



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 08:37 AM GMT

PDB ID : 1HWJ
Title : COMPLEX OF THE CATALYTIC PORTION OF HUMAN HMG-COA REDUCTASE WITH CERIVASTATIN
Authors : Istvan, E.S.; Deisenhofer, J.
Deposited on : 2001-01-09
Resolution : 2.26 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

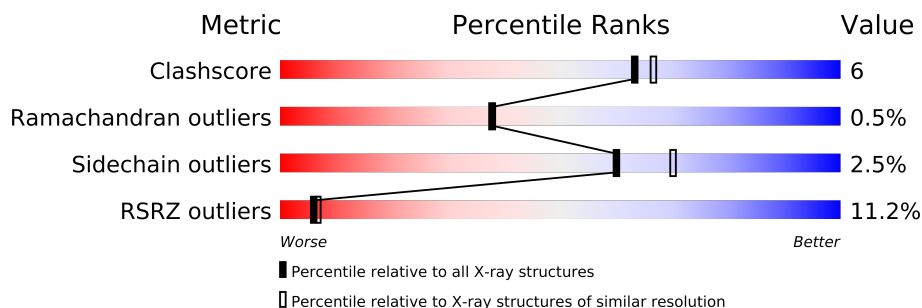
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.26 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))
Clashscore	79885	1326 (2.28-2.24)
Ramachandran outliers	78287	1291 (2.28-2.24)
Sidechain outliers	78261	1291 (2.28-2.24)
RSRZ outliers	66119	1110 (2.28-2.24)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	467	
1	B	467	
1	C	467	
1	D	467	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	ADP	B	105	-	X
2	ADP	D	106	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 12418 atoms, of which 0 are hydrogen and 0 are deuterium.

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 HMG-COA REDUCTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	410	Total	C	N	O	S	0	0	0
			3052	1900	537	585	30			
1	B	398	Total	C	N	O	S	0	0	0
			2952	1838	518	567	29			
1	C	408	Total	C	N	O	S	0	0	0
			3034	1888	533	583	30			
1	D	392	Total	C	N	O	S	0	0	0
			2900	1803	508	560	29			

There are 20 discrepancies between the modelled and reference sequences:

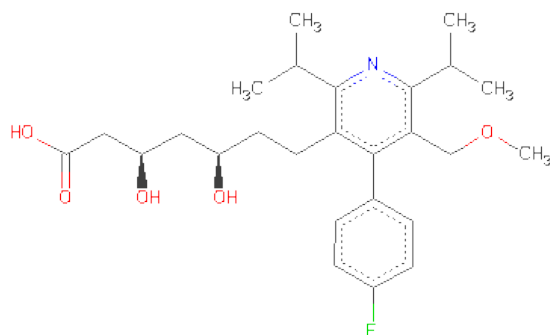
Chain	Residue	Modelled	Actual	Comment	Reference
A	422	GLY	-	INSERTION	UNP P04035
A	423	ALA	-	INSERTION	UNP P04035
A	424	MET	-	INSERTION	UNP P04035
A	425	ALA	-	INSERTION	UNP P04035
A	485	ILE	MET	ENGINEERED	UNP P04035
B	422	GLY	-	INSERTION	UNP P04035
B	423	ALA	-	INSERTION	UNP P04035
B	424	MET	-	INSERTION	UNP P04035
B	425	ALA	-	INSERTION	UNP P04035
B	485	ILE	MET	ENGINEERED	UNP P04035
C	422	GLY	-	INSERTION	UNP P04035
C	423	ALA	-	INSERTION	UNP P04035
C	424	MET	-	INSERTION	UNP P04035
C	425	ALA	-	INSERTION	UNP P04035
C	485	ILE	MET	ENGINEERED	UNP P04035
D	422	GLY	-	INSERTION	UNP P04035
D	423	ALA	-	INSERTION	UNP P04035
D	424	MET	-	INSERTION	UNP P04035
D	425	ALA	-	INSERTION	UNP P04035
D	485	ILE	MET	ENGINEERED	UNP P04035

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	C	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	C	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	D	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 3 is 7-[4-(4-FLUORO-PHENYL)-5-HYDROXYMETHYL-2,6-DIISOPROPYL-PYRIDIN-3-YL]-3,5-DIHYDROXY-HEPTANOICACID (three-letter code: 116) (formula: $C_{26}H_{36}FNO_5$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	1	Total	C	F	N	O	0	0
			33	26	1	1	5		
3	A	1	Total	C	F	N	O	0	0
			33	26	1	1	5		
3	D	1	Total	C	F	N	O	0	0
			33	26	1	1	5		
3	C	1	Total	C	F	N	O	0	0
			33	26	1	1	5		

- Molecule 4 is water.

