



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 10:54 PM UTC

PDB ID : 2QKB / pdb_00002qkb
Title : Human RNase H catalytic domain mutant D210N in complex with 20-mer RNA/DNA hybrid
Authors : Nowotny, M.; Gaidamakov, S.A.; Ghirlando, R.; Cerritelli, S.M.; Crouch, R.J.; Yang, W.
Deposited on : 2007-07-10
Resolution : 2.40 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

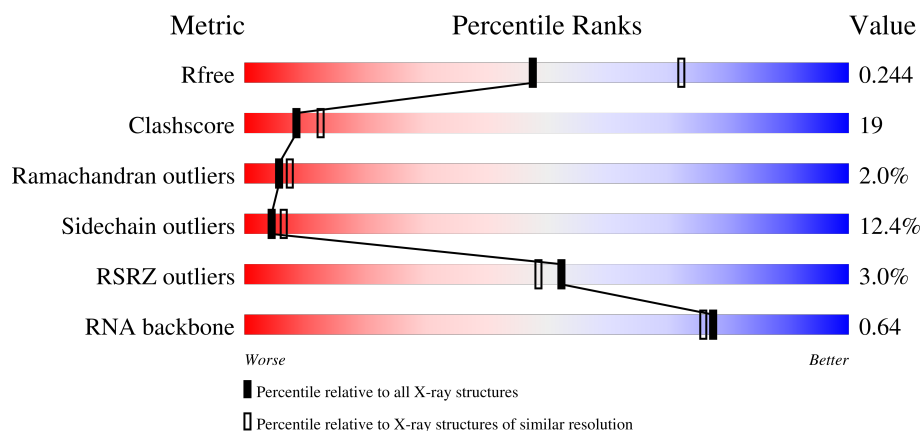
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)
RNA backbone	3983	1155 (2.70-2.10)

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	C	20	45% 45% 10%
2	D	20	35% 60% 5%
3	A	154	5% 60% 30% 6% ••
3	B	154	5% 58% 31% 7% ••

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 3244 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 RNA chain called 5'-R(*GP*GP*AP*GP*UP*GP*CP*GP*AP*CP*AP*C P*CP*UP*GP*AP*UP*UP*CP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	20	Total	C	N	O	P	0	0	1
			404	180	71	134	19			

- Molecule 2 is a DNA chain called 5'-D(*DGP*DGP*DAP*DAP*DTP*DCP*DAP*DGP*D GP*DTP*DGP*DTP*DCP*DGP*DCP*DAP*DCP*DTP*DCP*DT)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	20	Total	C	N	O	P	0	0	0
			408	195	75	119	19			

- Molecule 3 is a protein called Ribonuclease H1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	A	152	Total	C	N	O	S	Se	0	0	0
			1182	736	225	214	3	4			
3	B	150	Total	C	N	O	S	Se	0	0	0
			1174	732	223	212	3	4			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	133	GLY	-	expression tag	UNP O60930
A	134	SER	-	expression tag	UNP O60930
A	135	HIS	-	expression tag	UNP O60930
A	210	ASN	ASP	engineered mutation	UNP O60930
B	133	GLY	-	expression tag	UNP O60930
B	134	SER	-	expression tag	UNP O60930
B	135	HIS	-	expression tag	UNP O60930
B	210	ASN	ASP	engineered mutation	UNP O60930

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



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

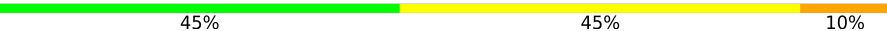
- Molecule 6 is water.

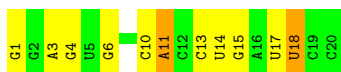
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	C	5	Total 5	O 5	0	0
6	D	8	Total 8	O 8	0	0
6	A	16	Total 16	O 16	0	0
6	B	29	Total 29	O 29	0	0

3 Residue-property plots [i](#)

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.

