



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 17, 2017 – 02:00 AM EDT

PDB ID : 5HHL
Title : Reverse transcriptase domain of group II intron maturase from Eubacterium rectale in P21 space group
Authors : Zhao, C.; Pyle, A.M.
Deposited on : unknown
Resolution : 2.10 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
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20029824
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	rb-20029824

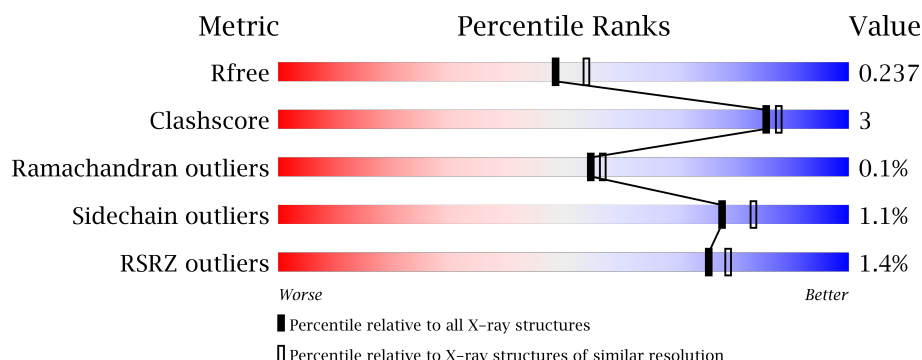
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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}	100719	4243 (2.10-2.10)
Clashscore	112137	4788 (2.10-2.10)
Ramachandran outliers	110173	4740 (2.10-2.10)
Sidechain outliers	110143	4741 (2.10-2.10)
RSRZ outliers	101464	4275 (2.10-2.10)

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 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	292	<div> <div>2%</div> <div> <div></div> <div>92%</div> <div>8%</div> <div>.</div> </div> </div>
1	B	292	<div> <div>%</div> <div> <div></div> <div>92%</div> <div>8%</div> <div>.</div> </div> </div>
1	C	292	<div> <div>2%</div> <div> <div></div> <div>90%</div> <div>8%</div> <div>..</div> </div> </div>
1	D	292	<div> <div>%</div> <div> <div></div> <div>92%</div> <div>7%</div> <div>.</div> </div> </div>
1	E	292	<div> <div>%</div> <div> <div></div> <div>94%</div> <div>5%</div> <div>.</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	F	292	<div><div></div><div>89%</div><div>8%</div><div>.</div></div>
1	G	292	<div><div>%</div><div></div><div>91%</div><div>7%</div><div>..</div></div>
1	H	292	<div><div>3%</div><div></div><div>83%</div><div>15%</div><div>.</div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 19755 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 Retron-type reverse transcriptase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	292	Total	C	N	O	S	0	3	0
			2343	1474	416	439	14			
1	B	290	Total	C	N	O	S	0	4	0
			2326	1467	414	432	13			
1	C	289	Total	C	N	O	S	0	1	0
			2314	1457	413	431	13			
1	D	289	Total	C	N	O	S	0	2	0
			2317	1460	410	433	14			
1	E	289	Total	C	N	O	S	0	1	0
			2308	1454	410	431	13			
1	F	283	Total	C	N	O	S	0	2	0
			2276	1436	404	423	13			
1	G	289	Total	C	N	O	S	0	1	0
			2307	1453	410	431	13			
1	H	287	Total	C	N	O	S	0	1	0
			2296	1445	408	430	13			

- Molecule 2 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	1	Total	K	0	0
			1	1		
2	D	1	Total	K	0	0
			1	1		
2	E	1	Total	K	0	0
			1	1		
2	H	1	Total	K	0	0
			1	1		
2	B	1	Total	K	0	0
			1	1		
2	C	1	Total	K	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total 1	K 1	0	0
2	F	1	Total 1	K 1	0	0

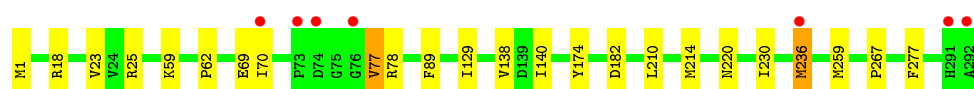
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	216	Total 218	O 218	0	2
3	B	197	Total 198	O 198	0	1
3	C	163	Total 163	O 163	0	0
3	D	169	Total 171	O 171	0	2
3	E	148	Total 148	O 148	0	0
3	F	155	Total 159	O 159	0	4
3	G	133	Total 133	O 133	0	0
3	H	69	Total 70	O 70	0	1

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.