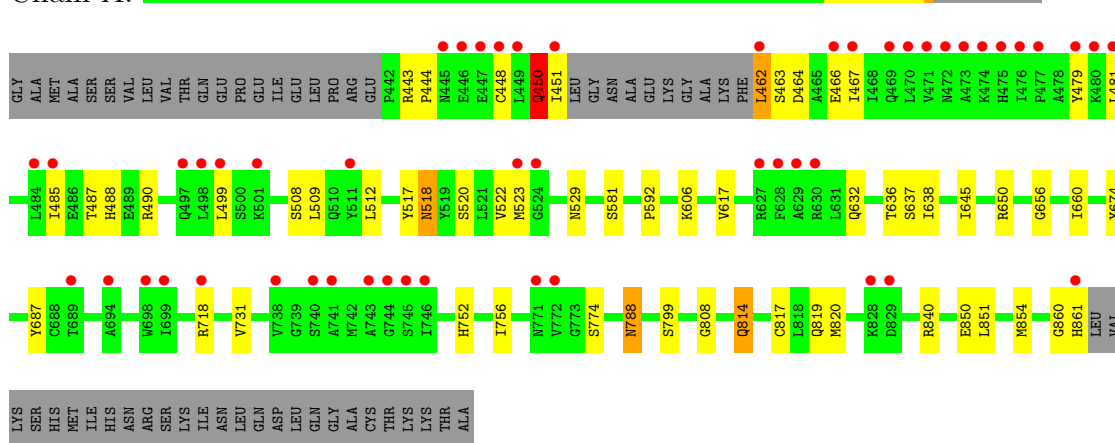
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	52	Total	O	0	0
			52	52		
4	B	52	Total	O	0	0
			52	52		
4	C	33	Total	O	0	0
			33	33		
4	D	49	Total	O	0	0
			49	49		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

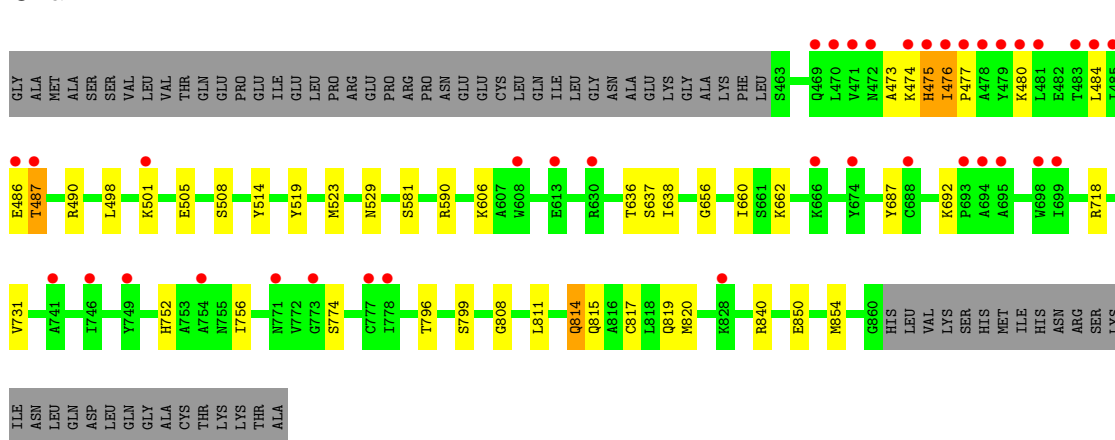
• Molecule 1: HMG-COA REDUCTASE

Chain A:



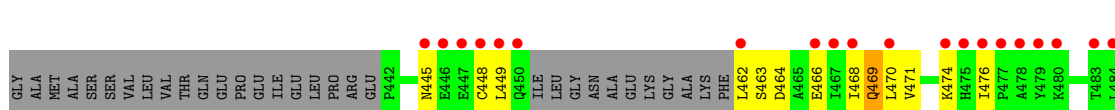
• Molecule 1: HMG-COA REDUCTASE

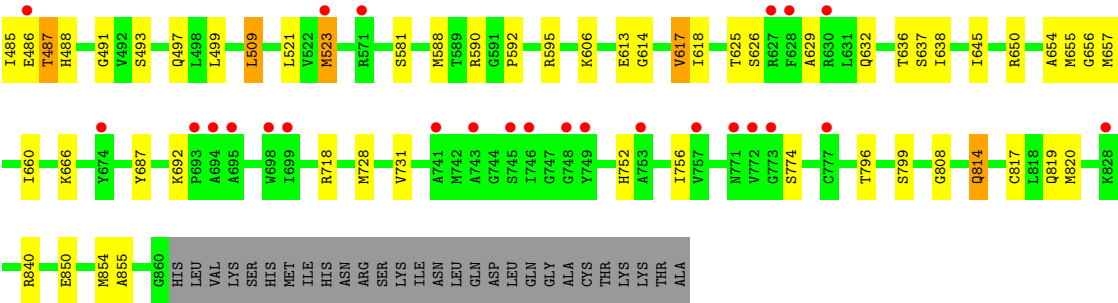
Chain B:



• Molecule 1: HMG-COA REDUCTASE

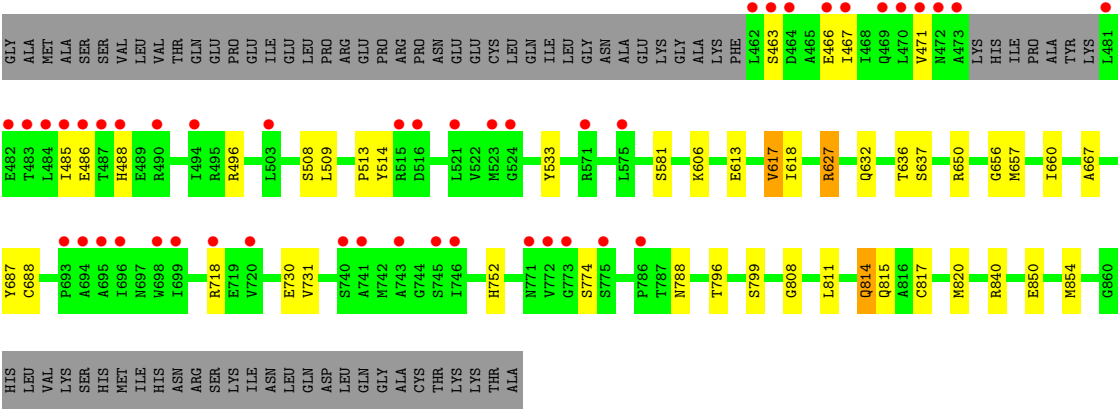
Chain C:





