



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 03:06 AM UTC

PDB ID : 6A8O / pdb_00006a8o
Title : Crystal structures of the serine protease domain of murine plasma kallikrein with peptide inhibitor mupain-1-16
Authors : Xu, M.; Jiang, L.; Huang, M.
Deposited on : 2018-07-09
Resolution : 2.77 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

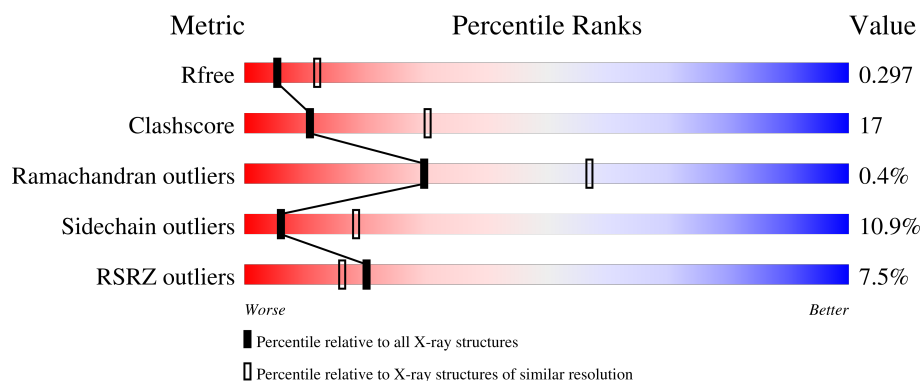
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.77 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	5248 (2.80-2.76)
Clashscore	190562	5693 (2.80-2.76)
Ramachandran outliers	187476	5590 (2.80-2.76)
Sidechain outliers	187428	5592 (2.80-2.76)
RSRZ outliers	180081	5251 (2.80-2.76)

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	248	
1	B	248	
2	P	10	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3821 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 Plasma kallikrein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	236	Total	C	N	O	S	0	0	0
			1858	1184	312	352	10			
1	B	236	Total	C	N	O	S	0	0	0
			1858	1184	312	352	10			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	122	SER	CYS	engineered mutation	UNP P26262
B	122	SER	CYS	engineered mutation	UNP P26262

- Molecule 2 is a protein called peptide inhibitor,.

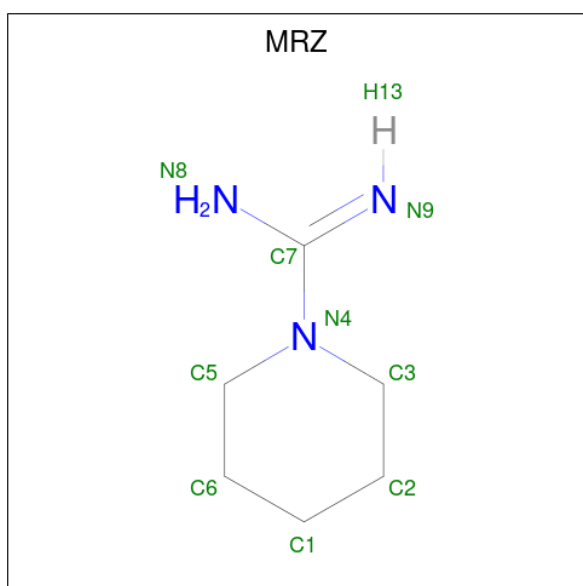
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	P	10	Total	C	N	O	S	0	0	0
			74	48	10	14	2			

- Molecule 3 is alpha-D-mannopyranose (CCD ID: MAN) (formula: C₆H₁₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			11	6	5		
3	B	1	Total	C	O	0	0
			11	6	5		

- Molecule 4 is piperidine-1-carboximidamide (CCD ID: MRZ) (formula: $C_6H_{13}N_3$).

