



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 06:04 AM UTC

PDB ID : 4HST / pdb_00004hst
Title : Crystal structure of a double mutant of a class III engineered cephalosporin acylase
Authors : Vrielink, A.; Golden, E.; Patterson, R.; Tie, W.J.; Anandan, A.; Flematti, G.; Molla, G.; Rosini, E.; Pollegioni, L.
Deposited on : 2012-10-30
Resolution : 1.57 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

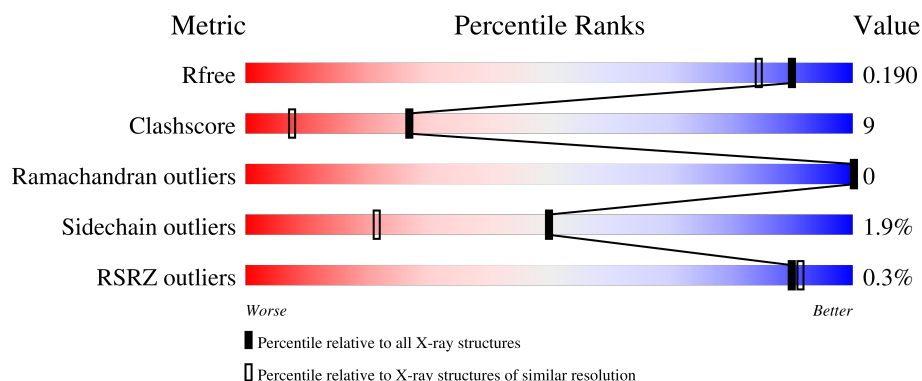
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.57 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	1094 (1.58-1.58)
Clashscore	190562	1105 (1.58-1.58)
Ramachandran outliers	187476	1082 (1.58-1.58)
Sidechain outliers	187428	1081 (1.58-1.58)
RSRZ outliers	180081	1094 (1.58-1.58)

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	229	
2	B	543	

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
3	GLJ	B	601	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6863 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.

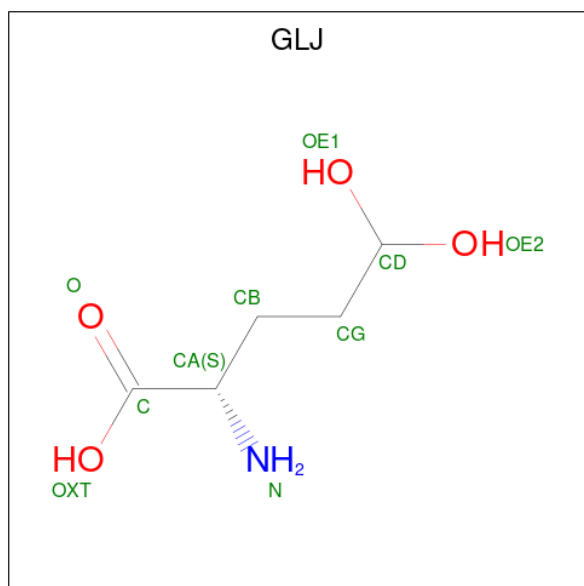
- Molecule 1 is a protein called glutaryl-7-aminocephalosporanic acid acylase alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	216	1686	1082	300	293	11	0	9	0

- Molecule 2 is a protein called glutaryl-7-aminocephalosporanic acid acylase beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	535	4191	2647	765	763	16	0	17	0

- Molecule 3 is 5,5-dihydroxy-L-norvaline (CCD ID: GLJ) (formula: $C_5H_{11}NO_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
3	B	1	10	5	1	4	0	0

- Molecule 4 is water.

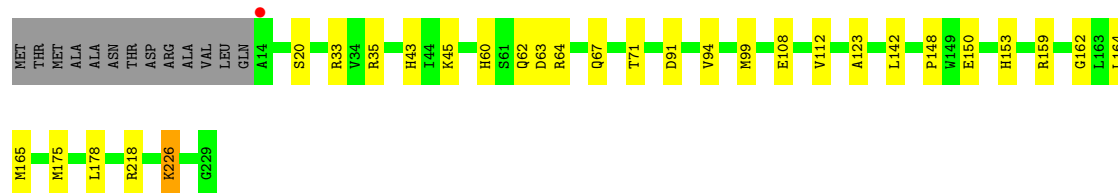
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	290	Total 290	O 290	0	0
4	B	686	Total 686	O 686	0	0

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.

