



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:38 PM GMT

PDB ID : 1XV8
Title : Crystal Structure of Human Salivary Alpha-Amylase Dimer
Authors : Fisher, S.Z.; Govindasamy, L.; Tu, C.K.; Silverman, D.N.; Rajaniemi, H.; McKenna, R.
Deposited on : 2004-10-27
Resolution : 3.00 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

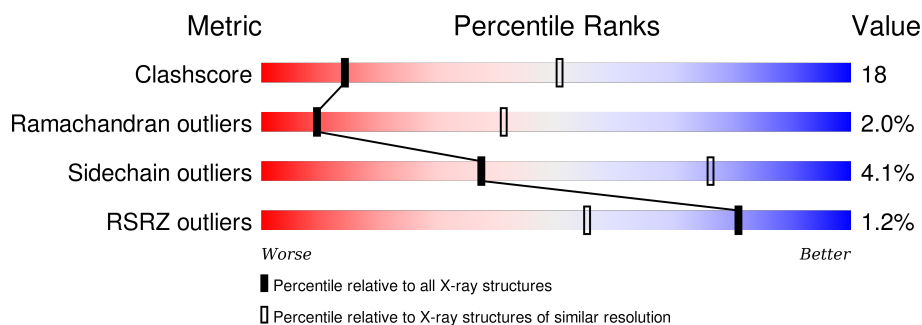
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (3.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	496	<div> <div style="width: 64%;"></div> <div style="width: 32%;"></div> <div style="width: 4%;"></div> </div>
1	B	496	<div> <div style="width: 2%;"></div> <div style="width: 66%;"></div> <div style="width: 31%;"></div> <div style="width: 1%;"></div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	CA	B	497	-	-	-	X

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 8024 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 Alpha-amylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	496	Total	C	N	O	S	0	0	0
			3945	2493	696	735	21			
1	B	496	Total	C	N	O	S	0	0	0
			3945	2493	696	735	21			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	PCA	GLN	MODIFIED RESIDUE	UNP P04745
B	1	PCA	GLN	MODIFIED RESIDUE	UNP P04745

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Ca	0	0
			1	1		
2	A	1	Total	Ca	0	0
			1	1		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Cl	0	0
			1	1		
3	A	1	Total	Cl	0	0
			1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	69	Total	O	0	0
			69	69		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	61	Total	O	0	0
			61	61		

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	150.81Å 72.27Å 91.11Å 90.00° 102.80° 90.00°	Depositor
Resolution (Å)	20.00 – 3.00 19.88 – 3.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-3.00) 96.3 (19.88-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.15	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.46 (at 2.98Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.228 , 0.271 0.259 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	19.3	Xtriage
Anisotropy	0.173	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 17.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	0 of 18583 reflections	Xtriage
F_o, F_c correlation	0.78	EDS
Total number of atoms	8024	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.86% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: CA, PCA, CL

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	0.42	0/4052	0.66	1/5502 (0.0%)
1	B	0.42	0/4052	0.66	1/5502 (0.0%)
All	All	0.42	0/8104	0.66	2/11004 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	57	PRO	CA-N-CD	-5.08	104.39	111.50
1	B	57	PRO	CA-N-CD	-5.04	104.45	111.50

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3945	0	3707	145	4
1	B	3945	0	3707	142	4
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	69	0	0	5	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	61	0	0	5	1
All	All	8024	0	7414	279	5

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 18.