• Molecule 1: HMG-COA REDUCTASE

Chain D:



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	74.62Å 172.98Å 80.16Å 90.00° 117.35° 90.00°	Depositor
Resolution (Å)	43.50 – 2.26 54.97 – 2.26	Depositor EDS
% Data completeness (in resolution range)	96.0 (43.50-2.26) 93.0 (54.97-2.26)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.23 (at 2.27Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.221 , 0.237 0.216 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	40.2	Xtriage
Anisotropy	0.292	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 40.2	EDS
Estimated twinning fraction	0.019 for h,-k,-h-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	4 of 80878 reflections (0.005%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	12418	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.81% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 116, ADP

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	0.49	0/3096	0.66	0/4186
1	B	0.46	0/2994	0.65	0/4049
1	C	0.47	0/3077	0.66	1/4160 (0.0%)
1	D	0.51	0/2938	0.65	0/3972
All	All	0.48	0/12105	0.65	1/16367 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	656	GLY	N-CA-C	5.29	126.34	113.10

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3052	0	3087	33	0
1	B	2952	0	2989	29	0
1	C	3034	0	3069	45	0
1	D	2900	0	2934	27	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	54	0	24	1	0
2	B	27	0	12	0	0
2	C	54	0	24	1	0
2	D	27	0	12	0	0
3	A	33	0	35	2	0
3	B	33	0	35	2	0
3	C	33	0	35	3	0
3	D	33	0	35	1	0
4	A	52	0	0	1	0
4	B	52	0	0	1	0
4	C	33	0	0	1	0
4	D	49	0	0	2	0
All	All	12418	0	12291	135	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 6.