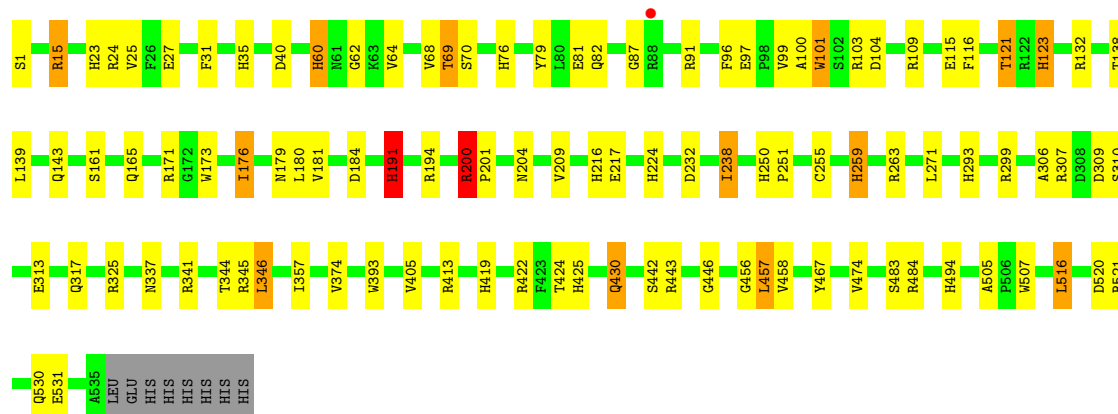
- Molecule 1: glutaryl-7-aminocephalosporanic acid acylase alpha chain

Chain A: 



- Molecule 2: glutaryl-7-aminocephalosporanic acid acylase beta chain

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	68.39Å 77.84Å 191.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	96.00 – 1.57 96.00 – 1.57	Depositor EDS
% Data completeness (in resolution range)	93.4 (96.00-1.57) 93.4 (96.00-1.57)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.51 (at 1.57Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.119 , 0.162 (Not available) , 0.190	Depositor DCC
R_{free} test set	6723 reflections (4.69%)	wwPDB-VP
Wilson B-factor (Å ²)	13.9	Xtriage
Anisotropy	0.204	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 49.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	6863	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.04% 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: GLJ

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	1.65	10/1753 (0.6%)	1.23	2/2375 (0.1%)
2	B	1.71	42/4366 (1.0%)	1.27	10/5968 (0.2%)
All	All	1.69	52/6119 (0.8%)	1.25	12/8343 (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
2	B	3	0

