



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 21, 2020 – 05:24 PM BST

PDB ID : 6OHB
Title : E. coli Guanine Deaminase
Authors : Shek, R.S.; French, J.B.
Deposited on : 2019-04-05
Resolution : 2.30 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.13.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.13.1

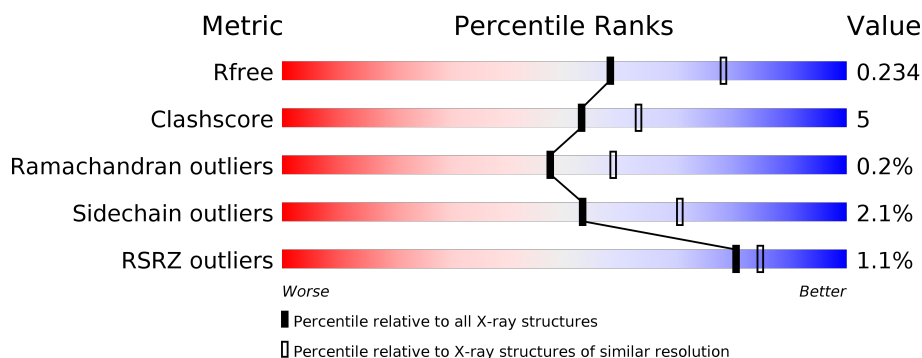
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 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	439	<div> <div>%</div> <div> <div></div> <div>86%</div> <div>12%</div> <div>..</div> </div> </div>
1	B	439	<div> <div></div> <div>87%</div> <div>10%</div> <div>..</div> </div>
1	C	439	<div> <div>%</div> <div> <div></div> <div>83%</div> <div>15%</div> <div>..</div> </div> </div>
1	D	439	<div> <div>2%</div> <div> <div></div> <div>80%</div> <div>17%</div> <div>..</div> </div> </div>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 13904 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 Guanine deaminase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	435	Total	C	N	O	S	0	1	0
			3434	2197	583	637	17			
1	B	435	Total	C	N	O	S	0	0	0
			3446	2204	585	640	17			
1	C	435	Total	C	N	O	S	0	1	0
			3456	2214	589	636	17			
1	D	432	Total	C	N	O	S	0	1	0
			3363	2154	574	618	17			

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		
2	D	1	Total	Zn	0	0
			1	1		
2	C	1	Total	Zn	0	0
			1	1		

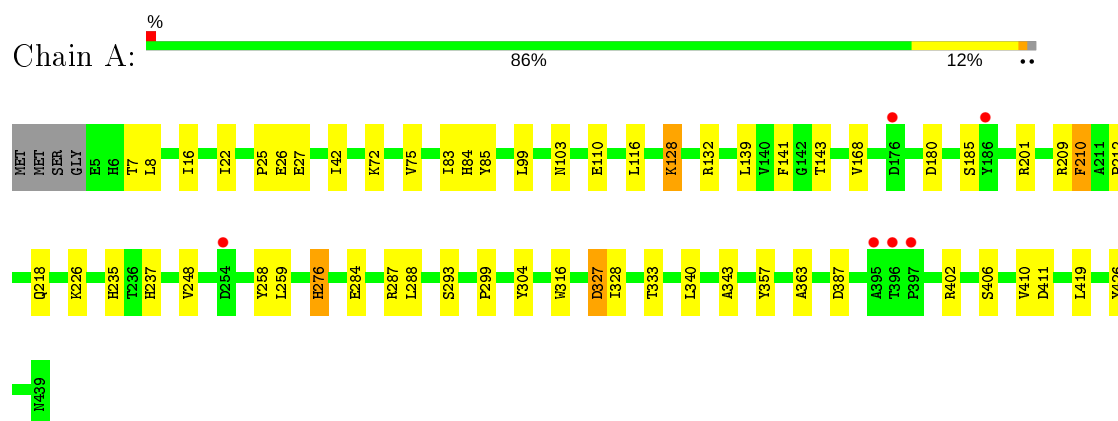
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	61	Total	O	0	0
			61	61		
3	B	65	Total	O	0	0
			65	65		
3	C	68	Total	O	0	0
			68	68		
3	D	7	Total	O	0	0
			7	7		