All (279) 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:A:56:ARG:HG3	1:A:56:ARG:HH11	1.16	1.09
1:B:56:ARG:HG3	1:B:56:ARG:HH11	1.16	1.06
1:B:56:ARG:CG	1:B:56:ARG:HH11	1.76	0.97
1:A:56:ARG:CG	1:A:56:ARG:HH11	1.76	0.96
1:B:56:ARG:NH1	1:B:56:ARG:HG3	1.76	0.92
1:A:319:ARG:HH11	1:A:319:ARG:HG2	1.37	0.90
1:B:7:GLN:OE1	1:B:90:GLY:HA2	1.73	0.89
1:B:319:ARG:HG2	1:B:319:ARG:HH11	1.37	0.89
1:B:7:GLN:O	1:B:10:ARG:HG3	1.74	0.88
1:A:7:GLN:OE1	1:A:90:GLY:HA2	1.73	0.87
1:A:7:GLN:O	1:A:10:ARG:HG3	1.74	0.87
1:A:285:GLY:HA3	1:B:152:ASN:ND2	1.90	0.87
1:A:56:ARG:HG3	1:A:56:ARG:NH1	1.76	0.86
1:A:276:TYR:CD1	1:B:149:GLU:HG2	2.13	0.84
1:B:349:GLU:O	1:B:350:ASN:ND2	2.17	0.77
1:A:217:LEU:HD12	1:A:226:SER:HB3	1.67	0.77
1:B:217:LEU:HD12	1:B:226:SER:HB3	1.67	0.77
1:A:285:GLY:CA	1:B:152:ASN:ND2	2.50	0.74
1:A:35:LYS:HD2	1:A:392:ARG:HD3	1.70	0.74
1:B:397:PHE:O	1:B:401:VAL:HG22	1.88	0.74
1:B:28:CYS:HA	1:B:32:LEU:HB2	1.69	0.74
1:A:349:GLU:O	1:A:350:ASN:ND2	2.17	0.74
1:A:397:PHE:O	1:A:401:VAL:HG22	1.88	0.73
1:A:418:ALA:HB2	1:A:428:VAL:HG13	1.71	0.73
1:A:349:GLU:C	1:A:350:ASN:HD22	1.91	0.72
1:B:35:LYS:HD2	1:B:392:ARG:HD3	1.70	0.72
1:B:349:GLU:C	1:B:350:ASN:HD22	1.91	0.72
1:A:28:CYS:HA	1:A:32:LEU:HB2	1.69	0.72
1:B:418:ALA:HB2	1:B:428:VAL:HG13	1.71	0.71
1:A:99:ILE:HB	1:A:179:ILE:HD13	1.73	0.70
1:B:99:ILE:HB	1:B:179:ILE:HD13	1.73	0.70
1:B:319:ARG:NH1	1:B:319:ARG:HG2	2.06	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:4:SER:O	1:B:5:ASN:HB2	1.94	0.66
1:A:4:SER:O	1:A:5:ASN:HB2	1.94	0.66
1:A:319:ARG:NH1	1:A:319:ARG:HG2	2.06	0.66
1:B:380:ASN:O	1:B:381:ASP:HB2	1.96	0.65
1:A:380:ASN:O	1:A:381:ASP:HB2	1.96	0.65
1:A:56:ARG:HD2	1:A:60:GLU:OE1	1.98	0.63
1:B:237:LEU:HD22	1:B:257:LYS:HG2	1.80	0.63
1:B:56:ARG:HD2	1:B:60:GLU:OE1	1.98	0.63
1:B:172:LYS:HB3	1:B:172:LYS:NZ	2.14	0.62
1:A:237:LEU:HD22	1:A:257:LYS:HG2	1.80	0.62
1:B:258:TYR:CE1	1:B:328:MET:HG3	2.35	0.62
1:A:258:TYR:CE1	1:A:328:MET:HG3	2.35	0.61
1:B:11:THR:H	1:B:399:ASN:HD21	1.49	0.61
1:A:172:LYS:NZ	1:A:172:LYS:HB3	2.14	0.61
1:A:293:LEU:HD23	1:A:335:PHE:O	2.01	0.61
1:A:153:ASP:HB3	1:A:156:GLN:HG2	1.83	0.60
1:B:293:LEU:HD23	1:B:335:PHE:O	2.01	0.60
1:B:153:ASP:HB3	1:B:156:GLN:HG2	1.83	0.60
1:B:199:SER:OG	1:B:232:GLN:HB3	2.01	0.60
1:A:224:GLU:OE1	1:A:224:GLU:HA	2.02	0.60
1:A:186:LEU:HB3	1:A:191:VAL:HG21	1.83	0.60
1:B:186:LEU:HB3	1:B:191:VAL:HG21	1.83	0.59