All (52) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	259	HIS	ND1-CE1	10.93	1.43	1.32
2	B	60	HIS	ND1-CE1	9.38	1.42	1.32
2	B	425	HIS	CE1-NE2	8.70	1.41	1.32
2	B	293	HIS	CE1-NE2	8.32	1.40	1.32
2	B	101	TRP	CD2-CE2	8.23	1.55	1.41
2	B	531	GLU	N-CA	8.23	1.56	1.46
2	B	200	ARG	N-CA	8.10	1.53	1.46
2	B	232	ASP	C-O	7.97	1.27	1.23
2	B	474	VAL	CA-CB	7.97	1.65	1.54
2	B	443	ARG	CZ-NH1	7.47	1.43	1.32
1	A	91	ASP	CG-OD1	7.40	1.39	1.25
2	B	430	GLN	CD-OE1	6.80	1.36	1.23
2	B	419	HIS	ND1-CE1	6.80	1.39	1.32
1	A	64	ARG	N-CA	6.77	1.53	1.46
2	B	191	HIS	ND1-CE1	6.71	1.39	1.32
1	A	178	LEU	N-CA	6.60	1.52	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	232	ASP	CA-C	6.53	1.58	1.53
2	B	15	ARG	CD-NE	6.43	1.55	1.46
1	A	45	LYS	C-O	6.24	1.31	1.24
2	B	87	GLY	C-N	-6.22	1.24	1.33
2	B	194	ARG	NE-CZ	6.22	1.39	1.33
2	B	259	HIS	CE1-NE2	6.18	1.38	1.32
2	B	123	HIS	ND1-CE1	6.17	1.38	1.32
2	B	458	VAL	N-CA	6.12	1.51	1.46
2	B	344	THR	N-CA	6.07	1.53	1.46
1	A	63	ASP	CG-OD2	-6.04	1.13	1.25
1	A	60	HIS	CE1-NE2	6.01	1.38	1.32
2	B	76	HIS	ND1-CE1	5.86	1.38	1.32
1	A	153	HIS	CE1-NE2	5.79	1.38	1.32
2	B	259	HIS	CG-CD2	5.73	1.42	1.35
1	A	43	HIS	ND1-CE1	5.71	1.38	1.32
2	B	100	ALA	N-CA	5.67	1.53	1.46
2	B	516	LEU	C-O	-5.54	1.17	1.24
2	B	60	HIS	CE1-NE2	5.50	1.38	1.32
2	B	255	CYS	N-CA	5.46	1.52	1.46
2	B	68	VAL	N-CA	5.45	1.52	1.46
2	B	374	VAL	CA-C	5.44	1.59	1.52
2	B	23	HIS	CE1-NE2	5.40	1.38	1.32
2	B	457	LEU	CB-CG	5.31	1.64	1.53
2	B	76	HIS	CG-CD2	5.30	1.41	1.35
2	B	357	ILE	CA-C	5.28	1.58	1.52
2	B	194	ARG	CA-CB	5.24	1.59	1.52
2	B	180	LEU	CA-CB	5.18	1.59	1.53
2	B	216	HIS	ND1-CE1	5.16	1.37	1.32
2	B	69	THR	C-O	5.14	1.30	1.23
2	B	446	GLY	C-O	5.08	1.28	1.24
2	B	209	VAL	CA-C	5.06	1.56	1.52
2	B	494	HIS	ND1-CE1	5.06	1.37	1.32
2	B	505	ALA	CA-C	5.05	1.58	1.52
2	B	507	TRP	NE1-CE2	5.05	1.43	1.37
1	A	165	MET	C-O	5.04	1.30	1.23
1	A	45	LYS	N-CA	5.04	1.52	1.46

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	271	LEU	N-CA-C	6.78	118.75	111.36
2	B	405	VAL	N-CA-C	-6.56	104.36	110.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	443	ARG	NE-CZ-NH2	-6.04	113.77	119.20
2	B	1	SER	CA-C-O	-5.40	111.62	120.80
1	A	123	ALA	N-CA-C	-5.37	105.43	111.28
2	B	97	GLU	CB-CA-C	5.37	117.12	108.91
2	B	309	ASP	N-CA-CB	-5.34	101.49	109.71
2	B	251	PRO	N-CA-C	5.32	120.83	113.65
2	B	309	ASP	CB-CA-C	5.15	117.99	110.26
1	A	165	MET	CG-SD-CE	-5.13	89.61	100.90
2	B	121	THR	CA-CB-OG1	5.11	117.27	109.60
2	B	191	HIS	CB-CG-CD2	-5.02	124.67	131.20

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	B	121	THR	CB
2	B	138	THR	CB
2	B	238[A]	ILE	CB

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	1686	0	1753	21	0
2	B	4191	0	4056	90	0
3	B	10	0	9	10	0
4	A	290	0	0	9	1
4	B	686	0	0	25	1
All	All	6863	0	5818	108	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 9.