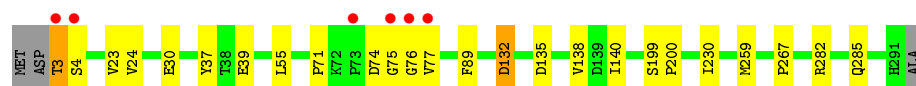
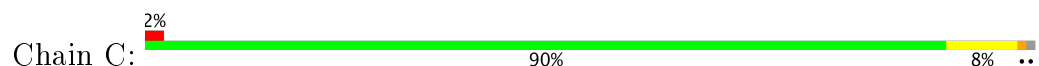
- Molecule 1: Retron-type reverse transcriptase



- Molecule 1: Retron-type reverse transcriptase



- Molecule 1: Retron-type reverse transcriptase



- Molecule 1: Retron-type reverse transcriptase



- Molecule 1: Retron-type reverse transcriptase

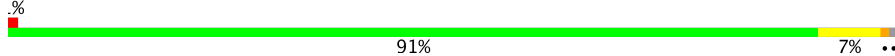


- Molecule 1: Retron-type reverse transcriptase

Chain F:  89% 8% .




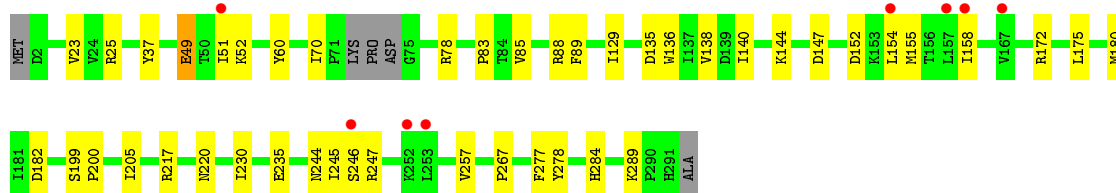
- Molecule 1: Retron-type reverse transcriptase

Chain G:  91% 7% ..



- Molecule 1: Retron-type reverse transcriptase

Chain H:  3% 83% 15% .



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	74.83Å 119.86Å 161.47Å 90.00° 92.08° 90.00°	Depositor
Resolution (Å)	49.14 – 2.10 49.14 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.6 (49.14-2.10) 99.2 (49.14-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.32 (at 2.10Å)	Xtriage
Refinement program	PHENIX (dev_2254: ???)	Depositor
R, R_{free}	0.201 , 0.238 0.200 , 0.237	Depositor DCC
R_{free} test set	8257 reflections (5.00%)	DCC
Wilson B-factor (Å ²)	36.4	Xtriage
Anisotropy	0.278	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 48.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.023 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	19755	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.06% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section:
K

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	0.24	0/2393	0.41	0/3221
1	B	0.24	0/2378	0.41	0/3202
1	C	0.24	0/2357	0.41	0/3173
1	D	0.24	0/2362	0.41	0/3179
1	E	0.24	0/2352	0.41	0/3167
1	F	0.24	0/2320	0.41	0/3121
1	G	0.24	0/2350	0.41	0/3164
1	H	0.23	0/2337	0.40	0/3145
All	All	0.24	0/18849	0.41	0/25372

There are no bond length outliers.

There are no bond angle outliers.

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	2343	0	2364	13	0
1	B	2326	0	2358	12	0
1	C	2314	0	2336	15	0
1	D	2317	0	2343	11	0
1	E	2308	0	2323	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	2276	0	2303	12	0
1	G	2307	0	2326	10	0
1	H	2296	0	2310	28	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
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
3	A	218	0	0	1	0
3	B	198	0	0	0	0
3	C	163	0	0	0	0
3	D	171	0	0	2	0
3	E	148	0	0	0	0
3	F	159	0	0	1	0
3	G	133	0	0	0	0
3	H	70	0	0	0	0
All	All	19755	0	18663	98	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 3.