1:A:331:HIS:CG	1:A:332:PRO:HD2	2.37	0.59
1:A:199:SER:OG	1:A:232:GLN:HB3	2.01	0.59
1:A:11:THR:H	1:A:399:ASN:HD21	1.49	0.59
1:A:11:THR:H	1:A:399:ASN:ND2	2.01	0.59
1:A:268:LYS:HD3	1:A:415:ASN:ND2	2.18	0.59
1:B:268:LYS:HD3	1:B:415:ASN:ND2	2.18	0.59
1:B:201:HIS:HB3	4:B:513:HOH:O	2.03	0.59
1:B:15:HIS:HD2	1:B:41:GLN:O	1.85	0.59
1:B:224:GLU:HA	1:B:224:GLU:OE1	2.02	0.59
1:B:331:HIS:CG	1:B:332:PRO:HD2	2.37	0.58
1:A:15:HIS:HD2	1:A:41:GLN:O	1.85	0.58
1:B:11:THR:H	1:B:399:ASN:ND2	2.01	0.58
1:A:201:HIS:HB3	4:A:510:HOH:O	2.03	0.58
1:B:75:ASN:OD1	1:B:77:ASP:HB2	2.05	0.57
1:A:101:HIS:HD2	1:A:102:MET:O	1.88	0.57
1:B:141:CYS:HB2	1:B:159:ASP:O	2.05	0.57
1:A:75:ASN:OD1	1:A:77:ASP:HB2	2.05	0.57
1:A:141:CYS:HB2	1:A:159:ASP:O	2.05	0.57
1:B:101:HIS:HD2	1:B:102:MET:O	1.88	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:31:TYR:OH	1:B:387:ARG:HA	2.06	0.56
1:B:303:ARG:HG3	1:B:303:ARG:HH11	1.71	0.56
1:A:303:ARG:HH11	1:A:303:ARG:HG3	1.71	0.56
1:B:417:VAL:HG22	1:B:429:PHE:HB2	1.88	0.56
1:B:58:TRP:NE1	1:B:356:ASP:O	2.39	0.56
1:A:46:ASN:O	1:A:71:THR:HG22	2.06	0.56
1:A:349:GLU:C	1:A:350:ASN:ND2	2.57	0.55
1:B:46:ASN:O	1:B:71:THR:HG22	2.06	0.55
1:A:58:TRP:NE1	1:A:356:ASP:O	2.39	0.55
1:B:349:GLU:C	1:B:350:ASN:ND2	2.57	0.55
1:A:231:TYR:CD2	1:A:293:LEU:HD12	2.41	0.55
1:B:231:TYR:CD2	1:B:293:LEU:HD12	2.41	0.55
1:A:31:TYR:OH	1:A:387:ARG:HA	2.06	0.55
1:A:54:PRO:HG2	1:A:57:PRO:HG3	1.89	0.55
1:A:417:VAL:HG22	1:A:429:PHE:HB2	1.88	0.55
1:B:54:PRO:HG2	1:B:57:PRO:HG3	1.89	0.54
1:A:217:LEU:HB2	1:A:226:SER:HB2	1.91	0.53
1:B:217:LEU:HB2	1:B:226:SER:HB2	1.91	0.53
1:B:111:THR:O	1:B:113:SER:N	2.42	0.52
1:A:56:ARG:NH2	4:A:535:HOH:O	2.25	0.52
1:A:111:THR:O	1:A:113:SER:N	2.42	0.52
1:A:330:ALA:HB2	1:A:394:MET:CE	2.41	0.51
1:A:470:SER:O	1:A:473:GLY:N	2.35	0.51
1:B:330:ALA:HB2	1:B:394:MET:CE	2.41	0.51
1:A:278:LYS:HE2	1:B:149:GLU:OE1	2.11	0.51
1:A:129:VAL:H	1:A:130:PRO:HA	1.75	0.51
1:B:2:TYR:CE1	1:B:251:GLY:HA2	2.46	0.51
1:B:470:SER:O	1:B:473:GLY:N	2.36	0.51
1:A:2:TYR:CE1	1:A:251:GLY:HA2	2.46	0.51
1:A:153:ASP:OD1	1:A:155:THR:HG22	2.11	0.51
1:B:153:ASP:OD1	1:B:155:THR:HG22	2.11	0.51
1:B:54:PRO:HB2	1:B:357:TRP:CZ3	2.46	0.51
1:B:129:VAL:H	1:B:130:PRO:HA	1.76	0.51
1:A:258:TYR:CZ	1:A:328:MET:HG3	2.46	0.51
1:A:54:PRO:HB2	1:A:357:TRP:CZ3	2.46	0.51
1:A:81:ASN:OD1	1:A:85:ARG:HD3	2.11	0.51
1:A:423:ASN:C	1:A:423:ASN:HD22	2.14	0.50
1:A:95:VAL:HG11	1:A:186:LEU:HD13	1.92	0.50
