Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 06:04 am BST

PDB ID	:	4CE7
Title	:	Crystal structure of a novel unsaturated beta-glucuronyl hydrolase enzyme, belonging to family GH105, involved in ulvan degradation
Authors	:	Nyvall Collen, P.; Jeudy, A.; Groisillier, A.; Coutinho, P.M.; Helbert, W.; Czjzek, M.
Deposited on	:	2013-11-09
Resolution	:	1.90 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:

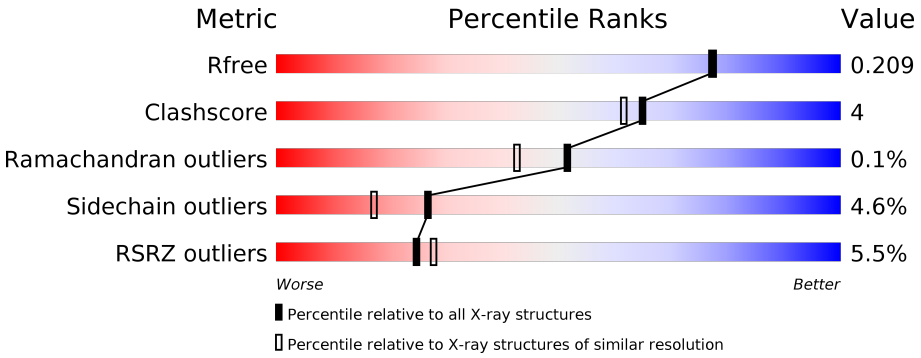
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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 <=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	370	7% 78% 12% • • 7%
1	B	370	4% 82% 11% • 6%
1	C	370	5% 78% 12% • • 8%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 9354 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 UNSATURATED 3S-RHAMNOGLYCURONYL HYDROLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	345	Total	C	N	O	S	0	3	0
			2868	1841	479	535	13			
1	B	348	Total	C	N	O	S	0	2	0
			2873	1843	480	537	13			
1	C	342	Total	C	N	O	S	0	5	0
			2855	1832	477	533	13			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	HIS	-	expression tag	UNP L7P9J4
A	-6	HIS	-	expression tag	UNP L7P9J4
A	-5	HIS	-	expression tag	UNP L7P9J4
A	-4	HIS	-	expression tag	UNP L7P9J4
A	-3	HIS	-	expression tag	UNP L7P9J4
A	-2	HIS	-	expression tag	UNP L7P9J4
A	-1	GLY	-	expression tag	UNP L7P9J4
A	0	SER	-	expression tag	UNP L7P9J4
B	-7	HIS	-	expression tag	UNP L7P9J4
B	-6	HIS	-	expression tag	UNP L7P9J4
B	-5	HIS	-	expression tag	UNP L7P9J4
B	-4	HIS	-	expression tag	UNP L7P9J4
B	-3	HIS	-	expression tag	UNP L7P9J4
B	-2	HIS	-	expression tag	UNP L7P9J4
B	-1	GLY	-	expression tag	UNP L7P9J4
B	0	SER	-	expression tag	UNP L7P9J4
C	-7	HIS	-	expression tag	UNP L7P9J4
C	-6	HIS	-	expression tag	UNP L7P9J4
C	-5	HIS	-	expression tag	UNP L7P9J4
C	-4	HIS	-	expression tag	UNP L7P9J4
C	-3	HIS	-	expression tag	UNP L7P9J4
C	-2	HIS	-	expression tag	UNP L7P9J4