All (135) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:485:ILE:HG22	1:C:486:GLU:H	1.37	0.89
1:A:487:THR:HG22	1:A:488:HIS:H	1.43	0.83
1:B:523:MET:HE1	4:B:1019:HOH:O	1.81	0.79
1:B:817:CYS:HA	1:B:820:MET:HE3	1.68	0.76
1:D:817:CYS:HA	1:D:820:MET:HE3	1.70	0.74
1:C:817:CYS:HA	1:C:820:MET:HE3	1.70	0.72
3:B:1:116:H91	3:B:1:116:H61	1.73	0.69
1:A:817:CYS:HA	1:A:820:MET:HE3	1.74	0.69
1:C:448:CYS:HB3	1:C:462:LEU:HD22	1.73	0.68
1:A:523:MET:HE1	4:A:1086:HOH:O	1.95	0.67
1:D:688:CYS:SG	4:D:1163:HOH:O	2.54	0.66
1:C:466:GLU:O	1:C:470:LEU:HG	1.96	0.65
1:D:467:ILE:O	1:D:471:VAL:HG23	1.97	0.64
1:C:655:MET:SD	1:C:657:MET:HG2	2.38	0.63
1:C:590:ARG:NH2	1:C:657:MET:HE3	2.14	0.62
1:C:485:ILE:HG22	1:C:486:GLU:N	2.11	0.62
1:C:632:GLN:HE21	1:C:650:ARG:HG3	1.65	0.60
3:A:2:116:H91	3:A:2:116:H61	1.84	0.60
1:D:656:GLY:O	1:D:660:ILE:HG12	2.03	0.58
1:C:588:MET:HB3	1:C:660:ILE:CD1	2.34	0.57
1:D:808:GLY:O	1:D:814:GLN:HG3	2.04	0.57
1:B:476:ILE:HG21	1:B:484:LEU:HD11	1.86	0.57
1:C:581:SER:OG	1:C:840:ARG:HD2	2.04	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:581:SER:OG	1:D:840:ARG:HD2	2.04	0.57
1:C:523:MET:HE1	4:C:1102:HOH:O	2.03	0.57
1:B:656:GLY:O	1:B:660:ILE:HG12	2.05	0.57
1:B:477:PRO:HD2	1:B:480:LYS:CG	2.35	0.57
1:A:581:SER:OG	1:A:840:ARG:HD2	2.05	0.56
1:C:808:GLY:O	1:C:814:GLN:HG3	2.07	0.55
1:A:656:GLY:O	1:A:660:ILE:HG12	2.06	0.54
1:B:581:SER:OG	1:B:840:ARG:HD2	2.08	0.54
3:C:4:116:H91	3:C:4:116:H61	1.90	0.53
3:D:3:116:H91	3:D:3:116:H61	1.90	0.53
1:C:463:SER:OG	1:C:466:GLU:HG2	2.08	0.53
1:C:654:ALA:HB1	2:C:104:ADP:O1B	2.09	0.52
1:B:519:TYR:O	1:B:523:MET:HG2	2.08	0.52
1:D:627:ARG:H	1:D:627:ARG:HD3	1.75	0.52
1:B:477:PRO:HD2	1:B:480:LYS:HG3	1.92	0.51
1:B:487:THR:HG23	1:B:490:ARG:HB3	1.93	0.51
1:C:625:THR:HG23	1:C:666:LYS:HD3	1.93	0.51
1:A:487:THR:HG22	1:A:488:HIS:N	2.21	0.50
1:B:808:GLY:O	1:B:814:GLN:HG3	2.11	0.50
1:C:632:GLN:NE2	1:C:650:ARG:HG3	2.26	0.50
1:C:626:SER:HB3	1:C:629:ALA:HB2	1.93	0.50
1:D:485:ILE:HG22	1:D:486:GLU:N	2.27	0.50
1:B:774:SER:HA	1:B:799:SER:O	2.13	0.48
1:A:774:SER:HA	1:A:799:SER:O	2.13	0.48
1:D:774:SER:HA	1:D:799:SER:O	2.14	0.48
1:C:637:SER:HB2	1:C:687:TYR:OH	2.13	0.48
1:C:606:LYS:HG3	1:C:636:THR:OG1	2.14	0.48
1:A:448:CYS:HA	1:A:462:LEU:HD21	1.95	0.47
1:A:632:GLN:HE21	1:A:650:ARG:HG3	1.79	0.47
1:A:479:TYR:HB3	1:A:529:ASN:OD1	2.15	0.47
1:C:464:ASP:O	1:C:468:ILE:HG12	2.15	0.47
1:C:487:THR:OG1	1:C:488:HIS:N	2.48	0.47
1:B:637:SER:HB2	1:B:687:TYR:OH	2.15	0.47
1:C:523:MET:CE	1:C:523:MET:HA	2.45	0.47
1:A:637:SER:HB2	1:A:687:TYR:OH	2.14	0.47
1:D:463:SER:HB2	1:D:466:GLU:HG3	1.97	0.47
1:D:637:SER:HB2	1:D:687:TYR:OH	2.15	0.47
1:B:473:ALA:O	1:B:475:HIS:N	2.49	0.46
1:A:518:ASN:ND2	1:A:520:SER:OG	2.49	0.46
1:C:499:LEU:HD23	1:C:509:LEU:HD21	1.98	0.46
1:A:808:GLY:O	1:A:814:GLN:HG3	2.15	0.46
1:A:820:MET:HB2	1:A:820:MET:HE2	1.73	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:850:GLU:O	1:C:854:MET:HG2	2.14	0.46
1:C:655:MET:CE	1:C:657:MET:HG2	2.45	0.46
1:C:474:LYS:HB2	1:C:476:ILE:HG13	1.98	0.46
1:A:463:SER:HB3	1:A:466:GLU:CD	2.37	0.45
1:C:613:GLU:O	1:C:617:VAL:HG12	2.17	0.45
1:C:774:SER:HA	1:C:799:SER:O	2.16	0.45
1:C:471:VAL:HG13	1:C:476:ILE:O	2.17	0.45
1:A:606:LYS:HG3	1:A:636:THR:OG1	2.17	0.45
1:A:467:ILE:HD11	1:A:490:ARG:HH21	1.82	0.45
1:C:820:MET:HB2	1:C:820:MET:HE2	1.86	0.44
1:B:850:GLU:O	1:B:854:MET:HG2	2.16	0.44
1:D:820:MET:HB2	1:D:820:MET:HE2	1.85	0.44
1:A:499:LEU:HD23	1:A:509:LEU:HD21	2.00	0.44
1:C:485:ILE:HD12	1:C:491:GLY:HA2	1.98	0.44
1:D:850:GLU:O	1:D:854:MET:HG2	2.17	0.44
1:B:638:ILE:O	1:C:796:THR:HG21	2.18	0.44
1:D:606:LYS:HG3	1:D:636:THR:OG1	2.18	0.43
1:C:592:PRO:HD2	1:C:645:ILE:O	2.17	0.43
1:C:731:VAL:HG12	1:C:854:MET:CE	2.48	0.43
1:C:614:GLY:O	1:C:618:ILE:HG12	2.17	0.43
1:A:850:GLU:O	1:A:854:MET:HG2	2.19	0.43
1:B:796:THR:HG21	1:C:638:ILE:O	2.18	0.43
1:D:731:VAL:HG12	1:D:854:MET:HE3	2.00	0.43
1:A:450:GLN:HB2	1:A:451:ILE:H	1.60	0.43
1:B:606:LYS:HG3	1:B:636:THR:OG1	2.18	0.43
1:B:731:VAL:HG12	1:B:854:MET:CE	2.48	0.43
1:C:595:ARG:HH22	1:D:730:GLU:HG2	1.84	0.43
1:A:638:ILE:O	1:D:796:THR:HG21	2.19	0.43
1:A:617:VAL:HG11	1:A:674:TYR:CE1	2.53	0.43
1:A:508:SER:HB2	1:B:819:GLN:CB	2.48	0.42
1:B:498:LEU:O	1:B:501:LYS:HG2	2.20	0.42
2:A:102:ADP:H2	1:B:529:ASN:HD21	1.66	0.42
1:D:618:ILE:HG23	1:D:667:ALA:HB1	2.02	0.42
1:C:692:LYS:HB2	1:C:692:LYS:HE2	1.91	0.42
1:B:486:GLU:O	1:B:487:THR:HB	2.20	0.42
1:A:731:VAL:HG12	1:A:854:MET:CE	2.49	0.42
1:B:590:ARG:HD3	1:B:590:ARG:HA	1.80	0.42
1:A:632:GLN:NE2	1:A:650:ARG:HG3	2.34	0.42
1:D:731:VAL:HG12	1:D:854:MET:CE	2.50	0.42
1:A:756:ILE:HD12	1:A:756:ILE:N	2.34	0.42
1:A:481:LEU:O	1:A:485:ILE:HG13	2.19	0.42
1:D:627:ARG:N	1:D:627:ARG:HD3	2.35	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:819:GLN:CB	1:B:508:SER:HB2	2.50	0.42
1:D:811:LEU:O	1:D:815:GLN:HG3	2.19	0.42
1:C:756:ILE:HD12	1:C:756:ILE:N	2.35	0.41
1:D:513:PRO:HB2	1:D:533:TYR:CZ	2.55	0.41
1:D:613:GLU:O	1:D:617:VAL:HG13	2.20	0.41
1:C:731:VAL:HG12	1:C:854:MET:HE3	2.01	0.41
1:B:692:LYS:HB2	1:B:692:LYS:HE2	1.94	0.41
1:C:590:ARG:HA	1:C:590:ARG:HD3	1.80	0.41
1:A:462:LEU:HB2	1:A:463:SER:H	1.52	0.41
1:A:788:ASN:HD22	1:A:788:ASN:HA	1.68	0.41
3:A:2:116:H72	3:A:2:116:C86	2.51	0.41
1:A:592:PRO:HD2	1:A:645:ILE:O	2.20	0.41
1:A:517:TYR:HE2	1:A:522:VAL:HG21	1.86	0.41
1:B:662:LYS:HB3	1:B:662:LYS:HE2	1.93	0.41
1:B:476:ILE:HG22	1:B:480:LYS:HB2	2.02	0.41
3:C:4:116:H72	3:C:4:116:C86	2.51	0.41
1:C:493:SER:O	1:C:497:GLN:HG3	2.21	0.41
1:D:496:ARG:NH2	1:D:509:LEU:O	2.53	0.41
1:B:811:LEU:O	1:B:815:GLN:HG3	2.21	0.41
3:B:1:116:C91	3:B:1:116:H61	2.46	0.40
3:C:4:116:C91	3:C:4:116:H61	2.51	0.40
1:C:728:MET:SD	1:C:855:ALA:HA	2.61	0.40
1:A:443:ARG:HA	1:A:444:PRO:HD3	1.86	0.40
1:B:756:ILE:HD12	1:B:756:ILE:N	2.36	0.40
1:D:632:GLN:HE21	1:D:650:ARG:HG3	1.85	0.40
1:D:657:MET:HG2	4:D:1186:HOH:O	2.21	0.40
1:C:445:ASN:OD1	1:C:469:GLN:HB3	2.21	0.40
1:C:819:GLN:HB2	1:D:508:SER:HB2	2.02	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	406/467 (87%)	386 (95%)	18 (4%)	2 (0%)	38 38