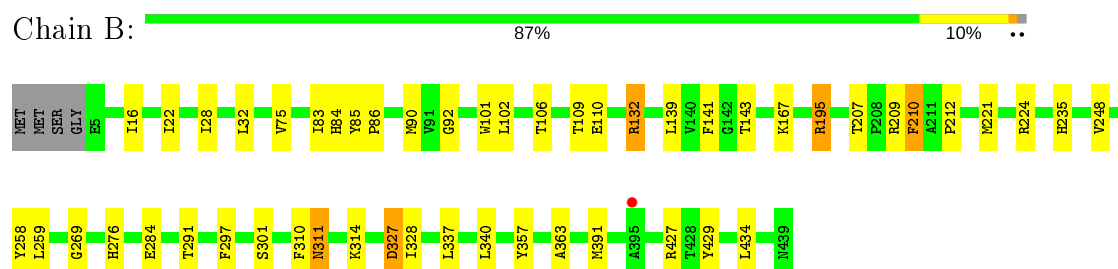
3 Residue-property plots

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.

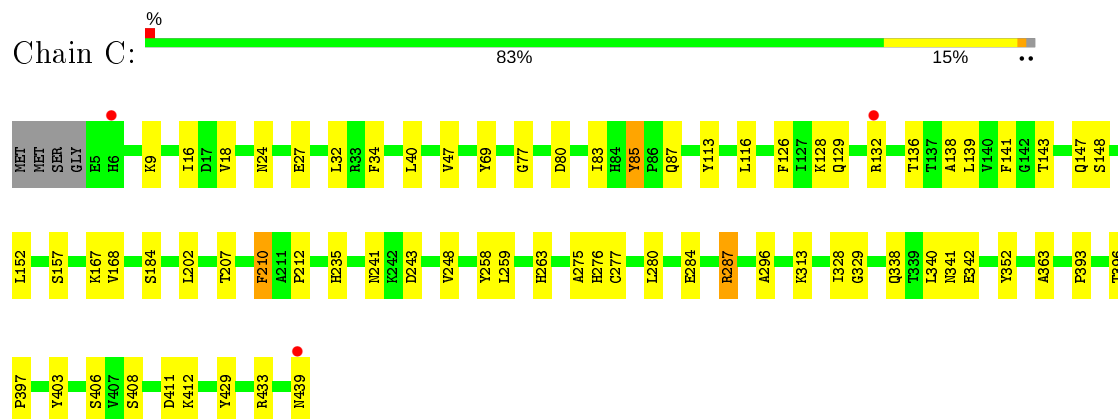
• Molecule 1: Guanine deaminase



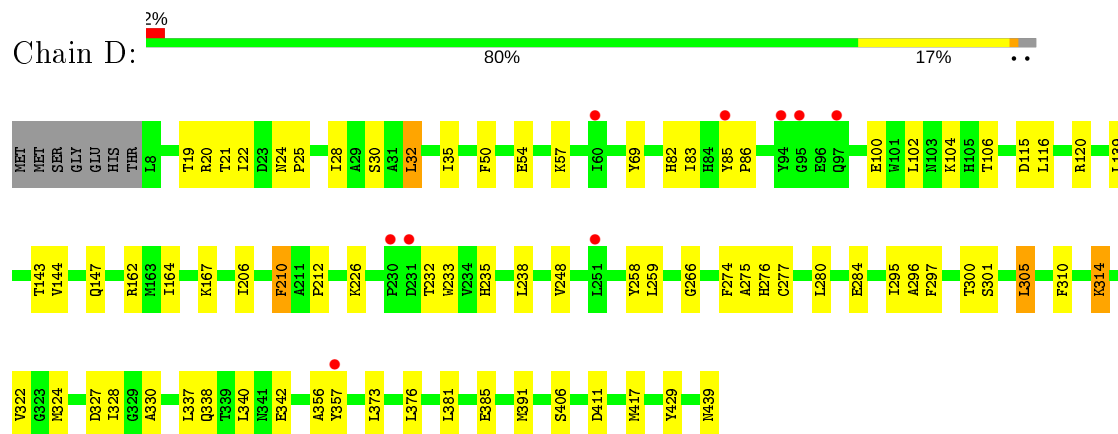
• Molecule 1: Guanine deaminase



• Molecule 1: Guanine deaminase



- Molecule 1: Guanine deaminase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	66.48Å 137.58Å 98.58Å 90.00° 102.16° 90.00°	Depositor
Resolution (Å)	55.99 – 2.30 55.99 – 2.30	Depositor EDS
% Data completeness (in resolution range)	96.0 (55.99-2.30) 91.6 (55.99-2.30)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.62 (at 2.29Å)	Xtriage
Refinement program	PHENIX 1.14_3260, PHENIX 1.14_3260	Depositor
R, R_{free}	0.186 , 0.234 0.185 , 0.234	Depositor DCC
R_{free} test set	3761 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	41.8	Xtriage
Anisotropy	0.357	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 38.4	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	0.96	EDS