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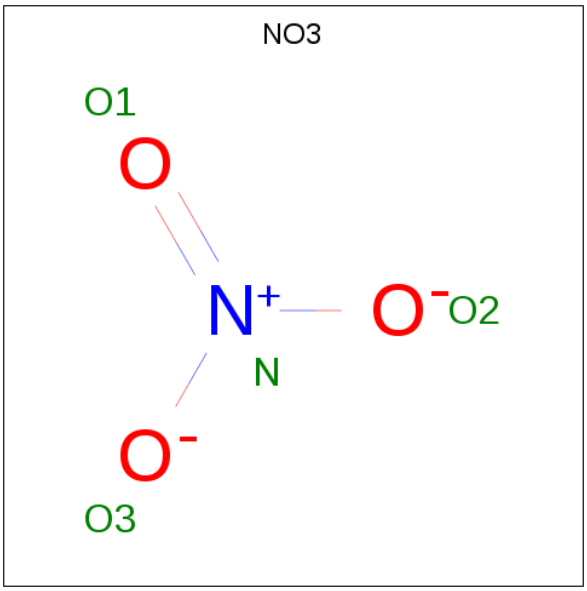
Chain	Residue	Modelled	Actual	Comment	Reference
C	-1	GLY	-	expression tag	UNP L7P9J4
C	0	SER	-	expression tag	UNP L7P9J4

- Molecule 2 is NITRATE ION (three-letter code: NO3) (formula: NO_3).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total N O 4 1 3	0	0
2	B	1	Total N O 4 1 3	0	0
2	C	1	Total N O 4 1 3	0	0
2	C	1	Total N O 4 1 3	0	0

- Molecule 3 is water.

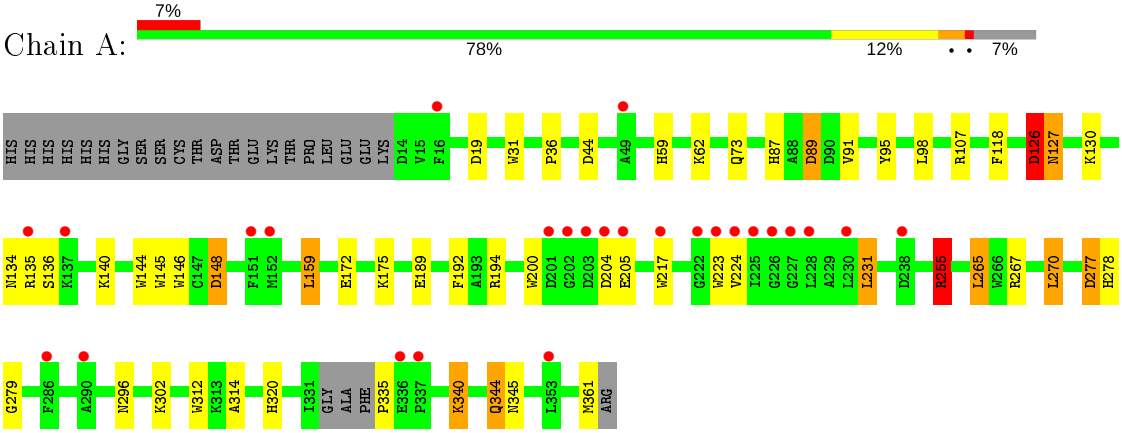
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	265	Total O 265 265	0	0
3	B	260	Total O 260 260	0	0
3	C	217	Total O 217 217	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA 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.

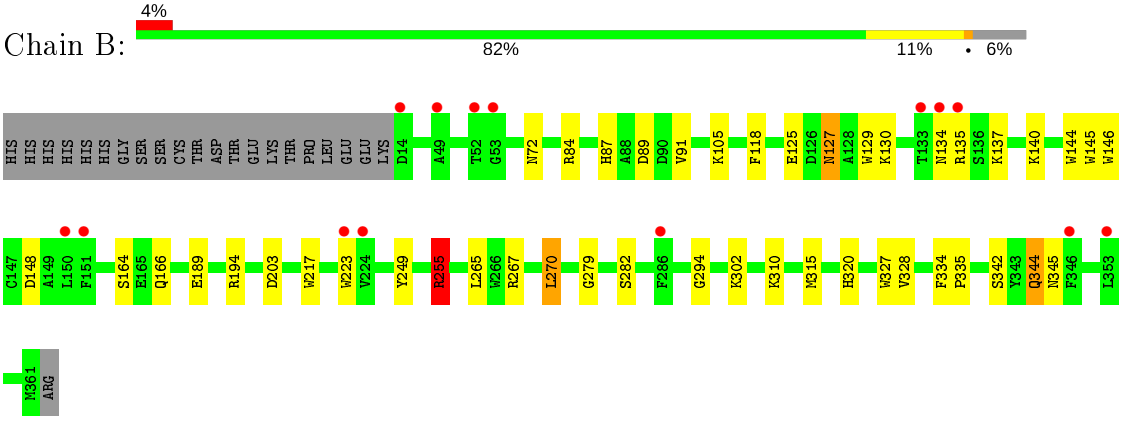
• Molecule 1: UNSATURATED 3S-RHAMNOGLYCURONYL HYDROLASE