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	396/467 (85%)	375 (95%)	17 (4%)	4 (1%)	22	17
1	C	404/467 (86%)	379 (94%)	24 (6%)	1 (0%)	56	63
1	D	388/467 (83%)	371 (96%)	16 (4%)	1 (0%)	50	54
All	All	1594/1868 (85%)	1511 (95%)	75 (5%)	8 (0%)	38	38

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	450	GLN
1	A	860	GLY
1	B	474	LYS
1	B	475	HIS
1	B	487	THR
1	C	449	LEU
1	B	514	TYR
1	D	514	TYR

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	328/375 (88%)	317 (97%)	11 (3%)	49	57
1	B	316/375 (84%)	311 (98%)	5 (2%)	75	84
1	C	326/375 (87%)	317 (97%)	9 (3%)	56	65
1	D	311/375 (83%)	304 (98%)	7 (2%)	63	72
All	All	1281/1500 (85%)	1249 (98%)	32 (2%)	60	70

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	450	GLN
1	A	462	LEU
1	A	464	ASP
1	A	512	LEU

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Mol	Chain	Res	Type
1	A	518	ASN
1	A	718	ARG
1	A	752	HIS
1	A	788	ASN
1	A	814	GLN
1	A	851	LEU
1	A	861	HIS
1	B	476	ILE
1	B	505	GLU
1	B	718	ARG
1	B	752	HIS
1	B	814	GLN
1	C	469	GLN
1	C	487	THR
1	C	509	LEU
1	C	521	LEU
1	C	523	MET
1	C	617	VAL
1	C	718	ARG
1	C	752	HIS
1	C	814	GLN
1	D	488	HIS
1	D	617	VAL
1	D	627	ARG
1	D	718	ARG
1	D	752	HIS
1	D	788	ASN
1	D	814	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	518	ASN
1	A	635	HIS
1	A	788	ASN
1	A	819	GLN
1	B	469	GLN
1	B	488	HIS
1	B	635	HIS
1	B	819	GLN
1	C	450	GLN
1	C	469	GLN

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Mol	Chain	Res	Type
1	C	510	GLN
1	C	632	GLN
1	C	635	HIS
1	C	819	GLN
1	D	469	GLN
1	D	635	HIS
1	D	788	ASN
1	D	819	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