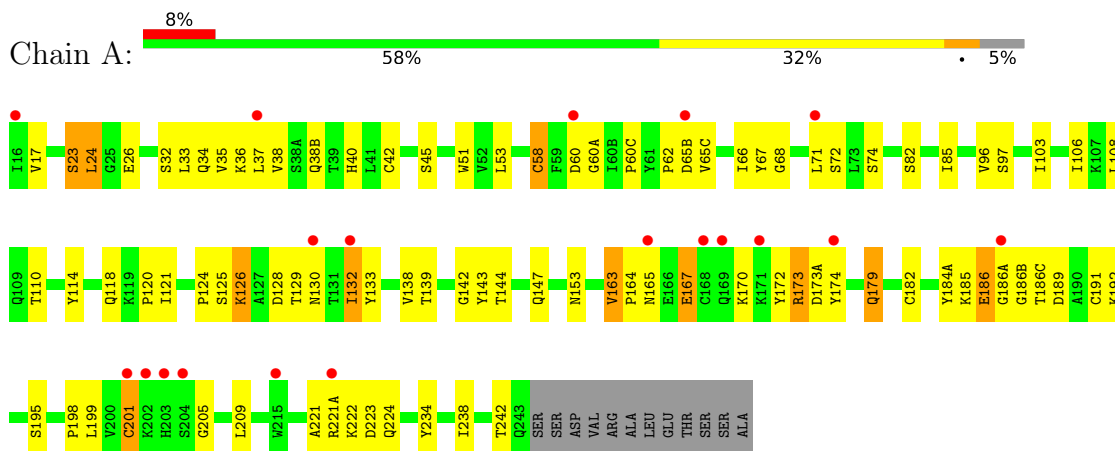


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	P	1	Total	C	N	0	0
			9	6	3		

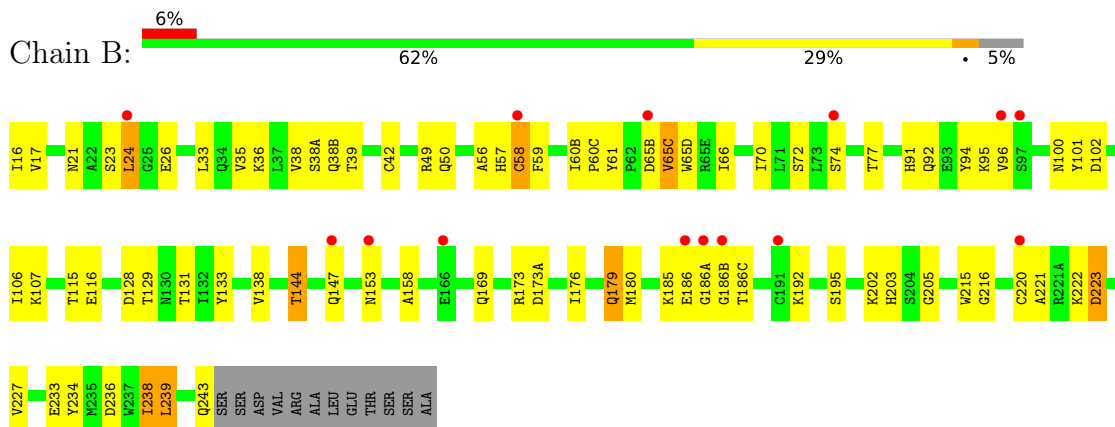
3 Residue-property plots

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.