Chain A:



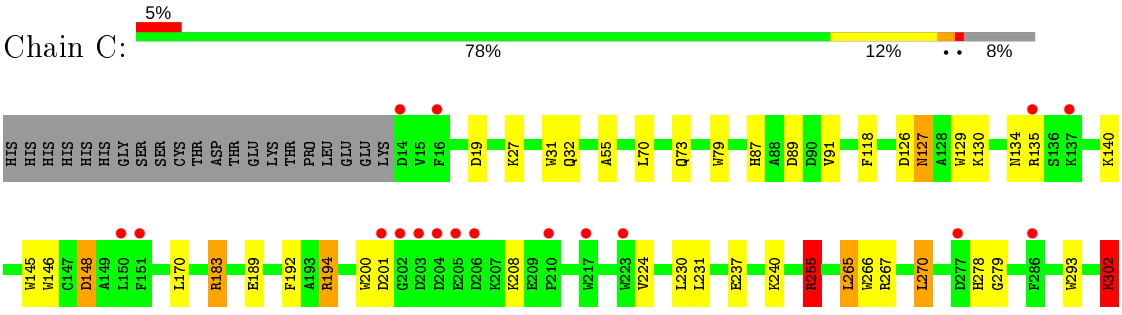
• Molecule 1: UNSATURATED 3S-RHAMNOGLYCURONYL HYDROLASE

Chain B:



• Molecule 1: UNSATURATED 3S-RHAMNOGLYCURONYL HYDROLASE

Chain C:




W312
K313
A314
H320
D321
D322
W327
I331
GLY
ALA
PHE
PRO
GLU
PRO
A338
S339
K340
D341
Q344
N345
T348
M361
ARG

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	92.98Å 93.33Å 156.76Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.19 – 1.90 34.81 – 1.90	Depositor EDS
% Data completeness (in resolution range)	98.9 (45.19-1.90) 98.9 (34.81-1.90)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.69 (at 1.89Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.162 , 0.207 0.164 , 0.209	Depositor DCC
R_{free} test set	5313 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å^2)	26.0	Xtriage
Anisotropy	0.040	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.38 , 45.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.023 for k,h,-l	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	9354	wwPDB-VP
Average B, all atoms (Å^2)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.31% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: NO3

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	1.09	7/2962 (0.2%)	1.03	14/4017 (0.3%)
1	B	1.07	6/2967 (0.2%)	0.95	9/4025 (0.2%)
1	C	1.08	7/2953 (0.2%)	1.02	16/4004 (0.4%)
All	All	1.08	20/8882 (0.2%)	1.00	39/12046 (0.3%)

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	145	TRP	CD2-CE2	7.56	1.50	1.41
1	B	145	TRP	CD2-CE2	7.33	1.50	1.41
1	A	223	TRP	CD2-CE2	7.13	1.50	1.41
1	C	79	TRP	CD2-CE2	7.02	1.49	1.41
1	C	266	TRP	CD2-CE2	6.75	1.49	1.41
1	A	200	TRP	CD2-CE2	6.53	1.49	1.41
1	B	146	TRP	CD2-CE2	6.14	1.48	1.41
1	C	312	TRP	CD2-CE2	6.04	1.48	1.41
1	A	144	TRP	CD2-CE2	5.92	1.48	1.41
1	B	327	TRP	CD2-CE2	5.68	1.48	1.41
1	B	144	TRP	CD2-CE2	5.53	1.48	1.41
1	C	31	TRP	CD2-CE2	5.44	1.47	1.41
1	A	146	TRP	CD2-CE2	5.43	1.47	1.41
1	B	223	TRP	CD2-CE2	5.42	1.47	1.41
1	C	327	TRP	CD2-CE2	5.34	1.47	1.41
1	A	31	TRP	CD2-CE2	5.34	1.47	1.41
1	B	217	TRP	CD2-CE2	5.18	1.47	1.41
1	C	145	TRP	CD2-CE2	5.10	1.47	1.41
1	A	312	TRP	CD2-CE2	5.06	1.47	1.41
1	C	293	TRP	CD2-CE2	5.01	1.47	1.41