All (108) 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)
2:B:520:ASP:HB2	4:B:795:HOH:O	1.32	1.26
4:A:358:HOH:O	2:B:176[A]:ILE:HD11	1.34	1.26
2:B:109[B]:ARG:NH1	2:B:430:GLN:HE22	1.49	1.10
2:B:109[B]:ARG:CZ	2:B:430:GLN:HE22	1.64	1.09
2:B:109[B]:ARG:NH1	2:B:430:GLN:NE2	2.10	0.98
2:B:40:ASP:HB3	4:B:1274:HOH:O	1.65	0.95
1:A:142:LEU:HD23	2:B:109[B]:ARG:NH1	1.83	0.93
2:B:121:THR:HG23	2:B:123:HIS:H	1.35	0.91
2:B:15:ARG:NH2	4:B:1055:HOH:O	2.02	0.91
2:B:81:GLU:OE2	2:B:121:THR:HG21	1.69	0.91
1:A:62:GLN:HE22	2:B:530:GLN:HE22	1.20	0.90
2:B:60:HIS:HD2	2:B:62:GLY:H	1.18	0.88
2:B:346:LEU:HD21	2:B:393:TRP:HZ3	1.38	0.87
1:A:142:LEU:CD2	2:B:109[B]:ARG:NH1	2.37	0.86
2:B:109[B]:ARG:CZ	2:B:430:GLN:NE2	2.36	0.86
1:A:142:LEU:HD23	2:B:109[B]:ARG:HH12	1.40	0.85
2:B:99[B]:VAL:CG1	2:B:101:TRP:CD1	2.61	0.84
2:B:70[B]:SER:H	3:B:601:GLJ:HG3	1.43	0.83
1:A:67:GLN:O	1:A:71[A]:THR:HG22	1.79	0.83
2:B:70[A]:SER:H	3:B:601:GLJ:HG3	1.45	0.81
2:B:299:ARG:NH2	4:B:1119:HOH:O	2.13	0.81
2:B:413:ARG:NH2	4:B:1116:HOH:O	2.01	0.80
2:B:299:ARG:HD3	4:B:779:HOH:O	1.81	0.79
2:B:346:LEU:HD21	2:B:393:TRP:CZ3	2.17	0.79
1:A:164:LEU:H	2:B:143:GLN:HE22	1.30	0.79
2:B:132:ARG:HD2	4:B:1326:HOH:O	1.85	0.76
2:B:181:VAL:HG22	2:B:238[A]:ILE:HD11	1.68	0.76
2:B:345:ARG:HD2	4:B:1296:HOH:O	1.87	0.75
2:B:99[B]:VAL:HG11	2:B:101:TRP:NE1	2.03	0.73
2:B:346:LEU:CD2	2:B:393:TRP:HZ3	2.01	0.73
2:B:70[B]:SER:H	3:B:601:GLJ:CG	2.01	0.73
2:B:109[A]:ARG:HD3	2:B:430:GLN:HE22	1.55	0.71
2:B:70[A]:SER:H	3:B:601:GLJ:CG	2.03	0.71
1:A:150:GLU:HG2	4:A:442:HOH:O	1.90	0.70
2:B:250:HIS:HE1	4:B:794:HOH:O	1.76	0.68
2:B:99[B]:VAL:HG11	2:B:101:TRP:HE1	1.57	0.68
2:B:123:HIS:HE1	2:B:217:GLU:OE2	1.78	0.67
2:B:307:ARG:N	2:B:307:ARG:HD2	2.10	0.66
1:A:226:LYS:NZ	4:A:521:HOH:O	2.29	0.66
2:B:27[B]:GLU:OE2	2:B:31:PHE:HE1	1.78	0.65
2:B:337:ASN:ND2	2:B:341:ARG:HE	1.97	0.62
2:B:521:ARG:HD3	4:B:900:HOH:O	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:60:HIS:HD2	2:B:62:GLY:N	1.96	0.62
2:B:99[B]:VAL:CG1	2:B:101:TRP:HD1	2.13	0.61
2:B:99[B]:VAL:HG13	2:B:101:TRP:CD1	2.34	0.61
2:B:173:TRP:NE1	4:B:1376:HOH:O	2.34	0.61
1:A:71[A]:THR:HG21	4:B:862:HOH:O	2.01	0.60
2:B:413:ARG:HD2	4:B:1176:HOH:O	2.00	0.60
2:B:24[B]:ARG:HE	3:B:601:GLJ:H	1.49	0.60
1:A:164:LEU:HD21	2:B:139:LEU:HD21	1.84	0.59
1:A:148:PRO:HB3	4:A:539:HOH:O	2.01	0.59
2:B:161:SER:H	2:B:165:GLN:NE2	2.00	0.59
2:B:81:GLU:OE2	2:B:138:THR:HG21	2.03	0.59