10 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	$\# Z > 2$
2	ADP	A	101	-	29,29,29	1.31	3 (10%)	45,45,45	0.89	1 (2%)
2	ADP	A	102	-	29,29,29	1.39	4 (13%)	45,45,45	0.80	0
3	116	A	2	-	34,34,34	1.83	7 (20%)	47,47,47	1.46	6 (12%)
3	116	B	1	-	34,34,34	1.93	8 (23%)	47,47,47	1.51	7 (14%)
2	ADP	B	105	-	29,29,29	1.72	9 (31%)	45,45,45	0.96	1 (2%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	ADP	C	103	-	29,29,29	1.31	2 (6%)	45,45,45	1.11	3 (6%)
2	ADP	C	104	-	29,29,29	1.48	5 (17%)	45,45,45	0.78	0
3	116	C	4	-	34,34,34	1.92	6 (17%)	47,47,47	1.52	6 (12%)
2	ADP	D	106	-	29,29,29	1.58	7 (24%)	45,45,45	0.92	2 (4%)
3	116	D	3	-	34,34,34	2.00	6 (17%)	47,47,47	1.41	5 (10%)

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	ADP	A	101	-	-	0/16/32/32	0/1/3/3
2	ADP	A	102	-	-	0/16/32/32	0/1/3/3
3	116	A	2	-	-	0/28/28/28	0/2/2/2
3	116	B	1	-	-	0/28/28/28	0/2/2/2
2	ADP	B	105	-	-	0/16/32/32	0/1/3/3
2	ADP	C	103	-	-	0/16/32/32	0/1/3/3
2	ADP	C	104	-	-	0/16/32/32	0/1/3/3
3	116	C	4	-	-	0/28/28/28	0/2/2/2
2	ADP	D	106	-	-	0/16/32/32	0/1/3/3
3	116	D	3	-	-	0/28/28/28	0/2/2/2

All (57) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	3	116	C12-C15	6.00	1.59	1.51
3	D	3	116	C8-C81	-5.90	1.38	1.50
3	C	4	116	C8-C81	-5.14	1.40	1.50
3	A	2	116	C12-C15	5.04	1.58	1.51
3	C	4	116	C12-C15	5.02	1.58	1.51
3	B	1	116	C8-C81	-4.91	1.40	1.50
3	B	1	116	C12-C15	4.83	1.58	1.51
3	A	2	116	C8-C81	-4.68	1.40	1.50
2	B	105	ADP	PA-O3A	4.61	1.68	1.59
3	C	4	116	C9-C91	4.21	1.57	1.51
3	B	1	116	C9-C91	4.18	1.57	1.51
3	D	3	116	C9-C91	4.07	1.56	1.51
2	D	106	ADP	PA-O3A	3.89	1.66	1.59
3	D	3	116	C6-C7	-3.89	1.33	1.53
3	A	2	116	C6-C7	-3.70	1.34	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	4	116	C6-C7	-3.69	1.34	1.53
3	A	2	116	C9-C91	3.67	1.56	1.51
2	C	104	ADP	PA-O3A	3.52	1.66	1.59
3	B	1	116	C13-C11	3.47	1.57	1.51
3	B	1	116	C6-C7	-3.40	1.35	1.53
2	C	103	ADP	C8-N7	-3.40	1.27	1.34
2	A	102	ADP	PA-O3A	3.12	1.65	1.59
2	D	106	ADP	C8-N7	-3.11	1.28	1.34
2	A	101	ADP	C8-N7	-3.04	1.28	1.34
2	B	105	ADP	O4'-C1'	3.03	1.46	1.41
2	A	102	ADP	C8-N7	-3.01	1.28	1.34
2	B	105	ADP	C8-N7	-2.95	1.28	1.34
2	A	101	ADP	PA-O3A	2.92	1.65	1.59
2	C	104	ADP	C8-N7	-2.90	1.28	1.34
2	B	105	ADP	O4'-C4'	2.87	1.51	1.45
3	C	4	116	C7-C10	2.85	1.57	1.52
2	A	102	ADP	C8-N9	-2.70	1.32	1.36
2	C	104	ADP	C8-N9	-2.68	1.32	1.36
2	C	104	ADP	O4'-C4'	2.59	1.51	1.45
2	D	106	ADP	O4'-C4'	2.56	1.51	1.45
2	D	106	ADP	C8-N9	-2.50	1.32	1.36
2	A	101	ADP	C8-N9	-2.46	1.33	1.36
2	B	105	ADP	C8-N9	-2.43	1.33	1.36
2	B	105	ADP	C2'-C3'	2.38	1.60	1.53
3	B	1	116	C8-C11	2.38	1.45	1.41
2	C	103	ADP	C8-N9	-2.35	1.33	1.36
3	B	1	116	C7-C10	2.34	1.56	1.52
3	D	3	116	C13-C11	2.29	1.55	1.51
3	A	2	116	C7-C10	2.28	1.56	1.52
2	D	106	ADP	C1'-N9	2.27	1.55	1.48
3	C	4	116	C83-C84	2.21	1.41	1.36
3	A	2	116	C13-C11	2.19	1.54	1.51
2	A	102	ADP	O4'-C4'	2.14	1.50	1.45
3	A	2	116	C4-C5	2.13	1.57	1.52
2	D	106	ADP	O4'-C1'	2.13	1.44	1.41
3	B	1	116	C8-C10	2.11	1.45	1.41
2	B	105	ADP	C1'-N9	2.10	1.55	1.48
3	D	3	116	C8-C11	2.10	1.45	1.41
2	D	106	ADP	C2'-C3'	2.10	1.59	1.53
2	B	105	ADP	O2'-C2'	2.06	1.47	1.43
2	B	105	ADP	C5-C4	-2.04	1.35	1.40
2	C	104	ADP	O4'-C1'	2.01	1.44	1.41