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	255	ARG	NE-CZ-NH2	-17.46	111.57	120.30
1	C	183	ARG	NE-CZ-NH2	-15.09	112.76	120.30
1	A	255	ARG	NE-CZ-NH1	11.61	126.10	120.30
1	C	183	ARG	NE-CZ-NH1	10.38	125.49	120.30
1	A	255	ARG	CG-CD-NE	-8.84	93.24	111.80
1	B	255	ARG	NE-CZ-NH2	-8.63	115.99	120.30
1	A	107	ARG	NE-CZ-NH1	8.20	124.40	120.30
1	C	126[A]	ASP	CB-CG-OD1	8.06	125.56	118.30
1	C	126[B]	ASP	CB-CG-OD1	8.06	125.56	118.30
1	A	44	ASP	CB-CG-OD1	7.75	125.28	118.30
1	C	255	ARG	CG-CD-NE	-7.08	96.94	111.80
1	A	277	ASP	CB-CG-OD2	6.98	124.59	118.30
1	C	255	ARG	NE-CZ-NH2	-6.66	116.97	120.30
1	A	255	ARG	CD-NE-CZ	6.46	132.65	123.60
1	B	267	ARG	NE-CZ-NH2	-6.44	117.08	120.30
1	A	231	LEU	CA-CB-CG	6.32	129.84	115.30
1	A	148	ASP	CB-CG-OD1	6.27	123.94	118.30
1	C	148	ASP	CB-CG-OD1	6.25	123.93	118.30
1	C	194	ARG	NE-CZ-NH1	6.13	123.37	120.30
1	A	107	ARG	NE-CZ-NH2	-6.11	117.25	120.30
1	A	270	LEU	CB-CG-CD1	6.09	121.35	111.00
1	A	126	ASP	CB-CA-C	5.93	122.25	110.40
1	B	203	ASP	CB-CG-OD1	5.82	123.54	118.30
1	C	322	ASP	CB-CG-OD1	5.82	123.53	118.30
1	C	183	ARG	CB-CG-CD	-5.72	96.72	111.60
1	B	270	LEU	CB-CG-CD1	5.69	120.68	111.00
1	B	255	ARG	CG-CD-NE	-5.66	99.91	111.80
1	C	194	ARG	CA-CB-CG	5.57	125.66	113.40
1	C	270	LEU	CB-CG-CD1	5.53	120.41	111.00
1	C	341	ASP	CB-CG-OD1	5.53	123.28	118.30
1	B	84	ARG	NE-CZ-NH1	-5.38	117.61	120.30
1	B	255	ARG	NE-CZ-NH1	5.37	122.98	120.30
1	C	148	ASP	CB-CG-OD2	-5.36	113.47	118.30
1	A	267	ARG	NE-CZ-NH1	5.32	122.96	120.30
1	B	84	ARG	NE-CZ-NH2	5.30	122.95	120.30
1	C	170	LEU	CB-CG-CD1	-5.25	102.07	111.00
1	B	194	ARG	NE-CZ-NH2	-5.17	117.72	120.30
1	A	89	ASP	CB-CG-OD1	5.11	122.90	118.30
1	C	230	LEU	CB-CG-CD1	-5.02	102.47	111.00

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	2868	0	2713	28	0
1	B	2873	0	2719	20	0
1	C	2855	0	2702	22	0
2	A	4	0	0	0	0
2	B	4	0	0	0	0
2	C	8	0	0	0	0
3	A	265	0	0	6	0
3	B	260	0	0	2	1
3	C	217	0	0	3	1
All	All	9354	0	8134	69	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 4.