Total number of atoms	13904	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.34% 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	A	0.29	0/3521	0.55	3/4779 (0.1%)
1	B	0.30	0/3530	0.54	4/4787 (0.1%)
1	C	0.31	0/3545	0.52	0/4811
1	D	0.28	0/3447	0.53	3/4687 (0.1%)
All	All	0.30	0/14043	0.54	10/19064 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	327	ASP	CB-CG-OD2	13.48	130.44	118.30
1	D	305	LEU	CB-CG-CD2	-8.59	96.39	111.00
1	B	195	ARG	NE-CZ-NH2	-8.21	116.19	120.30
1	B	195	ARG	CB-CG-CD	7.15	130.19	111.60
1	B	195	ARG	CD-NE-CZ	6.54	132.75	123.60
1	A	327	ASP	OD1-CG-OD2	-6.10	111.72	123.30
1	D	314	LYS	CD-CE-NZ	-5.70	98.59	111.70
1	B	327	ASP	CB-CG-OD2	5.55	123.30	118.30
1	D	314	LYS	CG-CD-CE	5.36	127.98	111.90
1	A	116	LEU	CB-CG-CD2	-5.29	102.02	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	195	ARG	Sidechain

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	3434	0	3296	30	0
1	B	3446	0	3338	29	0
1	C	3456	0	3343	41	0
1	D	3363	0	3181	49	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	61	0	0	1	0
3	B	65	0	0	0	0
3	C	68	0	0	1	0
3	D	7	0	0	0	0
All	All	13904	0	13158	146	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 5.