All (98) 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:F:235:GLU:OE1	1:H:37:TYR:OH	2.11	0.68
1:C:3:THR:OG1	1:C:4:SER:N	2.30	0.64
1:F:23:VAL:HG21	1:F:89:PHE:HA	1.79	0.64
1:H:23:VAL:HG21	1:H:89:PHE:HA	1.78	0.64
1:D:23:VAL:HG21	1:D:89:PHE:HA	1.81	0.63
1:C:138:VAL:HB	1:C:230:ILE:HB	1.81	0.63
1:D:131:ASN:O	1:E:27:LYS:NZ	2.32	0.61
1:G:138:VAL:HB	1:G:230:ILE:HB	1.83	0.61
1:E:23:VAL:HG21	1:E:89:PHE:HA	1.81	0.61
1:A:18:ARG:NH1	3:A:403:HOH:O	2.34	0.60
1:C:23:VAL:HG21	1:C:89:PHE:HA	1.82	0.60
1:H:152:ASP:OD1	1:H:172:ARG:NH2	2.35	0.60
1:D:25:ARG:NH1	1:E:132:ASP:OD2	2.35	0.60
1:F:138:VAL:HB	1:F:230:ILE:HB	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:155:MET:HG3	1:H:172:ARG:HD3	1.85	0.57
1:B:23:VAL:HG21	1:B:89:PHE:HA	1.86	0.56
1:A:78:ARG:NH1	1:A:182:ASP:OD1	2.39	0.56
1:A:23:VAL:HG21	1:A:89:PHE:HA	1.86	0.56
1:G:23:VAL:HG21	1:G:89:PHE:HA	1.87	0.56
1:H:158:ILE:HG22	1:H:205:ILE:HG21	1.89	0.55
1:H:217:ARG:HH12	1:H:244:ASN:HB3	1.71	0.55
1:C:285:GLN:NE2	1:H:25:ARG:O	2.40	0.55
1:H:154:LEU:HD23	1:H:175:LEU:HD11	1.90	0.54
1:F:153:LYS:NZ	3:F:412:HOH:O	2.42	0.52
1:B:266:ARG:NH1	1:E:135:ASP:OD2	2.43	0.52
1:E:138:VAL:HB	1:E:230:ILE:HB	1.92	0.52
1:C:39:GLU:OE2	1:D:243:ARG:NH2	2.43	0.51
1:H:129:ILE:HD13	1:H:220:ASN:HB3	1.93	0.50
1:H:70:ILE:O	1:H:78:ARG:N	2.43	0.50
1:A:138:VAL:HB	1:A:230:ILE:HB	1.93	0.50
1:D:138:VAL:HB	1:D:230:ILE:HB	1.93	0.50
1:G:72:LYS:HD3	1:G:78:ARG:HB3	1.94	0.49
1:H:138:VAL:HB	1:H:230:ILE:HB	1.94	0.49
1:H:49:GLU:HA	1:H:52:LYS:HB2	1.95	0.49
1:D:187:ASP:OD2	3:D:401:HOH:O	2.20	0.48
1:A:267:PRO:O	1:A:277:PHE:HB2	2.13	0.48
1:F:160:ARG:O	1:F:163[A]:LYS:NZ	2.46	0.48
1:H:246:SER:HB2	1:H:257:VAL:HG21	1.95	0.47
1:H:267:PRO:O	1:H:277:PHE:HB2	2.15	0.47
1:C:71:PRO:HA	1:C:77:VAL:HG12	1.97	0.47
1:G:267:PRO:O	1:G:277:PHE:HB2	2.15	0.47
1:A:129:ILE:HD13	1:A:220:ASN:HB3	1.97	0.46
1:D:267:PRO:O	1:D:277:PHE:HB2	2.14	0.46
1:B:271:LYS:HD2	1:B:291:HIS:CD2	2.51	0.46
1:B:138:VAL:HB	1:B:230:ILE:HB	1.96	0.46
1:D:184:GLU:HG2	1:D:189:ILE:HB	1.98	0.46
1:H:78:ARG:NH1	1:H:182:ASP:OD1	2.46	0.46
1:C:132:ASP:OD1	1:H:25:ARG:NE	2.47	0.45
1:C:37:TYR:OH	1:D:235:GLU:OE1	2.28	0.45
1:G:8:GLU:OE2	1:G:56:ARG:NH1	2.48	0.45
1:F:14:ASP:OD1	1:F:18:ARG:NH1	2.48	0.45
1:G:74:ASP:OD1	1:G:75:GLY:N	2.49	0.45
1:E:267:PRO:O	1:E:277:PHE:HB2	2.17	0.45
1:H:136:TRP:CZ3	1:H:235:GLU:HB2	2.52	0.45