1:A:218:ARG:HD2	4:A:569:HOH:O	2.03	0.58
2:B:224:HIS:HE1	4:B:842:HOH:O	1.85	0.58
2:B:259:HIS:HD2	4:B:1146:HOH:O	1.87	0.58
2:B:99[B]:VAL:HG11	2:B:101:TRP:CD1	2.38	0.57
1:A:67:GLN:O	1:A:71[B]:THR:HG23	2.03	0.57
2:B:81:GLU:OE2	2:B:123:HIS:HD2	1.87	0.57
2:B:191:HIS:HB2	2:B:238[A]:ILE:CD1	2.34	0.57
2:B:109[B]:ARG:NH2	4:B:854:HOH:O	2.39	0.55
4:A:306:HOH:O	2:B:35:HIS:HE1	1.89	0.55
3:B:601:GLJ:HA	3:B:601:GLJ:OE1	2.06	0.55
2:B:121:THR:HG23	2:B:123:HIS:N	2.13	0.54
2:B:313:GLU:O	2:B:317:GLN:HG3	2.07	0.54
2:B:161:SER:H	2:B:165:GLN:HE21	1.55	0.54
1:A:112[A]:VAL:HG13	4:A:451:HOH:O	2.08	0.53
1:A:162:GLY:HA3	2:B:176[A]:ILE:CD1	2.38	0.53
2:B:60:HIS:HE1	4:B:707:HOH:O	1.90	0.53
2:B:69:THR:HA	3:B:601:GLJ:HG3	1.91	0.52
2:B:103:ARG:HH21	2:B:115[B]:GLU:CD	2.17	0.52
2:B:171:ARG:HB2	2:B:171:ARG:NH1	2.25	0.51
2:B:179:ASN:OD1	2:B:191:HIS:HE1	1.94	0.49
2:B:306:ALA:C	2:B:307:ARG:HD2	2.37	0.49
2:B:346:LEU:C	2:B:346:LEU:HD23	2.38	0.49
2:B:82[A]:GLN:HE22	2:B:91:ARG:HH11	1.61	0.49
2:B:325:ARG:NH2	4:B:978:HOH:O	2.46	0.49
2:B:109[A]:ARG:NH2	4:B:917:HOH:O	2.42	0.48
1:A:33:ARG:HD2	1:A:35:ARG:NH2	2.29	0.47
2:B:456:GLY:HA3	2:B:467:TYR:CZ	2.50	0.47
2:B:25[A]:VAL:HG13	2:B:27[A]:GLU:HG3	1.97	0.46
2:B:259:HIS:CD2	2:B:263:ARG:HH21	2.33	0.46
2:B:99[B]:VAL:HG13	2:B:101:TRP:HD1	1.79	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:601:GLJ:OE1	3:B:601:GLJ:CA	2.63	0.46
2:B:181:VAL:HG13	2:B:238[A]:ILE:HD13	1.99	0.45
2:B:191:HIS:HB2	2:B:238[A]:ILE:HD11	1.97	0.45
1:A:33:ARG:HD2	1:A:35:ARG:CZ	2.47	0.45
2:B:442[A]:SER:HB2	4:B:813:HOH:O	2.16	0.45
2:B:483:SER:C	2:B:484:ARG:HG3	2.42	0.45
2:B:109[A]:ARG:HD3	2:B:430:GLN:NE2	2.28	0.44
4:A:313:HOH:O	2:B:191:HIS:HD2	2.00	0.44
2:B:64[B]:VAL:HG12	2:B:184:ASP:HB3	2.01	0.43
1:A:99[B]:MET:HE1	1:A:159:ARG:HB3	2.01	0.43
2:B:79:TYR:HB2	2:B:138:THR:CG2	2.49	0.43
2:B:299:ARG:HG3	4:B:1154:HOH:O	2.19	0.42
2:B:325:ARG:CZ	4:B:978:HOH:O	2.68	0.42
1:A:94:VAL:HB	1:A:99[B]:MET:HG3	2.02	0.42
2:B:104:ASP:HB3	2:B:116:PHE:CZ	2.54	0.41
2:B:516:LEU:HD23	2:B:516:LEU:HA	1.87	0.41
1:A:175[B]:MET:HE3	1:A:175[B]:MET:HB3	1.96	0.41
2:B:424[B]:THR:HG22	4:B:1198:HOH:O	2.20	0.41
2:B:176[A]:ILE:CD1	4:B:1378:HOH:O	2.68	0.41
2:B:200:ARG:HB2	2:B:201:PRO:CD	2.51	0.41
1:A:20[C]:SER:OG	4:A:351:HOH:O	2.22	0.40
2:B:250:HIS:HD2	4:B:1113:HOH:O	2.04	0.40
2:B:91:ARG:HB2	2:B:96:PHE:CZ	2.56	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:578:HOH:O	4:B:835:HOH:O[3_656]	1.14	1.06