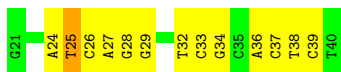
- Molecule 1: 5'-R(*GP*GP*AP*GP*UP*GP*CP*GP*AP*CP*AP*CP*CP*UP*GP*AP*UP*UP*CP*C)-3')

Chain C: 



- Molecule 2: 5'-D(*DGP*DGP*DAP*DAP*DTP*DCP*DAP*DGP*DGP*DTP*DGP*DTP*DCP*DGP*DCP*DAP*DCP*DTP*DCP*DT)-3'

Chain D: 



- Molecule 3: Ribonuclease H1

Chain A: 



- Molecule 3: Ribonuclease H1

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	109.99Å 109.99Å 67.25Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	40.00 – 2.40 40.00 – 2.40	Depositor EDS
% Data completeness (in resolution range)	90.8 (40.00-2.40) 90.8 (40.00-2.40)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.26 (at 2.29Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.208 , 0.255 0.195 , 0.244	Depositor DCC
R_{free} test set	1680 reflections (8.53%)	wwPDB-VP
Wilson B-factor (Å ²)	50.0	Xtriage
Anisotropy	0.555	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 34.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.038 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3244	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.72% 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: EDO, SO4

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	C	0.63	1/450 (0.2%)	0.88	0/700
2	D	0.46	0/457	0.89	0/704
3	A	0.82	1/1204 (0.1%)	1.19	7/1618 (0.4%)
3	B	0.85	1/1195 (0.1%)	1.27	14/1605 (0.9%)
All	All	0.77	3/3306 (0.1%)	1.14	21/4627 (0.5%)

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	C	0	3
2	D	0	1
All	All	0	4

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	162	VAL	CA-CB	6.44	1.63	1.54
3	A	162	VAL	CA-CB	5.60	1.61	1.54
1	C	1	DG	O3'-P	-5.58	1.52	1.61

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	261	VAL	CA-C-N	8.48	129.61	120.11
3	B	261	VAL	C-N-CA	8.48	129.61	120.11
3	B	189	ALA	N-CA-C	-6.69	104.07	111.36
3	B	168	HIS	CA-C-N	6.38	126.06	119.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	168	HIS	C-N-CA	6.38	126.06	119.56
3	B	256	ILE	N-CA-C	6.20	116.79	108.11
3	B	141	VAL	N-CA-C	6.09	118.12	108.87
3	A	283	GLN	N-CA-C	6.06	119.11	109.65
3	A	237	GLU	N-CA-C	5.98	119.38	109.46
3	B	220	ASN	N-CA-C	5.89	123.35	110.80
3	A	221	TRP	N-CA-C	5.86	123.27	110.80
3	B	165	GLY	CA-C-N	5.80	127.09	119.84
3	B	165	GLY	C-N-CA	5.80	127.09	119.84
3	A	189	ALA	N-CA-C	-5.70	105.20	111.82
3	A	170	LEU	N-CA-C	5.43	119.77	113.20
3	A	134	SER	N-CA-C	5.42	118.38	109.72
3	A	239	ILE	N-CA-C	5.39	117.79	111.05
3	B	173	GLY	N-CA-C	-5.30	100.61	113.18
3	B	138	ASP	N-CA-C	-5.19	106.80	113.23
3	B	262	PRO	N-CA-C	5.07	118.58	111.33
3	B	180	GLN	N-CA-C	5.01	116.47	108.41

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	11	A	Sidechain
1	C	13	C	Sidechain
1	C	18	U	Sidechain
2	D	25	DT	Sidechain

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	C	404	0	206	8	0
2	D	408	0	227	12	0
3	A	1182	0	1155	53	0
3	B	1174	0	1148	44	0
4	A	5	0	0	0	0
4	B	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	4	0	6	1	0
5	B	4	0	6	0	0
6	A	16	0	0	0	0
6	B	29	0	0	0	0
6	C	5	0	0	0	0
6	D	8	0	0	0	0
All	All	3244	0	2748	112	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 19.