All (146) 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:D:305:LEU:HD11	1:D:330:ALA:HB1	1.42	0.98
1:D:100:GLU:O	1:D:104:LYS:HB2	1.63	0.96
1:C:433:ARG:NH1	3:C:601:HOH:O	2.09	0.85
1:D:54:GLU:HA	1:D:57:LYS:HD2	1.59	0.83
1:C:340:LEU:HD11	1:C:363:ALA:HB2	1.59	0.82
1:B:340:LEU:HD11	1:B:363:ALA:HB2	1.67	0.77
1:B:90:MET:HE1	1:B:101:TRP:HE1	1.51	0.76
1:D:102:LEU:HD23	1:D:106:THR:HG21	1.67	0.74
1:B:84:HIS:HD2	1:B:327:ASP:OD1	1.70	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:305:LEU:CD1	1:D:330:ALA:HB1	2.16	0.73
1:B:427:ARG:HD2	1:B:434:LEU:HD11	1.70	0.73
1:B:83:ILE:HG13	1:B:328:ILE:HB	1.69	0.72
1:D:210:PHE:CD1	1:D:212:PRO:HD2	2.25	0.72
1:A:83:ILE:HG13	1:A:328:ILE:HB	1.71	0.71
1:D:102:LEU:HA	1:D:106:THR:HG22	1.72	0.71
1:D:83:ILE:HG13	1:D:328:ILE:HB	1.73	0.70
1:B:90:MET:HE1	1:B:101:TRP:NE1	2.06	0.70
1:C:83:ILE:HG13	1:C:328:ILE:HB	1.74	0.70
1:C:259:LEU:HB3	1:C:284:GLU:HG2	1.73	0.69
1:D:324:MET:HG3	1:D:340:LEU:HD21	1.75	0.68
1:D:162:ARG:HD3	1:D:373:LEU:HD22	1.73	0.68
1:A:210:PHE:CE2	1:A:212:PRO:HB2	2.30	0.66
1:D:210:PHE:CE1	1:D:212:PRO:HB2	2.34	0.63
1:A:210:PHE:CD2	1:A:212:PRO:HD2	2.35	0.62
1:A:84:HIS:HD2	1:A:327:ASP:OD1	1.80	0.60
1:D:301:SER:O	1:D:305:LEU:HB2	2.02	0.60
1:C:406:SER:HB3	1:C:411:ASP:HB3	1.86	0.58
1:C:275:ALA:HA	1:C:296:ALA:HB3	1.86	0.57
1:D:22:ILE:HG22	1:D:24:ASN:H	1.69	0.57
1:A:7:THR:HG21	1:D:115:ASP:OD2	2.04	0.57
1:A:340:LEU:HD11	1:A:363:ALA:HB2	1.86	0.57
1:B:92:GLY:HA3	1:B:101:TRP:CH2	2.40	0.57
1:C:116:LEU:HD13	1:C:147:GLN:HB3	1.86	0.56
1:D:144:VAL:O	1:D:167:LYS:NZ	2.33	0.56
1:D:25:PRO:HA	1:D:28:ILE:HD12	1.87	0.56
1:C:338:GLN:O	1:C:342:GLU:HG2	2.05	0.56
1:A:110:GLU:OE1	1:A:209:ARG:NH2	2.36	0.56
1:B:90:MET:HG3	1:B:109:THR:HG21	1.88	0.56
1:C:129:GLN:OE1	1:C:132[B]:ARG:NH1	2.39	0.55
1:B:212:PRO:HD3	1:B:248:VAL:HG22	1.88	0.54
1:A:210:PHE:HE2	1:A:212:PRO:HB2	1.73	0.53
1:D:210:PHE:HE1	1:D:212:PRO:HB2	1.73	0.53
1:D:259:LEU:HB3	1:D:284:GLU:HG2	1.90	0.53
1:C:18:VAL:HG22	1:C:32:LEU:HD12	1.91	0.53
1:D:32:LEU:HD13	1:D:357:TYR:HD1	1.74	0.53
1:A:180:ASP:OD1	1:A:218:GLN:NE2	2.41	0.53
1:D:373:LEU:HD23	1:D:376:LEU:HD12	1.92	0.52
1:D:275:ALA:HA	1:D:296:ALA:HB3	1.91	0.52
1:C:143:THR:OG1	1:C:148:SER:HB2	2.10	0.52