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	4	116	C7-C6-C5	5.45	126.68	115.40
3	D	3	116	C7-C6-C5	4.85	125.44	115.40
3	B	1	116	C7-C6-C5	4.70	125.14	115.40
3	A	2	116	C7-C6-C5	4.66	125.06	115.40
3	C	4	116	C10-C9-N1	-4.30	119.40	123.19
3	D	3	116	C11-C12-N1	-4.02	119.65	123.19
3	A	2	116	C10-C9-N1	-3.87	119.78	123.19
2	B	105	ADP	O4'-C1'-N9	3.85	112.02	108.44
3	B	1	116	C10-C9-N1	-3.83	119.82	123.19
3	B	1	116	C6-C7-C10	3.83	124.12	112.28
3	D	3	116	C10-C9-N1	-3.71	119.93	123.19
3	A	2	116	C6-C7-C10	3.70	123.73	112.28
3	A	2	116	C11-C12-N1	-3.67	119.96	123.19
2	D	106	ADP	O4'-C1'-N9	3.65	111.84	108.44
3	B	1	116	C11-C12-N1	-3.58	120.04	123.19
3	C	4	116	C11-C12-N1	-3.47	120.14	123.19
3	C	4	116	C6-C7-C10	3.39	122.78	112.28
2	C	103	ADP	O4'-C1'-N9	-3.39	105.28	108.44
3	D	3	116	C6-C7-C10	3.19	122.16	112.28
2	C	103	ADP	C4'-O4'-C1'	2.74	112.73	109.75
3	B	1	116	O13-C13-C11	2.61	112.58	108.29
2	A	101	ADP	O4'-C1'-N9	-2.59	106.03	108.44
3	B	1	116	C12-N1-C9	2.50	125.43	115.98
3	C	4	116	C12-N1-C9	2.45	125.27	115.98
3	A	2	116	C12-N1-C9	2.37	124.96	115.98
3	D	3	116	C12-N1-C9	2.36	124.93	115.98
3	A	2	116	O13-C13-C11	2.22	111.95	108.29
2	C	103	ADP	C3'-C2'-C1'	2.11	104.21	100.91
3	B	1	116	C85-C84-C83	-2.09	120.06	122.90
3	C	4	116	C92-C91-C9	2.07	114.99	111.28
2	D	106	ADP	N3-C4-N9	2.06	129.15	125.43

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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 ⓘ

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	410/467 (87%)	0.58	51 (12%) 5 5	27, 51, 99, 101	0
1	B	398/467 (85%)	0.57	38 (9%) 8 10	30, 50, 86, 101	0
1	C	408/467 (87%)	0.70	45 (11%) 6 7	30, 53, 100, 101	0
1	D	392/467 (83%)	0.53	46 (11%) 5 6	29, 48, 93, 101	0
All	All	1608/1868 (86%)	0.60	180 (11%) 6 7	27, 50, 99, 101	0

All (180) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	462	LEU	8.2
1	D	483	THR	7.9
1	A	473	ALA	7.6
1	C	475	HIS	7.4
1	A	861	HIS	7.1
1	B	471	VAL	6.9
1	D	484	LEU	6.8
1	A	484	LEU	6.6
1	A	449	LEU	6.2
1	D	486	GLU	6.2
1	B	475	HIS	6.1
1	C	483	THR	5.9
1	D	471	VAL	5.9
1	D	485	ILE	5.9
1	D	470	LEU	5.6
1	D	469	GLN	5.5
1	A	475	HIS	5.4
1	A	476	ILE	5.4
1	B	476	ILE	5.4
1	D	523	MET	5.3
1	B	485	ILE	5.3

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Mol	Chain	Res	Type	RSRZ
1	A	448	CYS	5.2
1	C	484	LEU	5.2
1	C	486	GLU	5.1
1	A	451	ILE	4.8
1	C	477	PRO	4.7
1	D	524	GLY	4.7
1	A	470	LEU	4.7
1	D	481	LEU	4.7
1	D	487	THR	4.6
1	D	467	ILE	4.5
1	D	473	ALA	4.4
1	B	470	LEU	4.4
1	B	486	GLU	4.4
1	A	479	TYR	4.2
1	B	613	GLU	4.2
1	A	471	VAL	4.1
1	B	487	THR	4.0
1	A	446	GLU	3.9
1	A	467	ILE	3.9
1	D	694	ALA	3.9
1	C	447	GLU	3.8
1	C	448	CYS	3.7
1	C	627	ARG	3.7
1	D	482	GLU	3.6
1	A	477	PRO	3.5
1	C	450	GLN	3.5
1	C	628	PHE	3.5
1	A	828	LYS	3.5
1	B	479	TYR	3.5
1	A	462	LEU	3.4
1	B	474	LYS	3.4
1	C	476	ILE	3.3
1	D	466	GLU	3.3
1	A	469	GLN	3.3
1	B	484	LEU	3.2
1	D	494	ILE	3.2
1	B	698	TRP	3.2
1	A	630	ARG	3.2
1	C	828	LYS	3.2
1	D	516	ASP	3.2
1	B	746	ILE	3.2
1	C	445	ASN	3.2