1:G:6:LEU:HD12	1:G:162:ILE:HG23	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:71:PRO:HA	1:F:77:VAL:HA	1.99	0.44
1:G:69:GLU:HB3	1:G:77:VAL:HG21	1.98	0.44
1:C:282:ARG:HH21	1:F:263:LYS:HB3	1.83	0.44
1:C:30:GLU:OE2	1:H:284:HIS:NE2	2.51	0.44
1:H:83:PRO:HG2	1:H:88:ARG:HG2	1.99	0.44
1:B:144:LYS:NZ	1:B:183:ASP:OD2	2.51	0.44
1:H:51:ILE:HD12	1:H:60:TYR:HE1	1.83	0.43
1:B:20:TYR:OH	1:B:41:LYS:HB2	2.18	0.43
1:D:271:LYS:HD2	1:D:291:HIS:CD2	2.54	0.43
1:H:37:TYR:HA	1:H:85:VAL:HG21	2.00	0.43
1:C:74:ASP:O	1:C:76:GLY:N	2.49	0.43
1:B:135:ASP:O	1:B:267:PRO:HD3	2.19	0.43
1:C:24:VAL:HG13	1:C:37:TYR:HD2	1.84	0.43
1:B:129:ILE:HD13	1:B:220:ASN:HB3	2.01	0.42
1:A:69:GLU:HB3	1:A:77:VAL:HG11	2.00	0.42
1:C:135:ASP:O	1:C:267:PRO:HD3	2.20	0.42
1:H:278:TYR:CE1	1:H:289:LYS:HD2	2.54	0.42
1:A:25[A]:ARG:HH21	1:B:128:ASN:HB3	1.84	0.42
1:E:70:ILE:HG21	1:E:78:ARG:HH21	1.83	0.42
1:F:144:LYS:NZ	1:F:183:ASP:OD2	2.51	0.42
1:F:199:SER:OG	1:F:200:PRO:HD3	2.20	0.42
1:F:267:PRO:O	1:F:277:PHE:HB2	2.20	0.42
1:B:3:THR:HB	1:B:6:LEU:HD11	2.01	0.42
1:A:210:LEU:HG	1:A:214:MET:HE2	2.02	0.41
1:G:57:THR:OG1	1:G:59:LYS:HD3	2.20	0.41
1:H:245:ILE:HG22	1:H:246:SER:H	1.85	0.41
1:E:183:ASP:HA	1:E:185:TYR:CE2	2.54	0.41
1:H:217:ARG:NH1	1:H:244:ASN:HB3	2.33	0.41
1:B:80:LEU:HD21	1:B:194:GLN:HB2	2.03	0.41
1:H:138:VAL:HG12	1:H:140:ILE:HG13	2.01	0.41
1:A:138:VAL:HG12	1:A:140:ILE:HG13	2.02	0.41
1:C:138:VAL:HG12	1:C:140:ILE:HG13	2.03	0.41
1:A:70:ILE:O	1:A:78:ARG:N	2.53	0.41
1:A:62:PRO:HB3	1:A:174:TYR:CE1	2.57	0.40
1:G:21:LEU:HA	1:G:21:LEU:HD12	1.91	0.40
1:H:135:ASP:O	1:H:267:PRO:HD3	2.22	0.40
1:A:236:MET:N	1:A:236:MET:SD	2.95	0.40
1:B:267:PRO:O	1:B:277:PHE:HB2	2.22	0.40
1:C:199:SER:OG	1:C:200:PRO:HD3	2.21	0.40
1:D:78:ARG:HD3	3:D:401:HOH:O	2.21	0.40
1:F:20:TYR:OH	1:F:41:LYS:HB3	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:144:LYS:HB2	1:H:147:ASP:HB2	2.03	0.40
1:H:199:SER:OG	1:H:200:PRO:HD3	2.21	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	293/292 (100%)	289 (99%)	4 (1%)	0	100	100
1	B	292/292 (100%)	288 (99%)	4 (1%)	0	100	100
1	C	288/292 (99%)	280 (97%)	7 (2%)	1 (0%)	44	44
1	D	287/292 (98%)	281 (98%)	5 (2%)	1 (0%)	44	44
1	E	288/292 (99%)	284 (99%)	4 (1%)	0	100	100
1	F	281/292 (96%)	277 (99%)	4 (1%)	0	100	100
1	G	288/292 (99%)	281 (98%)	6 (2%)	1 (0%)	44	44
1	H	284/292 (97%)	279 (98%)	5 (2%)	0	100	100
All	All	2301/2336 (98%)	2259 (98%)	39 (2%)	3 (0%)	55	57