1:B:423:ASN:HD22	1:B:423:ASN:C	2.14	0.50
1:B:81:ASN:OD1	1:B:85:ARG:HD3	2.11	0.50
1:B:191:VAL:HG12	1:B:193:GLY:H	1.77	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:330:ALA:HB2	1:A:394:MET:HE1	1.94	0.50
1:B:258:TYR:CZ	1:B:328:MET:HG3	2.46	0.49
1:B:353:ASP:HB3	1:B:356:ASP:HB3	1.94	0.49
1:A:362:ASN:HA	1:A:368:LYS:HG3	1.94	0.49
1:B:362:ASN:HA	1:B:368:LYS:HG3	1.94	0.49
1:B:95:VAL:HG11	1:B:186:LEU:HD13	1.92	0.49
1:A:353:ASP:HB3	1:A:356:ASP:HB3	1.94	0.49
1:A:191:VAL:HG12	1:A:193:GLY:H	1.77	0.49
1:B:63:GLN:N	1:B:64:PRO:CD	2.75	0.49
1:A:63:GLN:N	1:A:64:PRO:CD	2.75	0.49
1:B:49:VAL:HG12	1:B:50:ALA:N	2.27	0.49
1:B:20:ARG:HD2	4:B:504:HOH:O	2.12	0.49
1:A:439:THR:HA	1:A:476:HIS:HA	1.95	0.49
1:B:56:ARG:NH1	1:B:56:ARG:CG	2.45	0.49
1:B:208:LYS:HG3	4:B:546:HOH:O	2.13	0.49
1:B:28:CYS:HA	1:B:32:LEU:HD12	1.95	0.49
1:B:129:VAL:HB	1:B:130:PRO:HA	1.94	0.49
1:A:208:LYS:HG3	4:A:545:HOH:O	2.13	0.48
1:A:436:PHE:O	1:A:478:SER:HA	2.13	0.48
1:A:437:SER:O	1:A:438:LEU:HB2	2.13	0.48
1:A:49:VAL:HG12	1:A:50:ALA:N	2.27	0.48
1:B:437:SER:O	1:B:438:LEU:HB2	2.13	0.48
1:B:330:ALA:HB2	1:B:394:MET:HE1	1.95	0.48
1:A:28:CYS:HA	1:A:32:LEU:HD12	1.95	0.48
1:A:293:LEU:C	1:A:293:LEU:HD22	2.33	0.48
1:B:439:THR:HA	1:B:476:HIS:HA	1.95	0.48
1:B:63:GLN:N	1:B:64:PRO:HD3	2.29	0.48
1:A:129:VAL:HB	1:A:130:PRO:HA	1.94	0.48
1:B:436:PHE:O	1:B:478:SER:HA	2.13	0.48
1:B:31:TYR:CE1	1:B:35:LYS:HG3	2.49	0.48
1:A:346:ARG:HG3	1:A:353:ASP:OD2	2.14	0.47
1:A:20:ARG:HD2	4:A:501:HOH:O	2.12	0.47
1:B:124:ARG:HG2	1:B:136:PHE:CD1	2.50	0.47
1:B:346:ARG:HG3	1:B:353:ASP:OD2	2.14	0.47
1:A:319:ARG:CG	1:A:319:ARG:NH1	2.76	0.47
1:B:293:LEU:C	1:B:293:LEU:HD22	2.34	0.47
1:A:63:GLN:N	1:A:64:PRO:HD3	2.29	0.47
1:A:31:TYR:CE1	1:A:35:LYS:HG3	2.49	0.47
1:A:124:ARG:HG2	1:A:136:PHE:CD1	2.50	0.47
1:B:56:ARG:NH2	4:B:537:HOH:O	2.26	0.46
1:A:285:GLY:CA	1:B:152:ASN:HD21	2.27	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:352:LYS:HG2	1:B:353:ASP:N	2.30	0.46
1:A:352:LYS:HG2	1:A:353:ASP:N	2.30	0.46
1:A:150:ASN:C	1:A:152:ASN:H	2.19	0.46
1:A:442:THR:OG1	1:A:444:LEU:HB2	2.15	0.46
1:A:143:THR:HG23	1:A:160:CYS:SG	2.55	0.46
1:B:143:THR:HG23	1:B:160:CYS:SG	2.55	0.46
1:A:56:ARG:HD2	1:A:56:ARG:HA	1.59	0.46
1:A:129:VAL:HB	1:A:130:PRO:CA	2.45	0.46
1:B:183:MET:O	1:B:187:ILE:HG13	2.16	0.46
1:B:129:VAL:HB	1:B:130:PRO:CA	2.45	0.46
1:A:183:MET:O	1:A:187:ILE:HG13	2.16	0.46