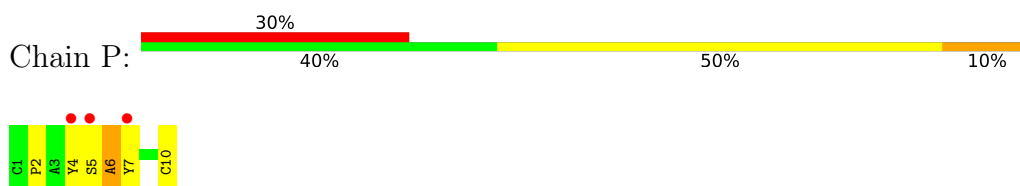
- Molecule 1: Plasma kallikrein



- Molecule 1: Plasma kallikrein



- Molecule 2: peptide inhibitor,



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	71.26Å 82.45Å 105.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	52.66 – 2.77 52.66 – 2.77	Depositor EDS
% Data completeness (in resolution range)	99.8 (52.66-2.77) 99.8 (52.66-2.77)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.43 (at 2.77Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.212 , 0.300 0.234 , 0.297	Depositor DCC
R_{free} test set	814 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	70.7	Xtriage
Anisotropy	0.027	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 48.5	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	0.93	EDS
Total number of atoms	3821	wwPDB-VP
Average B, all atoms (Å ²)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.25% 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: MRZ, MAN

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	1.05	2/1904 (0.1%)	1.42	10/2587 (0.4%)
1	B	1.04	2/1904 (0.1%)	1.40	3/2587 (0.1%)
2	P	2.70	5/76 (6.6%)	2.13	5/103 (4.9%)
All	All	1.10	9/3884 (0.2%)	1.43	18/5277 (0.3%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	P	6	ALA	C-O	-19.29	0.95	1.24
1	A	60(C)	PRO	C-O	-9.19	1.13	1.23
1	B	65(B)	ASP	CA-C	-6.35	1.44	1.52
1	B	65(C)	VAL	N-CA	6.26	1.54	1.46
2	P	6	ALA	CA-C	5.66	1.62	1.52
2	P	7	TYR	C-N	-5.50	1.26	1.33
2	P	4	TYR	C-N	5.43	1.41	1.33
1	A	65(B)	ASP	C-O	-5.29	1.16	1.24
2	P	5	SER	C-N	5.11	1.41	1.33

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	P	6	ALA	N-CA-C	10.65	125.86	113.19
1	A	38	VAL	N-CA-C	-8.41	104.33	111.56
2	P	6	ALA	CB-CA-C	-7.88	94.99	110.27
1	A	167	GLU	N-CA-C	-7.17	101.71	112.04
1	A	62	PRO	N-CA-C	6.75	122.45	113.40
1	A	130	ASN	CB-CA-C	5.97	115.75	109.83
2	P	6	ALA	CA-C-N	5.87	131.40	122.65
2	P	6	ALA	C-N-CA	5.87	131.40	122.65
1	A	165	ASN	CA-C-N	5.83	132.68	121.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	165	ASN	C-N-CA	5.83	132.68	121.54
1	A	110	THR	CB-CA-C	5.65	116.84	108.87
1	B	144	THR	CA-CB-OG1	-5.61	101.19	109.60
2	P	6	ALA	N-CA-CB	5.25	119.66	110.42
1	A	38	VAL	CB-CA-C	5.23	116.41	111.44
1	A	65(B)	ASP	CB-CA-C	5.23	120.03	110.36
1	A	62	PRO	CB-CA-C	-5.20	104.75	112.55
1	B	65(C)	VAL	N-CA-C	-5.16	104.52	111.44
1	B	205	GLY	O-C-N	5.12	128.94	122.43

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	1858	0	1827	67	0
1	B	1858	0	1829	61	0
2	P	74	0	65	1	0
3	A	11	0	10	1	0
3	B	11	0	10	2	0
4	P	9	0	11	2	0
All	All	3821	0	3752	126	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 17.

All (126) 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:B:21:ASN:HD22	3:B:301:MAN:C1	1.06	1.55
1:B:21:ASN:ND2	3:B:301:MAN:C1	1.81	1.40
1:B:42:CYS:CB	1:B:58:CYS:SG	2.30	1.20
1:A:184(A):TYR:HB2	1:A:186(A):GLY:HA3	1.16	1.14
1:A:184(A):TYR:CB	1:A:186(A):GLY:HA3	1.84	1.05