All (112) 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)
3:A:278:ARG:HH21	3:A:278:ARG:HB2	1.08	1.17
3:B:154:ARG:HH11	3:B:154:ARG:HB2	1.14	1.09
3:B:154:ARG:HG3	3:B:155:ARG:H	1.28	0.98
3:A:278:ARG:HH21	3:A:278:ARG:CB	1.81	0.93
3:A:208:TYR:HB3	3:A:261:VAL:HG21	1.57	0.87
3:A:232:THR:HG23	3:A:234:ALA:H	1.40	0.86
3:B:154:ARG:HB2	3:B:154:ARG:NH1	1.89	0.86
3:B:154:ARG:HG3	3:B:155:ARG:N	1.91	0.84
3:A:208:TYR:HB3	3:A:261:VAL:CG2	2.06	0.84
3:A:278:ARG:HB2	3:A:278:ARG:NH2	1.91	0.84
3:B:278:ARG:HG2	3:B:278:ARG:HH11	1.43	0.83
3:A:236:LYS:NZ	3:A:236:LYS:HB2	1.92	0.82
3:B:232:THR:HG23	3:B:234:ALA:H	1.43	0.82
3:B:155:ARG:O	3:B:156:PRO:C	2.29	0.74
3:A:230:TRP:CD2	3:A:241:LYS:HG3	2.23	0.73
3:A:196:GLN:O	3:A:199:THR:HB	1.88	0.73
3:B:140:VAL:HG11	3:B:202:ILE:HD13	1.71	0.71
3:A:162:VAL:HG13	3:A:172:VAL:HG22	1.72	0.71
3:A:230:TRP:CE3	3:A:241:LYS:HG3	2.26	0.70
3:B:232:THR:HG22	3:B:236:LYS:H	1.54	0.70
3:A:272:GLU:OE2	3:A:275:ARG:HD3	1.91	0.70
2:D:25:DT:OP2	3:B:179:ARG:NH1	2.25	0.68
3:A:232:THR:HG23	3:A:234:ALA:N	2.09	0.68
3:B:175:ARG:O	3:B:177:PRO:HD3	1.95	0.67
3:A:236:LYS:HG2	3:A:237:GLU:H	1.60	0.66
3:A:203:ASN:HD21	3:A:204:LYS:HE3	1.60	0.66
3:B:178:GLY:O	3:B:180:GLN:NE2	2.30	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:212:MSE:HG2	3:B:260:HIS:CD2	2.32	0.65
3:B:149:SER:O	3:B:150:SER:HB2	1.97	0.64
3:B:282:LYS:C	3:B:283:GLN:HG2	2.22	0.63
3:B:140:VAL:CG1	3:B:202:ILE:HD13	2.27	0.63
3:B:162:VAL:HG21	3:B:196:GLN:OE1	1.99	0.63
3:A:228:ASN:OD1	3:A:228:ASN:C	2.45	0.60
3:A:232:THR:HG22	3:A:236:LYS:N	2.16	0.60
2:D:37:DC:H2''	2:D:38:DT:H5'	1.83	0.60
2:D:38:DT:H2''	2:D:39:DC:O5'	2.01	0.60
3:A:236:LYS:HB2	3:A:236:LYS:HZ3	1.67	0.59
3:A:164:TRP:CD1	3:A:170:LEU:HB3	2.38	0.59
3:A:236:LYS:HB2	3:A:236:LYS:HZ2	1.67	0.59
3:A:230:TRP:CD1	3:A:241:LYS:HE3	2.38	0.59
3:B:230:TRP:CE3	3:B:241:LYS:HG3	2.38	0.59
3:A:140:VAL:HG11	3:A:202:ILE:HD13	1.84	0.58
3:A:278:ARG:HH21	3:A:278:ARG:CG	2.16	0.58
3:A:236:LYS:NZ	3:A:236:LYS:CB	2.64	0.58
3:A:264:HIS:N	3:A:264:HIS:CD2	2.69	0.57
3:B:247:LEU:O	3:B:251:THR:HG23	2.05	0.57
3:A:232:THR:HG22	3:A:236:LYS:H	1.70	0.56
3:B:154:ARG:NE	3:B:155:ARG:HG2	2.20	0.56
3:B:262:PRO:O	3:B:263:GLY:C	2.49	0.55
3:A:236:LYS:HZ3	3:A:236:LYS:CB	2.19	0.55
1:C:3:A:H2'	1:C:4:G:O4'	2.07	0.54
3:A:249:ARG:O	3:A:252:GLN:HG2	2.08	0.54
3:B:148:CYS:HA	3:B:157:ARG:O	2.08	0.54
3:B:272:GLU:O	3:B:276:LEU:HD23	2.08	0.54
1:C:17:U:H2'	1:C:18:U:O4'	2.08	0.53
3:B:252:GLN:HG3	3:B:252:GLN:O	2.07	0.53
3:A:179:ARG:HE	5:A:9003:EDO:H21	1.73	0.53