1:A:295:PHE:CE2	1:A:298:ASN:HB3	2.51	0.46
1:A:42:VAL:HG22	1:A:94:TYR:O	2.16	0.45
1:B:150:ASN:C	1:B:152:ASN:H	2.19	0.45
1:B:442:THR:OG1	1:B:444:LEU:HB2	2.15	0.45
1:B:42:VAL:HG22	1:B:94:TYR:O	2.16	0.45
1:A:417:VAL:O	1:A:418:ALA:HB2	2.17	0.45
1:B:417:VAL:O	1:B:418:ALA:HB2	2.17	0.45
1:B:295:PHE:CE2	1:B:298:ASN:HB3	2.51	0.45
1:A:153:ASP:O	1:A:155:THR:N	2.50	0.45
1:B:33:ALA:CB	1:B:89:VAL:HB	2.47	0.45
1:A:56:ARG:CG	1:A:56:ARG:NH1	2.45	0.45
1:B:7:GLN:OE1	1:B:90:GLY:CA	2.56	0.45
1:B:153:ASP:O	1:B:155:THR:N	2.50	0.45
1:A:255:GLU:OE1	1:A:257:LYS:HG3	2.17	0.45
1:A:344:TRP:O	1:A:346:ARG:N	2.50	0.45
1:A:20:ARG:HG2	1:A:73:SER:HA	1.98	0.45
1:A:33:ALA:CB	1:A:89:VAL:HB	2.47	0.45
1:B:56:ARG:HA	1:B:56:ARG:HD2	1.59	0.45
1:A:276:TYR:CE1	1:B:149:GLU:HG2	2.51	0.45
1:A:44:PRO:O	1:A:71:THR:HG21	2.17	0.45
1:B:20:ARG:HG2	1:B:73:SER:HA	1.98	0.45
1:A:338:VAL:HG13	4:A:537:HOH:O	2.17	0.45
1:B:344:TRP:O	1:B:346:ARG:N	2.50	0.44
1:B:338:VAL:HG13	4:B:539:HOH:O	2.17	0.44
1:B:44:PRO:O	1:B:71:THR:HG21	2.17	0.44
1:A:42:VAL:CG2	1:A:95:VAL:HA	2.47	0.44
1:A:268:LYS:CD	1:A:415:ASN:ND2	2.80	0.44
1:B:49:VAL:HG23	1:B:63:GLN:O	2.18	0.44
1:A:49:VAL:HG23	1:A:63:GLN:O	2.18	0.44
1:B:287:MET:HG3	1:B:292:ALA:HB2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:287:MET:HG3	1:A:292:ALA:HB2	1.99	0.44
1:B:42:VAL:CG2	1:B:95:VAL:HA	2.48	0.43
1:B:258:TYR:OH	1:B:328:MET:HG3	2.18	0.43
1:B:153:ASP:O	1:B:154:ALA:C	2.57	0.43
1:B:95:VAL:HG11	1:B:186:LEU:CD1	2.48	0.43
1:A:47:GLU:OE2	1:A:115:CYS:HB3	2.18	0.43
1:A:4:SER:O	1:A:5:ASN:CB	2.62	0.43
1:B:255:GLU:OE1	1:B:257:LYS:HG3	2.17	0.43
1:A:258:TYR:OH	1:A:328:MET:HG3	2.18	0.43
1:A:153:ASP:O	1:A:154:ALA:C	2.57	0.43
1:B:252:ARG:NH1	1:B:292:ALA:O	2.50	0.43
1:A:283:GLY:O	1:B:150:ASN:ND2	2.48	0.43
1:A:276:TYR:OH	1:B:147:ASP:HB3	2.18	0.43
1:B:11:THR:OG1	1:B:399:ASN:ND2	2.52	0.43
1:A:11:THR:OG1	1:A:399:ASN:ND2	2.52	0.43
1:A:344:TRP:HB2	1:A:345:PRO:HD2	2.00	0.43
1:A:252:ARG:NH1	1:A:292:ALA:O	2.50	0.43
1:B:430:ASN:HD21	1:B:481:ASN:HA	1.84	0.43
1:B:13:ILE:O	1:B:337:ARG:HA	2.19	0.43
1:A:95:VAL:HG11	1:A:186:LEU:CD1	2.48	0.43
1:B:312:ILE:HD13	1:B:312:ILE:HA	1.82	0.43
1:B:268:LYS:CD	1:B:415:ASN:ND2	2.80	0.42
1:A:64:PRO:HB2	1:A:102:MET:HA	2.01	0.42
1:A:423:ASN:ND2	1:A:423:ASN:C	2.72	0.42
1:A:48:ASN:OD1	1:A:114:THR:HG21	2.19	0.42
1:B:178:LYS:HA	1:B:178:LYS:HD3	1.84	0.42
1:B:414:SER:HB3	1:B:432:ASP:OD2	2.19	0.42