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:42:CYS:CB	1:A:58:CYS:SG	2.45	1.05
1:A:42:CYS:SG	1:A:58:CYS:SG	1.19	1.03
1:B:42:CYS:SG	1:B:58:CYS:SG	1.06	1.00
1:A:186(B):GLY:HA2	1:A:221:ALA:HB1	1.46	0.97
1:B:17:VAL:HG12	1:B:186(B):GLY:O	1.71	0.91
1:A:126:LYS:O	1:A:129:THR:HG23	1.70	0.88
1:A:186:GLU:HG2	1:A:222:LYS:HE3	1.57	0.86
1:A:186(B):GLY:CA	1:A:221:ALA:HB1	2.06	0.86
1:A:24:LEU:HD22	1:A:71:LEU:HD23	1.56	0.85
1:B:179:GLN:HB3	1:B:233:GLU:HG2	1.56	0.85
1:B:42:CYS:SG	1:B:58:CYS:CB	2.64	0.85
1:A:32:SER:HB2	1:A:40:HIS:HD2	1.42	0.84
1:A:184(A):TYR:HB2	1:A:186(A):GLY:CA	2.04	0.82
3:A:301:MAN:O6	3:A:301:MAN:O4	1.98	0.80
1:A:42:CYS:SG	1:A:58:CYS:CB	2.70	0.80
1:A:186(B):GLY:C	1:A:221:ALA:HB1	2.08	0.79
1:A:124:PRO:HD3	1:A:209:LEU:O	1.83	0.77
1:A:32:SER:HB2	1:A:40:HIS:CD2	2.19	0.76
1:A:96:VAL:HG21	1:B:239:LEU:HD22	1.66	0.76
1:B:57:HIS:HE1	1:B:195:SER:OG	1.69	0.76
1:A:60:ASP:OD1	1:A:60(A):GLY:N	2.19	0.76
1:B:72:SER:HA	1:B:153:ASN:O	1.90	0.71
1:B:169:GLN:HE22	1:B:176:ILE:H	1.40	0.70
1:B:186:GLU:HG3	1:B:222:LYS:HG2	1.74	0.69
1:A:186:GLU:CG	1:A:222:LYS:HE3	2.24	0.68
1:A:164:PRO:HG2	1:A:167:GLU:HG3	1.77	0.67
1:B:91:HIS:CE1	1:B:101:TYR:CD1	2.84	0.65
1:B:128:ASP:O	1:B:131:THR:HG22	1.96	0.65
1:B:33:LEU:HD22	1:B:66:ILE:HG12	1.79	0.65
1:A:185:LYS:O	1:A:222:LYS:HA	1.97	0.63
1:B:180:MET:HE1	1:B:215:TRP:HZ2	1.63	0.63
1:A:71:LEU:HD12	1:A:72:SER:HB3	1.80	0.62
1:A:184(A):TYR:CB	1:A:186(A):GLY:CA	2.71	0.62
1:A:222:LYS:O	1:A:223:ASP:HB2	2.01	0.61
1:B:57:HIS:HD2	1:B:102:ASP:OD2	1.84	0.60
1:B:216:GLY:N	4:P:101:MRZ:H5	2.15	0.60
1:A:173:ARG:CZ	1:A:173:ARG:HB2	2.31	0.60
1:A:51:TRP:CE2	1:A:242:THR:HG22	2.37	0.60
1:B:186(B):GLY:HA2	1:B:221:ALA:HB1	1.86	0.58
1:A:133:TYR:HB3	1:A:201:CYS:SG	2.44	0.58
1:A:132:ILE:HG23	1:A:132:ILE:O	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:114:TYR:CZ	1:A:120:PRO:HD3	2.38	0.57
1:A:163:VAL:HG13	1:A:182:CYS:HB2	1.85	0.57
1:B:220:CYS:O	1:B:221:ALA:HB3	2.05	0.57
1:B:216:GLY:CA	4:P:101:MRZ:H5	2.34	0.57
1:A:96:VAL:HG21	1:B:239:LEU:CD2	2.35	0.56
1:A:164:PRO:HG2	1:A:167:GLU:CG	2.35	0.56
1:B:42:CYS:HB3	1:B:58:CYS:SG	2.40	0.56
1:B:24:LEU:HD12	1:B:24:LEU:O	2.06	0.56
1:A:17:VAL:HG23	1:A:144:THR:O	2.08	0.54
1:B:17:VAL:O	1:B:186(C):THR:HA	2.06	0.54
1:B:131:THR:CG2	1:B:133:TYR:CZ	2.91	0.54
1:B:169:GLN:NE2	1:B:176:ILE:H	2.06	0.54
1:A:67:TYR:CE1	1:A:82:SER:HB3	2.42	0.54
1:A:96:VAL:CG2	1:B:239:LEU:HD22	2.35	0.54
1:A:186(B):GLY:C	1:A:221:ALA:CB	2.79	0.53
1:B:59:PHE:HZ	1:B:106:ILE:HD11	1.73	0.53
1:A:179:GLN:CD	1:A:179:GLN:N	2.67	0.53
1:B:180:MET:CE	1:B:227:VAL:HG11	2.38	0.53
1:B:180:MET:HE3	1:B:215:TRP:HE1	1.74	0.53
1:B:239:LEU:HD23	1:B:239:LEU:N	2.23	0.52
1:A:71:LEU:HD12	1:A:72:SER:N	2.26	0.51
1:B:131:THR:HG23	1:B:133:TYR:CE2	2.46	0.50