All (69) 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:62:LYS:HE2	3:A:2064:HOH:O	1.07	1.24
1:A:126:ASP:HB2	3:A:2139:HOH:O	1.36	1.23
1:C:27:LYS:HE3	3:C:2005:HOH:O	1.75	0.86
1:A:278:HIS:HD2	1:A:279:GLY:O	1.72	0.72
1:A:189:GLU:OE1	1:A:255:ARG:NH2	2.24	0.70
1:A:335:PRO:HD3	3:A:2263:HOH:O	1.92	0.69
1:B:189:GLU:OE1	1:B:255:ARG:NH2	2.25	0.69
1:C:87:HIS:HD2	1:C:89:ASP:H	1.45	0.63
1:C:265:LEU:HD13	1:C:314:ALA:HB3	1.82	0.61
1:A:127:ASN:C	1:A:127:ASN:HD22	2.04	0.59
1:A:172:GLU:OE2	1:A:175:LYS:HE2	2.03	0.58
1:A:126:ASP:CB	3:A:2139:HOH:O	2.14	0.58
1:B:310:LYS:NZ	3:B:2212:HOH:O	2.37	0.57
1:B:127:ASN:ND2	1:B:130:LYS:H	2.02	0.57
1:A:265:LEU:HB3	1:A:279:GLY:HA3	1.89	0.55
1:C:134:ASN:ND2	1:C:140:LYS:H	2.05	0.55
1:A:98:LEU:HD13	1:A:159:LEU:HB3	1.89	0.54
1:C:302:LYS:HA	1:C:302:LYS:HE3	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:265:LEU:HD13	1:A:314:ALA:HB3	1.88	0.54
1:C:55:ALA:HA	1:C:70:LEU:HD11	1.89	0.54
1:B:134:ASN:ND2	1:B:140:LYS:H	2.06	0.53
1:A:344:GLN:HE21	1:A:345:ASN:H	1.57	0.53
1:C:265:LEU:HD13	1:C:314:ALA:CB	2.39	0.53
1:C:265:LEU:HB3	1:C:279:GLY:HA3	1.90	0.52
1:A:87:HIS:HD2	1:A:89:ASP:H	1.58	0.51
1:C:344:GLN:HE21	1:C:345:ASN:H	1.59	0.51
1:C:278:HIS:HD2	1:C:279:GLY:O	1.94	0.51
1:C:189:GLU:OE1	1:C:255:ARG:NH2	2.44	0.50
1:B:164:SER:CB	1:B:166:GLN:HE21	2.25	0.50
1:C:140:LYS:NZ	3:C:2124:HOH:O	2.38	0.50
1:B:342:SER:HB2	3:B:2249:HOH:O	2.12	0.50
1:A:62:LYS:CE	3:A:2064:HOH:O	1.92	0.49
1:C:32:GLN:HB2	1:C:348:THR:HG21	1.93	0.49
1:A:127:ASN:ND2	1:A:130:LYS:H	2.11	0.48
1:B:334:PHE:HB2	1:B:335:PRO:HD2	1.96	0.48
1:C:320:HIS:HE1	1:C:340:LYS:O	1.97	0.48
1:B:334:PHE:HB2	1:B:335:PRO:CD	2.43	0.47
1:B:344:GLN:HE21	1:B:345:ASN:H	1.63	0.47
1:B:87:HIS:HD2	1:B:89:ASP:H	1.63	0.47
1:A:59:HIS:HD2	3:A:2060:HOH:O	1.97	0.47
1:B:265:LEU:HB3	1:B:279:GLY:HA3	1.98	0.46
1:B:282:SER:HB3	1:B:328:VAL:HG13	1.98	0.45
1:C:237:GLU:HB2	3:C:2178:HOH:O	2.16	0.45
1:C:200:TRP:CZ2	1:C:208:LYS:HG2	2.52	0.45
1:A:36:PRO:HG2	1:B:72[B]:ASN:ND2	2.32	0.45
1:A:91:VAL:HB	1:A:118:PHE:CE1	2.52	0.45
1:A:194:ARG:HB2	1:A:217[A]:TRP:HD1	1.82	0.44
1:B:127:ASN:HD22	1:B:129:TRP:H	1.65	0.44
1:A:344:GLN:HE21	1:A:345:ASN:N	2.16	0.44
1:B:249:TYR:OH	1:B:294:GLY:HA3	2.18	0.44
1:A:320:HIS:HE1	1:A:340:LYS:O	2.01	0.43
1:B:315:MET:HE2	1:B:315:MET:HB2	1.63	0.43
1:A:192:PHE:CE1	1:A:224:VAL:HG11	2.52	0.43
1:A:134:ASN:ND2	1:A:136:SER:H	2.17	0.42
1:A:296:ASN:OD1	1:A:361:MET:HA	2.19	0.42
1:B:91:VAL:HB	1:B:118:PHE:CE1	2.53	0.42
1:B:127:ASN:HD22	1:B:129:TRP:N	2.17	0.42
1:C:91:VAL:HB	1:C:118:PHE:CE1	2.55	0.42
1:A:127:ASN:C	1:A:127:ASN:ND2	2.71	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:105:LYS:HD2	1:B:105:LYS:HA	1.79	0.42
1:C:127:ASN:ND2	1:C:130:LYS:H	2.17	0.42
1:A:134:ASN:ND2	1:A:140:LYS:H	2.17	0.41
1:C:265:LEU:O	1:C:267:ARG:NH1	2.51	0.41
1:C:192:PHE:CE1	1:C:224:VAL:HG11	2.55	0.41
1:B:265:LEU:HD12	1:B:265:LEU:HA	1.94	0.41
1:C:146:TRP:HA	1:C:194:ARG:O	2.21	0.41
1:C:127:ASN:HD22	1:C:129:TRP:H	1.68	0.40
1:A:265:LEU:HD13	1:A:314:ALA:CB	2.52	0.40
1:A:95:TYR:CE2	1:A:159:LEU:HD22	2.57	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)
3:B:2103:HOH:O	3:C:2178:HOH:O[3_655]	2.15	0.05