1:D:102:LEU:HA	1:D:106:THR:CG2	2.37	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:162:ARG:NH2	1:D:385:GLU:OE2	2.43	0.51
1:A:212:PRO:HD3	1:A:248:VAL:HG22	1.92	0.50
1:A:237:HIS:HD2	1:A:276:HIS:CD2	2.30	0.50
1:A:25:PRO:HD3	1:A:316:TRP:CD2	2.47	0.49
1:C:24:ASN:HD21	1:C:27:GLU:HG3	1.77	0.49
1:D:206:ILE:HD12	1:D:233:TRP:O	2.13	0.49
1:A:406:SER:HB3	1:A:411:ASP:HB3	1.94	0.49
1:D:86:PRO:HG3	1:D:143:THR:HG23	1.94	0.49
1:B:259:LEU:HB3	1:B:284:GLU:HG2	1.95	0.48
1:C:403:TYR:CE1	1:C:412:LYS:HD2	2.47	0.48
1:A:99:LEU:O	1:A:103:ASN:ND2	2.47	0.48
1:D:212:PRO:HD3	1:D:248:VAL:HG22	1.95	0.48
1:B:110:GLU:OE1	1:B:209:ARG:NH2	2.45	0.48
1:D:139:LEU:HG	1:D:235:HIS:CE1	2.49	0.48
1:D:297:PHE:CE1	1:D:310:PHE:HB3	2.50	0.47
1:B:16:ILE:HD11	1:B:32:LEU:HD11	1.95	0.47
1:C:69:TYR:OH	1:C:429:TYR:OH	2.19	0.47
1:A:22:ILE:HG21	1:A:27:GLU:HB2	1.97	0.47
1:C:24:ASN:ND2	1:C:27:GLU:HG3	2.29	0.47
1:C:408:SER:O	1:C:412:LYS:HG2	2.15	0.47
1:C:157:SER:HB2	1:C:202:LEU:HD11	1.96	0.46
1:A:402[B]:ARG:NH1	1:A:419:LEU:HD11	2.31	0.46
1:A:139:LEU:HG	1:A:235:HIS:CE1	2.49	0.46
1:B:16:ILE:HG23	1:B:75:VAL:HG12	1.98	0.46
1:C:16:ILE:HG12	1:C:34:PHE:HD1	1.81	0.46
1:D:277:CYS:HB3	1:D:280:LEU:HD11	1.96	0.46
1:B:337:LEU:HD11	1:B:391:MET:HE3	1.97	0.46
1:C:212:PRO:HD3	1:C:248:VAL:HG22	1.96	0.46
1:C:403:TYR:CZ	1:C:412:LYS:HD2	2.51	0.46
1:B:139:LEU:HG	1:B:235:HIS:CE1	2.51	0.46
1:A:72:LYS:HE2	1:A:426:TYR:CD2	2.51	0.45
1:D:210:PHE:CD1	1:D:212:PRO:CD	2.98	0.45
1:D:21:THR:C	1:D:22:ILE:HG13	2.37	0.45
1:D:338:GLN:O	1:D:342:GLU:HG2	2.16	0.45
1:B:311:ASN:ND2	1:B:314:LYS:H	2.15	0.45
1:D:300:THR:HG22	1:D:342:GLU:HG3	1.98	0.45
1:B:22:ILE:HG12	1:B:28:ILE:HG12	1.99	0.45
1:D:164:ILE:HD11	1:D:373:LEU:HD11	1.99	0.45
1:C:87:GLN:OE1	1:C:329:GLY:HA3	2.17	0.45
1:A:201:ARG:NH2	1:A:387:ASP:OD2	2.35	0.45
1:A:287:ARG:HG2	1:A:287:ARG:HH11	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:20[A]:ARG:NH2	1:D:30:SER:O	2.50	0.44
1:D:337:LEU:HD11	1:D:391:MET:HE3	1.99	0.44
1:C:128:LYS:HA	1:C:128:LYS:HD3	1.81	0.44
1:D:406:SER:HB3	1:D:411:ASP:HB3	1.99	0.44
1:C:393:PRO:HB2	1:C:403:TYR:CD1	2.53	0.44
1:B:221:MET:HE1	1:B:224:ARG:HH11	1.83	0.44
1:B:32:LEU:HD22	1:B:357:TYR:HD2	1.82	0.44
1:C:167:LYS:NZ	1:C:168:VAL:O	2.50	0.44
1:C:40:LEU:HD11	1:C:47:VAL:HG22	1.99	0.44
1:C:77:GLY:HA3	1:C:136:THR:OG1	2.18	0.44