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Mol	Chain	Res	Type	RSRZ
1	C	446	GLU	3.1
1	A	481	LEU	3.1
1	B	630	ARG	3.1
1	D	695	ALA	3.1
1	D	743	ALA	3.1
1	D	490	ARG	3.1
1	D	463	SER	3.1
1	A	718	ARG	3.1
1	B	694	ALA	3.1
1	D	488	HIS	3.0
1	D	718	ARG	3.0
1	A	741	ALA	3.0
1	C	749	TYR	3.0
1	B	483	THR	3.0
1	D	693	PRO	3.0
1	C	523	MET	3.0
1	D	472	ASN	3.0
1	B	608	TRP	3.0
1	C	630	ARG	3.0
1	B	666	LYS	3.0
1	A	472	ASN	3.0
1	C	695	ALA	3.0
1	A	627	ARG	2.9
1	C	698	TRP	2.8
1	B	688	CYS	2.8
1	C	479	TYR	2.8
1	A	698	TRP	2.7
1	C	470	LEU	2.7
1	C	699	ILE	2.7
1	C	693	PRO	2.7
1	B	699	ILE	2.7
1	C	694	ALA	2.7
1	C	480	LYS	2.7
1	A	699	ILE	2.7
1	C	745	SER	2.7
1	A	474	LYS	2.7
1	A	497	GLN	2.6
1	C	771	ASN	2.6
1	A	447	GLU	2.6
1	C	746	ILE	2.6
1	D	699	ILE	2.6
1	A	628	PHE	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	501	LYS	2.5
1	B	693	PRO	2.5
1	D	773	GLY	2.5
1	A	485	ILE	2.5
1	B	754	ALA	2.5
1	A	466	GLU	2.5
1	D	698	TRP	2.5
1	C	773	GLY	2.5
1	D	746	ILE	2.5
1	C	757	VAL	2.5
1	B	501	LYS	2.5
1	D	515	ARG	2.4
1	B	741	ALA	2.4
1	C	478	ALA	2.4
1	A	523	MET	2.4
1	C	466	GLU	2.4
1	B	695	ALA	2.4
1	C	753	ALA	2.4
1	D	464	ASP	2.4
1	B	478	ALA	2.4
1	A	772	VAL	2.4
1	D	772	VAL	2.4
1	A	744	GLY	2.4
1	C	571	ARG	2.4
1	A	445	ASN	2.4
1	D	521	LEU	2.4
1	D	575	LEU	2.3
1	B	771	ASN	2.3
1	B	480	LYS	2.3
1	D	571	ARG	2.3
1	C	674	TYR	2.3
1	A	629	ALA	2.3
1	A	745	SER	2.3
1	A	771	ASN	2.3
1	A	511	TYR	2.3
1	D	741	ALA	2.3
1	A	689	THR	2.3
1	C	741	ALA	2.3
1	B	773	GLY	2.3
1	D	696	ILE	2.3
1	A	524	GLY	2.2
1	C	777	CYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	743	ALA	2.2
1	B	828	LYS	2.2
1	A	694	ALA	2.2
1	C	474	LYS	2.2
1	D	740	SER	2.2
1	A	746	ILE	2.2
1	A	829	ASP	2.2
1	A	499	LEU	2.2
1	B	469	GLN	2.1
1	C	467	ILE	2.1
1	C	449	LEU	2.1
1	D	720	VAL	2.1
1	B	777	CYS	2.1
1	C	743	ALA	2.1
1	B	472	ASN	2.1
1	D	745	SER	2.1
1	B	674	TYR	2.1
1	B	749	TYR	2.1
1	A	480	LYS	2.1
1	C	748	GLY	2.1
1	A	738	VAL	2.1
1	C	468	ILE	2.1
1	A	740	SER	2.1
1	A	498	LEU	2.1
1	D	503	LEU	2.1
1	C	772	VAL	2.1
1	B	477	PRO	2.0
1	D	771	ASN	2.0
1	D	786	PRO	2.0
1	D	775	SER	2.0
1	B	778	ILE	2.0
1	B	481	LEU	2.0
1	C	462	LEU	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	ADP	B	105	27/27	0.30	4.85	99,100,100,100	0
2	ADP	D	106	27/27	0.29	3.43	99,100,100,100	0
2	ADP	C	103	27/27	0.31	1.34	91,95,100,100	0
2	ADP	C	104	27/27	0.23	1.00	98,100,100,100	0
2	ADP	A	102	27/27	0.23	0.64	86,91,100,100	0
3	116	C	4	33/33	0.16	0.42	39,45,48,50	0
3	116	B	1	33/33	0.16	0.41	42,49,54,55	0
2	ADP	A	101	27/27	0.20	0.14	88,92,100,100	0
3	116	D	3	33/33	0.15	0.06	39,46,49,49	0
3	116	A	2	33/33	0.14	-0.28	35,40,44,46	0

6.5 Other polymers ⓘ

There are no such residues in this entry.