1:A:237:LEU:CD2	1:A:257:LYS:HG2	2.47	0.42
1:B:47:GLU:OE2	1:B:115:CYS:HB3	2.18	0.42
1:A:414:SER:HB3	1:A:432:ASP:OD2	2.19	0.42
1:A:245:SER:HA	1:A:248:PHE:CE2	2.54	0.42
1:B:245:SER:HA	1:B:248:PHE:CE2	2.54	0.42
1:A:13:ILE:O	1:A:337:ARG:HA	2.19	0.42
1:A:9:GLY:O	1:A:11:THR:HG23	2.19	0.42
1:B:344:TRP:HB2	1:B:345:PRO:HD2	2.00	0.42
1:A:244:SER:HB3	1:A:287:MET:HE1	2.02	0.42
1:A:118:TYR:CG	1:A:119:PHE:N	2.88	0.42
1:A:446:ALA:HB2	1:A:471:ASP:N	2.35	0.42
1:B:423:ASN:ND2	1:B:423:ASN:C	2.72	0.42
1:B:86:CYS:O	1:B:89:VAL:HG23	2.20	0.42
1:B:9:GLY:O	1:B:11:THR:HG23	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:64:PRO:HB2	1:B:102:MET:HA	2.01	0.41
1:A:86:CYS:O	1:A:89:VAL:HG23	2.20	0.41
1:B:244:SER:HB3	1:B:287:MET:HE1	2.02	0.41
1:B:48:ASN:OD1	1:B:114:THR:HG21	2.19	0.41
1:A:355:ASN:C	1:A:357:TRP:H	2.24	0.41
1:A:430:ASN:HD21	1:A:481:ASN:HA	1.84	0.41
1:B:4:SER:O	1:B:5:ASN:CB	2.62	0.41
1:B:33:ALA:HB2	1:B:89:VAL:HB	2.02	0.41
1:B:237:LEU:CD2	1:B:257:LYS:HG2	2.47	0.41
1:B:273:LYS:HB2	1:B:415:ASN:OD1	2.21	0.41
1:A:416:GLN:HG3	1:A:432:ASP:OD2	2.21	0.41
1:B:68:LYS:O	1:B:70:CYS:N	2.54	0.41
1:A:14:VAL:HG22	1:A:15:HIS:N	2.36	0.41
1:A:13:ILE:HG23	1:A:335:PHE:CE1	2.56	0.41
1:B:13:ILE:HG23	1:B:335:PHE:CE1	2.56	0.41
1:A:68:LYS:O	1:A:70:CYS:N	2.54	0.41
1:B:446:ALA:HB2	1:B:471:ASP:N	2.35	0.41
1:B:161:ARG:HE	1:B:161:ARG:HB2	1.68	0.41
1:A:347:TYR:O	1:A:354:VAL:HG22	2.20	0.41
1:B:150:ASN:ND2	1:B:152:ASN:HB3	2.36	0.41
1:A:44:PRO:HA	1:A:45:PRO:HD2	2.00	0.41
1:A:45:PRO:O	1:A:69:LEU:HA	2.21	0.41
1:B:118:TYR:CG	1:B:119:PHE:N	2.88	0.41
1:B:319:ARG:CG	1:B:319:ARG:NH1	2.76	0.40
1:A:172:LYS:HZ2	1:A:172:LYS:HB3	1.85	0.40
1:B:347:TYR:O	1:B:354:VAL:HG22	2.20	0.40
1:A:58:TRP:HD1	1:A:357:TRP:O	2.04	0.40
1:A:340:SER:HA	1:A:385:GLU:OE1	2.22	0.40
1:A:4:SER:OG	1:A:6:THR:HG23	2.22	0.40
1:B:355:ASN:C	1:B:357:TRP:H	2.24	0.40
1:B:229:PHE:CE1	1:B:252:ARG:HD2	2.57	0.40
1:B:416:GLN:HG3	1:B:432:ASP:OD2	2.21	0.40
1:A:451:ASP:OD1	1:A:488:ILE:HG23	2.21	0.40
1:A:7:GLN:OE1	1:A:90:GLY:CA	2.56	0.40
1:A:273:LYS:HB2	1:A:415:ASN:OD1	2.21	0.40
1:B:14:VAL:HG22	1:B:15:HIS:N	2.36	0.40
1:A:129:VAL:N	1:A:130:PRO:HA	2.34	0.40
1:A:33:ALA:HB2	1:A:89:VAL:HB	2.02	0.40
1:B:430:ASN:ND2	1:B:481:ASN:HA	2.37	0.40
1:B:8:GLN:HA	1:B:8:GLN:OE1	2.22	0.40
1:A:302:GLN:HB2	1:A:314:THR:CG2	2.52	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:16:LEU:HD12	1:A:24:ILE:HG23	2.04	0.40