1:A:66:ILE:HD11	1:A:106:ILE:HG21	1.93	0.50
1:B:35:VAL:HG11	1:B:60(B):ILE:HD12	1.94	0.50
1:A:23:SER:H	1:A:26:GLU:HG3	1.76	0.50
1:A:172:TYR:HB3	1:A:174:TYR:HB2	1.93	0.50
1:B:16:ILE:O	1:B:144:THR:HA	2.13	0.49
1:A:38(B):GLN:NE2	1:A:67:TYR:OH	2.44	0.49
1:B:17:VAL:CG1	1:B:186(B):GLY:O	2.55	0.49
1:B:57:HIS:CE1	1:B:195:SER:OG	2.57	0.49
1:B:23:SER:H	1:B:26:GLU:CD	2.21	0.49
1:B:185:LYS:O	1:B:223:ASP:N	2.45	0.48
1:A:36:LYS:HG3	1:A:38(B):GLN:HG3	1.95	0.48
1:A:138:VAL:HG22	1:A:199:LEU:HD13	1.95	0.48
1:B:180:MET:HE3	1:B:227:VAL:HG11	1.95	0.48
1:A:37:LEU:N	1:A:37:LEU:HD12	2.29	0.47
1:A:66:ILE:HD12	1:A:108:LEU:CD2	2.45	0.47
1:A:185:LYS:O	1:A:222:LYS:CA	2.61	0.47
1:B:147:GLN:HA	1:B:147:GLN:OE1	2.15	0.47
1:A:125:SER:O	1:A:128:ASP:HB2	2.15	0.47
1:A:126:LYS:O	1:A:126:LYS:HD3	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:72:SER:HA	1:A:153:ASN:O	2.15	0.46
1:B:95:LYS:HD3	1:B:95:LYS:HA	1.75	0.46
1:A:234:TYR:O	1:A:238:ILE:HG13	2.16	0.46
1:A:68:GLY:C	1:A:118:GLN:HE21	2.24	0.46
1:B:57:HIS:CE1	1:B:195:SER:HB3	2.50	0.46
1:A:198:PRO:HB3	1:A:209:LEU:CD1	2.46	0.46
1:A:221(A):ARG:HB2	1:A:224:GLN:HB2	1.99	0.45
1:B:158:ALA:HB1	1:B:186(C):THR:CG2	2.46	0.45
1:B:56:ALA:HB1	1:B:94:TYR:CD2	2.51	0.45
1:B:186(A):GLY:O	1:B:221:ALA:O	2.34	0.45
1:B:57:HIS:CE1	1:B:195:SER:CB	2.99	0.45
1:B:57:HIS:HE1	1:B:195:SER:CB	2.30	0.45
1:B:234:TYR:O	1:B:238:ILE:HG13	2.16	0.45
1:B:60(C):PRO:HD2	1:B:61:TYR:CE2	2.52	0.45
1:B:35:VAL:HG13	1:B:65(C):VAL:HG13	1.99	0.44
1:A:35:VAL:HG13	1:A:65(C):VAL:HG13	2.00	0.44
1:B:50:GLN:O	1:B:107:LYS:HA	2.18	0.44
1:A:45:SER:OG	1:A:198:PRO:HB3	2.18	0.43
1:B:131:THR:HG23	1:B:133:TYR:CZ	2.53	0.43
1:A:142:GLY:HA2	1:A:192:LYS:O	2.18	0.43
1:B:243:GLN:OE1	1:B:243:GLN:HA	2.19	0.43
1:A:85:ILE:HG12	1:A:106:ILE:CG2	2.48	0.43
1:A:167:GLU:O	1:A:170:LYS:HB2	2.19	0.43
1:A:66:ILE:HD12	1:A:108:LEU:HD21	2.01	0.43
1:B:192:LYS:HA	2:P:6:ALA:O	2.18	0.42
1:A:53:LEU:HD11	1:A:103:ILE:HD11	2.00	0.42
1:A:34:GLN:HG2	1:A:40:HIS:HA	2.01	0.42
1:A:189:ASP:HB2	1:A:221:ALA:HB2	2.02	0.42
1:B:158:ALA:HB1	1:B:186(C):THR:HG23	2.00	0.42
1:B:131:THR:CG2	1:B:133:TYR:CE2	3.02	0.41
1:A:23:SER:N	1:A:26:GLU:HG3	2.35	0.41
1:A:33:LEU:HD23	1:A:66:ILE:HG12	2.02	0.41
1:B:180:MET:CE	1:B:215:TRP:CZ2	3.04	0.41
1:A:143:TYR:HB2	1:A:191:CYS:HB3	2.02	0.41
1:A:142:GLY:CA	1:A:192:LYS:O	2.68	0.41
1:A:184(A):TYR:CG	1:A:186(A):GLY:HA3	2.50	0.40
1:B:180:MET:CE	1:B:215:TRP:HZ2	2.32	0.40
1:B:96:VAL:O	1:B:96:VAL:HG12	2.20	0.40
1:B:65(C):VAL:HG12	1:B:65(D):TRP:CD1	2.57	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	234/248 (94%)	209 (89%)	24 (10%)	1 (0%)	30	57
1	B	234/248 (94%)	216 (92%)	18 (8%)	0	100	100
2	P	8/10 (80%)	4 (50%)	3 (38%)	1 (12%)	0	0
All	All	476/506 (94%)	429 (90%)	45 (10%)	2 (0%)	30	57