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	344/370 (93%)	339 (98%)	5 (2%)	0	100	100
1	B	348/370 (94%)	338 (97%)	10 (3%)	0	100	100
1	C	343/370 (93%)	335 (98%)	7 (2%)	1 (0%)	41	31
All	All	1035/1110 (93%)	1012 (98%)	22 (2%)	1 (0%)	51	43

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	302	LYS

5.3.2 Protein sidechains (i)

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	298/317 (94%)	281 (94%)	17 (6%)	20	11
1	B	298/317 (94%)	288 (97%)	10 (3%)	37	28
1	C	297/317 (94%)	283 (95%)	14 (5%)	26	16
All	All	893/951 (94%)	852 (95%)	41 (5%)	27	17

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	ASP
1	A	73	GLN
1	A	126	ASP
1	A	127	ASN
1	A	135	ARG
1	A	148	ASP
1	A	159	LEU
1	A	204	ASP
1	A	205	GLU
1	A	231	LEU
1	A	255	ARG
1	A	265	LEU
1	A	270	LEU
1	A	277	ASP
1	A	302	LYS
1	A	340	LYS
1	A	344	GLN
1	B	125	GLU
1	B	127	ASN
1	B	135	ARG
1	B	137	LYS
1	B	148	ASP
1	B	255	ARG
1	B	270	LEU
1	B	302	LYS
1	B	320	HIS

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Mol	Chain	Res	Type
1	B	344	GLN
1	C	19	ASP
1	C	73	GLN
1	C	127	ASN
1	C	135	ARG
1	C	148	ASP
1	C	183	ARG
1	C	201	ASP
1	C	231	LEU
1	C	240	LYS
1	C	255	ARG
1	C	265	LEU
1	C	270	LEU
1	C	302	LYS
1	C	344	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	73	GLN
1	A	87	HIS
1	A	108	ASN
1	A	127	ASN
1	A	134	ASN
1	A	166	GLN
1	A	247	ASN
1	A	278	HIS
1	A	287	HIS
1	A	320	HIS
1	A	344	GLN
1	B	87	HIS
1	B	127	ASN
1	B	134	ASN
1	B	163	GLN
1	B	166	GLN
1	B	287	HIS
1	B	344	GLN
1	C	73	GLN
1	C	87	HIS
1	C	127	ASN
1	C	134	ASN
1	C	166	GLN

