184:LEU:HD12	1:H:184:LEU:C	2.38	0.47
1:H:12:VAL:HG13	1:H:16:ALA:HB3	1.96	0.46
1:H:17:SER:CB	1:H:84:SER:HA	2.44	0.46
3:P:2:GLN:O	3:P:3:LEU:CB	2.64	0.46
3:P:5:PRO:C	3:P:9:THR:HG22	2.41	0.46
1:H:211:THR:HG22	1:H:212:LYS:N	2.31	0.46
2:L:95:SER:H	3:P:3:LEU:C	2.22	0.46
1:H:106:TYR:HA	2:L:36:ASN:HD21	1.81	0.46
2:L:93:TRP:CZ2	3:P:9:THR:HA	2.52	0.45
2:L:23:ARG:HG2	2:L:72:LYS:HG2	1.97	0.45
1:H:163:SER:H	1:H:203:ASN:ND2	2.09	0.45
2:L:20:LEU:HD12	2:L:20:LEU:N	2.30	0.45
2:L:125:SER:HA	2:L:128:LEU:HB2	1.99	0.45
1:H:126:PRO:HB3	1:H:152:TYR:HB3	1.99	0.45
1:H:165:SER:C	1:H:166:ILE:HD12	2.41	0.45
2:L:113:LYS:HG3	2:L:144:PRO:HD3	1.99	0.45
2:L:4:VAL:CG1	2:L:22:CYS:SG	3.04	0.44
1:H:165:SER:HB2	5:H:2054:HOH:O	2.17	0.44
2:L:154:ASP:HA	2:L:193:GLN:HB3	2.00	0.44
1:H:35:HIS:CE1	1:H:50:ARG:HB2	2.53	0.44
2:L:191:HIS:HB2	2:L:194:TYR:OH	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:114:SER:O	2:L:142:PHE:HA	2.18	0.44
1:H:162:ASN:CG	1:H:166:ILE:HD13	2.44	0.43
1:H:18:VAL:HG22	1:H:86:LEU:HD21	2.00	0.43
2:L:22:CYS:HB3	2:L:73:ALA:HB3	2.01	0.43
1:H:62:GLU:OE1	1:H:65:LYS:NZ	2.51	0.43
1:H:12:VAL:CG1	1:H:13:LYS:N	2.82	0.43
2:L:87:ILE:HG12	2:L:105:LYS:HD3	2.00	0.42
2:L:52:GLY:O	2:L:53:THR:HB	2.19	0.42
2:L:96:ASN:ND2	2:L:97:HIS:ND1	2.67	0.42
1:H:165:SER:HB3	1:H:166:ILE:HD12	2.00	0.42
2:L:29:VAL:HG11	2:L:73:ALA:HB2	2.00	0.42
2:L:131:ASN:C	2:L:132:LYS:HD3	2.44	0.42
1:H:1:GLN:NE2	5:H:2001:HOH:O	2.52	0.42
1:H:46:GLU:HG2	2:L:1:GLN:HE22	1.85	0.42
1:H:12:VAL:HG11	1:H:86:LEU:HD23	2.02	0.42
1:H:30:THR:O	1:H:54:ASN:HB2	2.20	0.42
1:H:63:LYS:HG2	1:H:63:LYS:O	2.20	0.42
2:L:189:GLU:O	2:L:211:ARG:NH2	2.53	0.42
3:P:5:PRO:CA	3:P:9:THR:HG22	2.49	0.41
2:L:43:ASP:HA	5:L:2011:HOH:O	2.20	0.41
2:L:94:TYR:HA	3:P:3:LEU:CA	2.49	0.41
1:H:99:TYR:CG	1:H:106:TYR:O	2.74	0.41
2:L:23:ARG:CG	2:L:72:LYS:HD3	2.50	0.41
1:H:152:TYR:CE2	1:H:157:VAL:HG13	2.56	0.41
2:L:26:THR:HG22	3:P:1:TYR:H1	1.85	0.41
2:L:143:TYR:HA	2:L:144:PRO:C	2.46	0.41
3:P:5:PRO:HA	3:P:9:THR:HG22	2.02	0.41
1:H:101:TYR:CD1	1:H:102:TYR:HD1	2.38	0.41
1:H:177:LEU:HD13	1:H:182:TYR:CE1	2.55	0.41
2:L:96:ASN:HD22	2:L:97:HIS:N	2.17	0.41
1:H:51:ILE:CD1	1:H:71:GLN:HA	2.51	0.41
1:H:71:GLN:CG	1:H:72:VAL:N	2.81	0.40