3:A:148:CYS:HA	3:A:157:ARG:O	2.09	0.53
3:B:278:ARG:HG2	3:B:278:ARG:NH1	2.18	0.53
3:A:246:ALA:O	3:A:249:ARG:HG2	2.10	0.52
3:A:229:GLY:O	3:A:231:LYS:HD3	2.11	0.50
3:B:157:ARG:HG2	3:B:157:ARG:HH11	1.76	0.50
3:B:278:ARG:HH11	3:B:278:ARG:CG	2.17	0.49
2:D:32:DT:H2'	2:D:33:DC:C6	2.48	0.49
3:A:267:PHE:O	3:A:271:GLU:HG3	2.13	0.48
3:B:179:ARG:NH1	3:B:181:THR:HG21	2.29	0.48
3:A:232:THR:CG2	3:A:236:LYS:H	2.27	0.48
1:C:18:U:H5'	3:B:210:ASN:O	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:228:ASN:OD1	3:A:228:ASN:O	2.33	0.47
2:D:28:DG:H2'	2:D:29:DG:H5''	1.96	0.47
3:B:149:SER:O	3:B:150:SER:CB	2.63	0.47
3:B:207:LEU:C	3:B:207:LEU:HD23	2.40	0.47
3:B:157:ARG:HG2	3:B:157:ARG:NH1	2.31	0.46
3:A:168:HIS:CE1	3:A:170:LEU:HB2	2.50	0.46
3:A:187:ILE:HD11	3:A:213:PHE:HE2	1.81	0.46
3:A:241:LYS:O	3:A:245:VAL:HG23	2.17	0.45
2:D:39:DC:OP1	2:D:39:DC:H4'	2.17	0.45
3:B:217:GLY:HA2	3:B:221:TRP:HB2	1.99	0.44
3:B:139:PHE:HB3	3:B:206:VAL:HG23	1.99	0.44
3:B:158:ALA:O	3:B:175:ARG:HD2	2.18	0.44
1:C:6:G:C2	2:D:36:DA:C2	3.05	0.44
3:A:187:ILE:HD11	3:A:213:PHE:CE2	2.52	0.44
3:A:278:ARG:HA	3:A:278:ARG:HD3	1.78	0.44
3:A:156:PRO:HB2	3:A:180:GLN:HB2	2.00	0.43
1:C:10:C:H2'	1:C:11:A:O4'	2.18	0.43
3:B:150:SER:O	3:B:151:ASN:C	2.61	0.43
3:A:208:TYR:CD2	3:A:261:VAL:HG22	2.54	0.43
2:D:24:DA:H2''	2:D:25:DT:O5'	2.18	0.43
3:B:283:GLN:O	3:B:284:SER:O	2.37	0.42
3:A:187:ILE:HG21	3:A:243:ASP:HB3	2.00	0.42
3:A:160:ILE:HD11	3:A:174:ILE:HD11	2.01	0.42
3:A:230:TRP:O	3:A:231:LYS:HD2	2.19	0.42
3:B:154:ARG:CG	3:B:155:ARG:H	2.16	0.42
3:B:175:ARG:NH1	3:B:284:SER:HA	2.35	0.42
2:D:33:DC:H2'	2:D:34:DG:C8	2.55	0.42
1:C:4:G:N2	2:D:38:DT:C2	2.88	0.42
3:A:236:LYS:HG2	3:A:237:GLU:N	2.30	0.42
3:A:264:HIS:N	3:A:264:HIS:HD2	2.16	0.42
3:B:207:LEU:HD13	3:B:258:TRP:CZ3	2.55	0.42
3:B:232:THR:CG2	3:B:234:ALA:H	2.22	0.41
3:A:194:ILE:HD12	3:A:256:ILE:HD11	2.03	0.41
3:B:158:ALA:HB3	3:B:176:LEU:HB3	2.02	0.41
1:C:14:U:H2'	1:C:15:G:O4'	2.21	0.41
2:D:27:DA:H2'	2:D:28:DG:C8	2.54	0.41
2:D:25:DT:H2'	2:D:26:DC:C6	2.56	0.41
3:B:230:TRP:CD2	3:B:241:LYS:HG3	2.55	0.41
1:C:10:C:H4'	3:A:212:MSE:HG2	2.03	0.41
3:A:278:ARG:NH2	3:A:278:ARG:CG	2.77	0.41
3:B:188:HIS:HE1	3:B:243:ASP:OD1	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:210:ASN:HA	3:A:261:VAL:O	2.20	0.40
3:A:139:PHE:CD1	3:A:139:PHE:N	2.89	0.40
3:A:234:ALA:O	3:A:236:LYS:N	2.54	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	
3	A	150/154 (97%)	135 (90%)	13 (9%)	2 (1%)	9	14
3	B	146/154 (95%)	132 (90%)	10 (7%)	4 (3%)	4	4
All	All	296/308 (96%)	267 (90%)	23 (8%)	6 (2%)	6	7