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	224/229 (98%)	221 (99%)	3 (1%)	0	100	100
2	B	550/543 (101%)	537 (98%)	13 (2%)	0	100	100
All	All	774/772 (100%)	758 (98%)	16 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	171/171 (100%)	169 (99%)	2 (1%)	63	40
2	B	430/421 (102%)	419 (97%)	11 (3%)	40	11
All	All	601/592 (102%)	588 (98%)	13 (2%)	50	15

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	108	GLU
1	A	226	LYS
2	B	176[A]	ILE
2	B	176[B]	ILE
2	B	191	HIS
2	B	200	ARG
2	B	204	ASN
2	B	238[A]	ILE
2	B	238[B]	ILE
2	B	310	SER
2	B	346	LEU
2	B	422	ARG
2	B	457	LEU

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

Mol	Chain	Res	Type
2	B	35	HIS
2	B	60	HIS
2	B	94	ASN
2	B	123	HIS
2	B	143	GLN
2	B	165	GLN
2	B	191	HIS
2	B	204	ASN
2	B	224	HIS
2	B	250	HIS
2	B	259	HIS
2	B	286	HIS
2	B	337	ASN
2	B	408	GLN
2	B	430	GLN
2	B	479	ASN
2	B	530	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](#)

1 ligand is 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
3	GLJ	B	601	-	8,9,9	1.91	2 (25%)	7,11,11	2.34	2 (28%)

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
3	GLJ	B	601	-	-	5/9/9/9	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	601	GLJ	CG-CB	-3.91	1.42	1.53
3	B	601	GLJ	CB-CA	-2.20	1.48	1.53

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	601	GLJ	OXT-C-O	4.33	133.90	124.08
3	B	601	GLJ	CB-CA-N	3.74	119.86	110.12

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	601	GLJ	N-CA-CB-CG
3	B	601	GLJ	C-CA-CB-CG
3	B	601	GLJ	CA-CB-CG-CD
3	B	601	GLJ	OXT-C-CA-N
3	B	601	GLJ	O-C-CA-N

There are no ring outliers.

1 monomer is involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	601	GLJ	10	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	216/229 (94%)	-0.62	1 (0%) 87 90	4, 9, 22, 42	9 (4%)
2	B	535/543 (98%)	-0.48	1 (0%) 91 92	3, 9, 22, 40	17 (3%)
All	All	751/772 (97%)	-0.52	2 (0%) 90 91	3, 9, 22, 42	26 (3%)

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	88	ARG	2.7
1	A	14	ALA	2.5

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	GLJ	B	601	10/10	0.96	0.09	6,14,29,35	0

6.5 Other polymers [i](#)

There are no such residues in this entry.