All (5) 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:A:220:ASN:O	1:B:8:GLN:OE1[3_445]	1.35	0.85
4:A:499:HOH:O	4:B:547:HOH:O[3_445]	1.62	0.58
1:A:111:THR:OG1	1:B:484:GLU:OE2[2_655]	2.04	0.16
1:A:220:ASN:O	1:B:8:GLN:CD[3_445]	2.13	0.07
1:A:221:TRP:CA	1:B:8:GLN:NE2[3_445]	2.16	0.04

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	494/496 (100%)	427 (86%)	57 (12%)	10 (2%)	9	41
1	B	494/496 (100%)	428 (87%)	56 (11%)	10 (2%)	9	41
All	All	988/992 (100%)	855 (86%)	113 (11%)	20 (2%)	9	41

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	154	ALA
1	A	380	ASN
1	B	154	ALA
1	B	380	ASN
1	A	6	THR
1	A	114	THR
1	B	6	THR
1	B	114	THR

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Mol	Chain	Res	Type
1	A	418	ALA
1	B	418	ALA
1	A	44	PRO
1	A	334	GLY
1	B	44	PRO
1	B	334	GLY
1	A	151	TYR
1	A	345	PRO
1	B	151	TYR
1	B	345	PRO
1	A	271	GLY
1	B	271	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	418/418 (100%)	401 (96%)	17 (4%)	37	76
1	B	418/418 (100%)	401 (96%)	17 (4%)	37	76
All	All	836/836 (100%)	802 (96%)	34 (4%)	37	76