All (1) 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)
1:H:3:GLN:NE2	2:L:150:ASP:OD2[3_445]	2.18	0.02

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	H	211/220 (96%)	196 (93%)	13 (6%)	2 (1%)	14	22
2	L	209/211 (99%)	198 (95%)	9 (4%)	2 (1%)	12	20
3	P	7/12 (58%)	2 (29%)	0	5 (71%)	0	0
All	All	427/443 (96%)	396 (93%)	22 (5%)	9 (2%)	5	7

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	102	TYR
1	H	105	SER
3	P	2	GLN
3	P	3	LEU
3	P	4	ARG
3	P	5	PRO
3	P	6	ASN
2	L	43	ASP
2	L	3	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	H	182/185 (98%)	157 (86%)	25 (14%)	3	5
2	L	178/178 (100%)	160 (90%)	18 (10%)	7	11
3	P	7/11 (64%)	5 (71%)	2 (29%)	0	0

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	367/374 (98%)	322 (88%)	45 (12%)	4 6

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	1	GLN
1	H	11	LEU
1	H	13	LYS
1	H	23	LYS
1	H	50	ARG
1	H	58	THR
1	H	62	GLU
1	H	74	LYS
1	H	78	THR
1	H	86	LEU
1	H	89	GLU
1	H	98	ARG
1	H	100	ASP
1	H	102	TYR
1	H	106	TYR
1	H	108	ASP
1	H	115	THR
1	H	147	CYS
1	H	156	PRO
1	H	166	ILE
1	H	167	SER
1	H	170	VAL
1	H	178	GLN
1	H	181	LEU
1	H	203	ASN
2	L	1	GLN
2	L	30	THR
2	L	31	THR
2	L	42	PRO
2	L	58	PRO
2	L	77	ILE
2	L	81	GLN
2	L	92	LEU
2	L	96	ASN
2	L	115	SER
2	L	117	SER
2	L	132	LYS

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Mol	Chain	Res	Type
2	L	160	GLN
2	L	162	MET
2	L	168	SER
2	L	186	ARG
2	L	188	TRP
2	L	193	GLN
3	P	2	GLN
3	P	6	ASN

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

Mol	Chain	Res	Type
1	H	1	GLN
1	H	6	GLN
1	H	54	ASN
1	H	162	ASN
1	H	171	HIS
1	H	203	ASN
2	L	1	GLN
2	L	6	GLN
2	L	36	ASN
2	L	44	HIS
2	L	81	GLN
2	L	96	ASN
2	L	170	GLN
2	L	191	HIS
2	L	193	GLN
2	L	200	HIS
3	P	6	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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 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 [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	H	215/220 (97%)	0.40	14 (6%) 25 21	16, 27, 48, 56	0
2	L	211/211 (100%)	0.22	9 (4%) 40 36	13, 23, 41, 48	0
3	P	9/12 (75%)	3.97	9 (100%) 0 0	59, 66, 72, 72	0
All	All	435/443 (98%)	0.39	32 (7%) 20 17	13, 25, 47, 72	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	P	7	ALA	6.1
2	L	1	GLN	5.4
3	P	5	PRO	5.0
3	P	1	TYR	4.6
1	H	105	SER	4.4
3	P	9	THR	4.1
3	P	3	LEU	4.1
2	L	2	ALA	3.9
3	P	6	ASN	3.6
1	H	166	ILE	3.6
1	H	41	PRO	3.5
1	H	141	SER	3.5
1	H	134	GLY	3.1
1	H	106	TYR	3.0
3	P	8	GLU	3.0
3	P	2	GLN	2.9
1	H	103	GLY	2.8
1	H	102	TYR	2.7
2	L	3	VAL	2.7
2	L	160	GLN	2.6
1	H	133	PRO	2.6
1	H	199	THR	2.6
2	L	130	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	H	168	SER	2.4
3	P	4	ARG	2.4
1	H	142	MET	2.4
2	L	58	PRO	2.3
2	L	131	ASN	2.3
2	L	109	LEU	2.3
1	H	193	SER	2.2
1	H	197	SER	2.2
2	L	59	GLY	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($\text{\AA}^2$)	Q<0.9
4	ZN	H	1221	1/1	0.99	0.05	31,31,31,31	0
4	ZN	P	1221	1/1	0.99	0.13	17,17,17,17	0

6.5 Other polymers [i](#)

There are no such residues in this entry.