1:B:102:LEU:HA	1:B:106:THR:OG1	2.18	0.43
1:B:132:ARG:HH22	1:C:439:ASN:CB	2.32	0.43
1:A:8:LEU:HD12	1:A:42:ILE:O	2.18	0.43
1:C:113:TYR:O	1:C:148:SER:OG	2.27	0.43
1:B:297:PHE:CE1	1:B:310:PHE:HB3	2.54	0.43
1:B:86:PRO:HG3	1:B:143:THR:HG23	2.01	0.42
1:D:35:ILE:HD12	1:D:50:PHE:CZ	2.54	0.42
1:B:210:PHE:CD1	1:B:212:PRO:HD2	2.54	0.42
1:B:167:LYS:O	1:B:207:THR:HG23	2.19	0.42
1:D:82:HIS:CD2	1:D:327:ASP:OD1	2.71	0.42
1:C:168:VAL:HA	1:C:207:THR:O	2.19	0.42
1:D:69:TYR:OH	1:D:429:TYR:OH	2.24	0.42
1:C:85:TYR:HA	1:C:126:PHE:CE1	2.55	0.42
1:A:16:ILE:HG23	1:A:75:VAL:HG12	2.00	0.42
1:A:259:LEU:HB3	1:A:284:GLU:HG2	2.01	0.42
1:B:210:PHE:CE1	1:B:212:PRO:HB2	2.55	0.42
1:B:337:LEU:HD11	1:B:391:MET:CE	2.50	0.42
1:C:210:PHE:CD2	1:C:212:PRO:HD2	2.55	0.42
1:D:295:ILE:HB	1:D:322:VAL:HG22	2.01	0.42
1:A:128:LYS:HD3	3:A:620:HOH:O	2.18	0.42
1:C:85:TYR:CD1	1:C:152:LEU:HD13	2.55	0.42
1:D:206:ILE:HD11	1:D:232:THR:OG1	2.20	0.42
1:D:226:LYS:NZ	1:D:266:GLY:O	2.41	0.41
1:A:299:PRO:HG3	1:A:343:ALA:HB2	2.02	0.41
1:C:277:CYS:HB3	1:C:280:LEU:HD11	2.01	0.41
1:D:25:PRO:HA	1:D:28:ILE:CD1	2.50	0.41
1:C:139:LEU:HG	1:C:235:HIS:CE1	2.56	0.41
1:A:143:THR:O	1:A:168:VAL:N	2.37	0.41
1:B:429:TYR:CE1	1:B:434:LEU:HD13	2.55	0.41
1:D:238:LEU:HB3	1:D:274:PHE:HB3	2.03	0.41
1:C:396:THR:HG23	1:C:397:PRO:HD2	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:328:ILE:HD11	1:A:333:THR:C	2.41	0.41
1:C:263:HIS:CG	1:C:287:ARG:HD3	2.55	0.41
1:D:116:LEU:O	1:D:120:ARG:HG3	2.21	0.41
1:A:288:LEU:HD22	1:A:293:SER:HB2	2.02	0.41
1:D:25:PRO:O	1:D:28:ILE:HD12	2.20	0.41
1:C:80:ASP:O	1:C:138:ALA:HA	2.21	0.41
1:C:313:LYS:HB2	1:C:352:TYR:CD1	2.56	0.41
1:D:19:THR:O	1:D:381:LEU:HD11	2.22	0.40
1:B:269:GLY:HA2	1:B:291:THR:O	2.21	0.40
1:C:9:LYS:HB2	1:C:9:LYS:HE3	1.86	0.40
1:A:304:TYR:OH	1:C:341:ASN:ND2	2.41	0.40
1:A:357:TYR:CE1	1:A:410:VAL:HG12	2.56	0.40
1:C:241:ASN:OD1	1:C:243:ASP:N	2.55	0.40
1:D:356:ALA:HA	1:D:417:MET:HE1	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	434/439 (99%)	423 (98%)	10 (2%)	1 (0%)	47	58
1	B	433/439 (99%)	420 (97%)	12 (3%)	1 (0%)	47	58
1	C	434/439 (99%)	419 (96%)	14 (3%)	1 (0%)	47	58
1	D	431/439 (98%)	417 (97%)	13 (3%)	1 (0%)	47	58
All	All	1732/1756 (99%)	1679 (97%)	49 (3%)	4 (0%)	47	58