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	228	ASN
3	B	283	GLN
3	A	235	GLY
3	B	263	GLY
3	B	156	PRO
3	B	266	GLY

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	
3	A	121/120 (101%)	103 (85%)	18 (15%)	3	3
3	B	121/120 (101%)	109 (90%)	12 (10%)	7	11
All	All	242/240 (101%)	212 (88%)	30 (12%)	4	6

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	134	SER
3	A	138	ASP
3	A	140	VAL
3	A	154	ARG
3	A	162	VAL
3	A	170	LEU
3	A	179	ARG
3	A	187	ILE
3	A	199	THR
3	A	201	ASN
3	A	228	ASN
3	A	236	LYS
3	A	257	GLN
3	A	264	HIS
3	A	274	ASP
3	A	275	ARG
3	A	276	LEU
3	A	278	ARG
3	B	140	VAL
3	B	154	ARG
3	B	162	VAL
3	B	170	LEU
3	B	199	THR
3	B	220	ASN
3	B	232	THR
3	B	250	LEU
3	B	252	GLN
3	B	261	VAL
3	B	282	LYS
3	B	283	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	203	ASN
3	A	216	ASN
3	A	264	HIS
3	B	188	HIS
3	B	203	ASN
3	B	216	ASN
3	B	220	ASN
3	B	240	ASN
3	B	252	GLN
3	B	257	GLN
3	B	283	GLN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	C	18/20 (90%)	0	0

There are no RNA backbone outliers to report.

There are no RNA pucker outliers to report.

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](#)

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
5	EDO	B	9004	-	3,3,3	0.40	0	2,2,2	0.35	0
4	SO4	A	9001	-	4,4,4	0.31	0	6,6,6	0.29	0
5	EDO	A	9003	-	3,3,3	0.39	0	2,2,2	0.38	0
4	SO4	B	9002	-	4,4,4	0.43	0	6,6,6	0.11	0

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
5	EDO	B	9004	-	-	1/1/1/1	-
5	EDO	A	9003	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	9003	EDO	O1-C1-C2-O2
5	B	9004	EDO	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	9003	EDO	1	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	C	20/20 (100%)	-0.44	0	100	100	41, 55, 90, 102	0
2	D	20/20 (100%)	-0.04	0	100	100	41, 53, 124, 145	0
3	A	148/154 (96%)	0.04	2 (1%)	73	69	34, 53, 82, 104	1 (0%)
3	B	146/154 (94%)	0.01	8 (5%)	30	27	33, 51, 98, 143	2 (1%)
All	All	334/348 (95%)	-0.01	10 (2%)	52	48	33, 53, 93, 145	3 (0%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	A	284	SER	5.4
3	B	135	HIS	5.0
3	B	151	ASN	3.4
3	B	263	GLY	3.1
3	B	134	SER	2.9
3	B	149	SER	2.8
3	B	137	GLY	2.3
3	B	284	SER	2.2
3	A	234	ALA	2.0
3	B	264	HIS	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 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($\text{\AA}^2$)	Q<0.9
4	SO4	B	9002	5/5	0.61	0.11	142,142,142,142	0
4	SO4	A	9001	5/5	0.88	0.08	99,100,101,101	0
5	EDO	B	9004	4/4	0.91	0.12	68,70,70,70	0
5	EDO	A	9003	4/4	0.95	0.08	63,63,63,63	0

6.5 Other polymers [i](#)

There are no such residues in this entry.