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	P	2	PRO
1	A	205	GLY

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	203/213 (95%)	185 (91%)	18 (9%)	9	26
1	B	203/213 (95%)	177 (87%)	26 (13%)	4	13
2	P	8/8 (100%)	7 (88%)	1 (12%)	4	13
All	All	414/434 (95%)	369 (89%)	45 (11%)	6	18

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

Mol	Chain	Res	Type
1	A	23	SER

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Mol	Chain	Res	Type
1	A	24	LEU
1	A	58	CYS
1	A	74	SER
1	A	97	SER
1	A	121	ILE
1	A	126	LYS
1	A	132	ILE
1	A	139	THR
1	A	147	GLN
1	A	163	VAL
1	A	173	ARG
1	A	173(A)	ASP
1	A	179	GLN
1	A	186	GLU
1	A	186(C)	THR
1	A	195	SER
1	A	201	CYS
1	B	24	LEU
1	B	36	LYS
1	B	38	VAL
1	B	38(A)	SER
1	B	38(B)	GLN
1	B	39	THR
1	B	49	ARG
1	B	58	CYS
1	B	70	ILE
1	B	74	SER
1	B	77	THR
1	B	92	GLN
1	B	100	ASN
1	B	115	THR
1	B	116	GLU
1	B	129	THR
1	B	138	VAL
1	B	173	ARG
1	B	173(A)	ASP
1	B	179	GLN
1	B	202	LYS
1	B	203	HIS
1	B	223	ASP
1	B	236	ASP
1	B	238	ILE