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	276	HIS

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Mol	Chain	Res	Type
1	B	276	HIS
1	C	276	HIS
1	D	276	HIS

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	354/379 (93%)	345 (98%)	9 (2%)	47	65
1	B	360/379 (95%)	353 (98%)	7 (2%)	57	73
1	C	359/379 (95%)	353 (98%)	6 (2%)	60	76
1	D	336/379 (89%)	329 (98%)	7 (2%)	53	70
All	All	1409/1516 (93%)	1380 (98%)	29 (2%)	53	70

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	GLU
1	A	85	TYR
1	A	128	LYS
1	A	132	ARG
1	A	141	PHE
1	A	185	SER
1	A	210	PHE
1	A	226	LYS
1	A	258	TYR
1	B	85	TYR
1	B	132	ARG
1	B	141	PHE
1	B	210	PHE
1	B	258	TYR
1	B	301	SER
1	B	311	ASN
1	C	85	TYR
1	C	141	PHE

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Mol	Chain	Res	Type
1	C	184	SER
1	C	210	PHE
1	C	258	TYR
1	C	287	ARG
1	D	32	LEU
1	D	85	TYR
1	D	147	GLN
1	D	210	PHE
1	D	258	TYR
1	D	314	LYS
1	D	439	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	311	ASN
1	C	24	ASN
1	D	59	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 monosaccharides 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 [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	435/439 (99%)	0.07	6 (1%) 75 80	30, 48, 67, 83	0
1	B	435/439 (99%)	0.07	1 (0%) 95 96	28, 45, 66, 82	0
1	C	435/439 (99%)	-0.06	3 (0%) 87 91	30, 45, 59, 82	0
1	D	432/439 (98%)	0.30	9 (2%) 63 70	47, 63, 77, 92	0
All	All	1737/1756 (98%)	0.09	19 (1%) 80 85	28, 50, 72, 92	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	94	TYR	4.5
1	C	439	ASN	3.7
1	A	254	ASP	3.4
1	D	95	GLY	3.3
1	D	60	ILE	3.0
1	A	186	TYR	2.9
1	D	357	TYR	2.8
1	C	132[A]	ARG	2.7
1	D	231	ASP	2.5
1	A	395	ALA	2.5
1	C	6	HIS	2.5
1	A	176	ASP	2.4
1	D	97	GLN	2.4
1	B	395	ALA	2.3
1	A	396	THR	2.2
1	A	397	PRO	2.2
1	D	230	PRO	2.2
1	D	251	LEU	2.1
1	D	85	TYR	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 monosaccharides 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(\AA^2)	Q<0.9
2	ZN	A	501	1/1	0.83	0.11	51,51,51,51	0
2	ZN	B	501	1/1	0.92	0.09	46,46,46,46	0
2	ZN	C	501	1/1	0.98	0.09	38,38,38,38	0
2	ZN	D	501	1/1	0.98	0.08	54,54,54,54	0

6.5 Other polymers [i](#)

There are no such residues in this entry.