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Mol	Chain	Res	Type
1	C	278	HIS
1	C	287	HIS
1	C	320	HIS
1	C	344	GLN

5.3.3 RNA

There are no RNA molecules 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

4 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	NO3	A	1362	-	1,3,3	3.96	1 (100%)	0,3,3	0.00	-
2	NO3	B	1362	-	1,3,3	3.54	1 (100%)	0,3,3	0.00	-
2	NO3	C	1362	-	1,3,3	3.67	1 (100%)	0,3,3	0.00	-
2	NO3	C	1363	-	1,3,3	3.56	1 (100%)	0,3,3	0.00	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1362	NO3	O1-N	3.96	1.42	1.24
2	C	1362	NO3	O1-N	3.67	1.41	1.24
2	C	1363	NO3	O1-N	3.56	1.40	1.24
2	B	1362	NO3	O1-N	3.54	1.40	1.24

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, 95^{th} 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	345/370 (93%)	0.11	26 (7%)	14	15	16, 26, 45, 89	0
1	B	348/370 (94%)	0.04	14 (4%)	38	41	17, 25, 41, 57	0
1	C	342/370 (92%)	0.09	17 (4%)	28	32	15, 25, 48, 76	0
All	All	1035/1110 (93%)	0.08	57 (5%)	25	28	15, 25, 46, 89	0

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	202	GLY	5.7
1	C	204	ASP	4.7
1	A	204	ASP	4.3
1	A	336	GLU	4.3
1	A	217[A]	TRP	4.0
1	A	203	ASP	3.8
1	C	205	GLU	3.7
1	A	205	GLU	3.7
1	C	210	PRO	3.6
1	C	217[A]	TRP	3.6
1	A	337	PRO	3.5
1	A	223	TRP	3.3
1	A	151	PHE	3.2
1	A	202	GLY	3.1
1	C	201	ASP	3.1
1	B	223	TRP	3.1
1	B	151	PHE	3.1
1	C	203	ASP	3.0
1	A	224	VAL	3.0
1	A	201	ASP	2.9
1	B	224	VAL	2.9
1	C	14	ASP	2.8
1	A	230	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
1	C	151	PHE	2.8
1	B	353	LEU	2.8
1	C	277	ASP	2.8
1	A	137	LYS	2.8
1	B	134	ASN	2.8
1	C	16	PHE	2.6
1	B	135	ARG	2.6
1	B	49	ALA	2.6
1	C	137	LYS	2.6
1	A	286	PHE	2.5
1	B	346	PHE	2.5
1	A	225	ILE	2.5
1	B	133	THR	2.4
1	C	286	PHE	2.4
1	C	150	LEU	2.3
1	B	14	ASP	2.3
1	B	150	LEU	2.3
1	C	223	TRP	2.3
1	A	222	GLY	2.2
1	A	226	GLY	2.2
1	A	227	GLY	2.2
1	A	49	ALA	2.2
1	B	286	PHE	2.2
1	A	228	LEU	2.2
1	B	52	THR	2.1
1	A	238	ASP	2.1
1	C	135	ARG	2.1
1	A	353	LEU	2.1
1	B	53	GLY	2.1
1	A	16	PHE	2.1
1	A	135	ARG	2.0
1	A	152	MET	2.0
1	A	290	ALA	2.0
1	C	206	ASP	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 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. The B-factors column lists the minimum, median, 95^{th} 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
2	NO3	A	1362	4/4	0.84	0.16	58,62,67,70	0
2	NO3	C	1362	4/4	0.89	0.10	56,64,64,66	0
2	NO3	C	1363	4/4	0.89	0.12	61,61,65,70	0
2	NO3	B	1362	4/4	0.91	0.11	64,69,70,74	0

6.5 Other polymers (i)

There are no such residues in this entry.