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Mol	Chain	Res	Type
1	B	239	LEU
2	P	10	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38(B)	GLN
1	A	40	HIS
1	A	118	GLN
1	A	147	GLN
1	A	179	GLN
1	B	21	ASN
1	B	57	HIS
1	B	91	HIS
1	B	169	GLN
1	B	179	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

3 ligands 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
4	MRZ	P	101	-	8,9,9	0.65	0	11,11,11	1.26	1 (9%)
3	MAN	B	301	-	11,11,12	0.39	0	15,15,17	0.91	0
3	MAN	A	301	1	11,11,12	0.60	0	15,15,17	1.37	4 (26%)

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
4	MRZ	P	101	-	-	0/4/12/12	0/1/1/1
3	MAN	B	301	-	-	1/2/19/22	0/1/1/1
3	MAN	A	301	1	-	1/2/19/22	0/1/1/1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	P	101	MRZ	C3-N4-C5	2.69	118.17	112.68
3	A	301	MAN	O3-C3-C4	-2.43	104.65	110.38
3	A	301	MAN	O5-C5-C6	2.25	112.05	107.66
3	A	301	MAN	C1-C2-C3	2.14	112.75	109.64
3	A	301	MAN	C6-C5-C4	-2.04	108.01	113.02

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	301	MAN	O5-C5-C6-O6
3	B	301	MAN	O5-C5-C6-O6

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	P	101	MRZ	2	0
3	B	301	MAN	2	0
3	A	301	MAN	1	0

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	236/248 (95%)	0.50	19 (8%)	18 14	47, 74, 124, 164	2 (0%)
1	B	236/248 (95%)	0.36	14 (5%)	28 23	39, 64, 102, 131	3 (1%)
2	P	10/10 (100%)	1.96	3 (30%)	1 1	43, 94, 118, 147	0
All	All	482/506 (95%)	0.46	36 (7%)	20 16	39, 70, 117, 164	5 (1%)

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	65(B)	ASP	5.4
1	B	97	SER	5.1
2	P	5	SER	5.0
1	A	60	ASP	4.2
1	A	130	ASN	3.9
2	P	7	TYR	3.8
1	A	165	ASN	3.5
1	A	203	HIS	3.4
1	B	186(B)	GLY	3.1
1	A	215	TRP	3.1
1	A	71	LEU	3.0
1	B	220	CYS	3.0
1	B	153	ASN	2.9
1	B	186	GLU	2.8
1	A	202	LYS	2.7
1	B	96	VAL	2.7
1	B	74	SER	2.6
2	P	4	TYR	2.6
1	A	132	ILE	2.5
1	A	186(A)	GLY	2.4
1	A	201	CYS	2.4
1	B	65(B)	ASP	2.4
1	B	58	CYS	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	147	GLN	2.3
1	A	37	LEU	2.3
1	A	221(A)	ARG	2.3
1	A	16	ILE	2.3
1	A	169	GLN	2.3
1	A	204	SER	2.2
1	A	174	TYR	2.2
1	A	171	LYS	2.1
1	B	24	LEU	2.1
1	B	186(A)	GLY	2.1
1	A	168	CYS	2.0
1	B	191	CYS	2.0
1	B	166	GLU	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
3	MAN	B	301	11/12	0.68	0.13	91,109,114,117	0
3	MAN	A	301	11/12	0.78	0.11	54,71,81,82	0
4	MRZ	P	101	9/9	0.87	0.14	56,61,63,73	0

6.5 Other polymers [i](#)

There are no such residues in this entry.