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	VAL
1	A	56	ARG
1	A	77	ASP
1	A	115	CYS
1	A	124	ARG
1	A	125	ASP
1	A	172	LYS
1	A	224	GLU
1	A	290	ASP
1	A	293	LEU
1	A	319	ARG

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Mol	Chain	Res	Type
1	A	323	MET
1	A	341	SER
1	A	350	ASN
1	A	395	VAL
1	A	423	ASN
1	A	442	THR
1	B	22	VAL
1	B	56	ARG
1	B	77	ASP
1	B	115	CYS
1	B	124	ARG
1	B	125	ASP
1	B	172	LYS
1	B	224	GLU
1	B	290	ASP
1	B	293	LEU
1	B	319	ARG
1	B	323	MET
1	B	341	SER
1	B	350	ASN
1	B	395	VAL
1	B	423	ASN
1	B	442	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	HIS
1	A	185	HIS
1	A	216	ASN
1	A	270	ASN
1	A	350	ASN
1	A	399	ASN
1	A	408	ASN
1	A	423	ASN
1	A	459	ASN
1	A	476	HIS
1	B	15	HIS
1	B	152	ASN
1	B	185	HIS
1	B	216	ASN
1	B	270	ASN

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Mol	Chain	Res	Type
1	B	350	ASN
1	B	399	ASN
1	B	408	ASN
1	B	423	ASN
1	B	459	ASN
1	B	476	HIS

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	$\# Z > 2$
1	PCA	A	1	1	7,8,9	2.04	2 (28%)	9,10,12	2.24	3 (33%)
1	PCA	B	1	1	7,8,9	2.04	2 (28%)	9,10,12	2.24	3 (33%)

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
1	PCA	A	1	1	-	0/0/11/13	0/1/1/1
1	PCA	B	1	1	-	0/0/11/13	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	1	PCA	CA-N	3.66	1.51	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1	PCA	CA-N	3.69	1.51	1.46
1	A	1	PCA	CD-N	3.79	1.46	1.33
1	B	1	PCA	CD-N	3.81	1.46	1.33

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1	PCA	CB-CA-C	-4.38	106.77	112.76
1	B	1	PCA	CB-CA-C	-4.37	106.78	112.76
1	B	1	PCA	CA-N-CD	-3.12	103.34	113.81
1	A	1	PCA	CA-N-CD	-3.11	103.37	113.81
1	A	1	PCA	OE-CD-CG	-2.56	121.08	126.81
1	B	1	PCA	OE-CD-CG	-2.56	121.10	126.81

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 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 ⓘ

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	495/496 (99%)	-0.03	2 (0%) 93 80	7, 19, 33, 50	0
1	B	495/496 (99%)	0.42	10 (2%) 68 39	7, 19, 33, 50	0
All	All	990/992 (99%)	0.19	12 (1%) 81 55	7, 19, 33, 50	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	8	GLN	3.1
1	B	163	SER	2.7
1	B	462	CYS	2.7
1	B	364	ASN	2.6
1	B	9	GLY	2.4
1	B	446	ALA	2.4
1	B	90	GLY	2.2
1	B	23	ASP	2.2
1	A	88	ASN	2.2
1	A	134	TRP	2.2
1	B	471	ASP	2.1
1	B	145	SER	2.0

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

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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
1	PCA	B	1	8/9	0.83	0.24	-	23,29,30,32	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
1	PCA	A	1	8/9	0.92	0.20	-	23,29,30,32	0

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. 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	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	CA	B	497	1/1	0.56	0.30	2.08	12,12,12,12	0
2	CA	A	497	1/1	0.79	0.12	-2.25	12,12,12,12	0
3	CL	A	498	1/1	0.95	0.15	-5.94	16,16,16,16	0
3	CL	B	498	1/1	0.94	0.09	-9.80	16,16,16,16	0

6.5 Other polymers [i](#)

There are no such residues in this entry.