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	2	ASP
1	G	74	ASP
1	C	75	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	258/255 (101%)	253 (98%)	5 (2%)	62	68
1	B	256/255 (100%)	255 (100%)	1 (0%)	93	95
1	C	254/255 (100%)	250 (98%)	4 (2%)	68	74
1	D	256/255 (100%)	253 (99%)	3 (1%)	75	81
1	E	253/255 (99%)	252 (100%)	1 (0%)	93	95
1	F	251/255 (98%)	250 (100%)	1 (0%)	93	95
1	G	253/255 (99%)	249 (98%)	4 (2%)	68	74
1	H	252/255 (99%)	249 (99%)	3 (1%)	75	81
All	All	2033/2040 (100%)	2011 (99%)	22 (1%)	78	83

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	59	LYS
1	A	77	VAL
1	A	236	MET
1	A	259	MET
1	B	282	ARG
1	C	3	THR
1	C	55	LEU
1	C	132	ASP
1	C	259	MET
1	D	17	ASN
1	D	172	ARG
1	D	216	LYS
1	E	18	ARG
1	F	289	LYS
1	G	14	ASP
1	G	21	LEU
1	G	67	ARG
1	G	128	ASN

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Mol	Chain	Res	Type
1	H	49	GLU
1	H	180	MET
1	H	247	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	5	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 8 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	292/292 (100%)	-0.23	7 (2%) 59 64	25, 35, 70, 209	0
1	B	290/292 (99%)	-0.22	2 (0%) 87 89	24, 38, 71, 125	0
1	C	289/292 (98%)	-0.17	6 (2%) 64 68	24, 40, 71, 115	0
1	D	289/292 (98%)	-0.18	3 (1%) 82 85	22, 39, 74, 136	0
1	E	289/292 (98%)	-0.35	2 (0%) 87 89	29, 42, 73, 133	0
1	F	283/292 (96%)	-0.29	1 (0%) 92 93	27, 40, 65, 89	0
1	G	289/292 (98%)	-0.17	3 (1%) 82 85	23, 45, 81, 135	0
1	H	287/292 (98%)	0.19	8 (2%) 53 60	29, 60, 96, 126	0
All	All	2308/2336 (98%)	-0.18	32 (1%) 75 79	22, 41, 81, 209	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	73	PRO	5.7
1	H	158	ILE	4.0
1	C	76	GLY	3.6
1	B	73	PRO	3.4
1	D	72	LYS	3.3
1	D	1	MET	3.0
1	A	76	GLY	2.9
1	H	154	LEU	2.9
1	H	253	LEU	2.9
1	C	3	THR	2.8
1	G	57	THR	2.8
1	A	236	MET	2.8
1	H	252	LYS	2.6
1	C	73	PRO	2.6
1	E	73	PRO	2.6
1	H	157	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	H	246	SER	2.5
1	F	291	HIS	2.5
1	A	292	ALA	2.3
1	A	70	ILE	2.3
1	A	74	ASP	2.3
1	A	291	HIS	2.3
1	G	73	PRO	2.3
1	C	75	GLY	2.2
1	H	167	VAL	2.2
1	D	4	SER	2.1
1	C	4	SER	2.1
1	H	51	ILE	2.1
1	G	51	ILE	2.0
1	E	76	GLY	2.0
1	B	4	SER	2.0
1	C	77	VAL	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. 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
2	K	G	301	1/1	0.99	0.11	0.07	44,44,44,44	0
2	K	D	301	1/1	0.99	0.10	-0.65	44,44,44,44	0
2	K	E	301	1/1	0.93	0.08	-0.82	57,57,57,57	0
2	K	C	301	1/1	1.00	0.07	-2.08	47,47,47,47	0
2	K	A	301	1/1	0.97	0.06	-2.50	51,51,51,51	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	K	F	301	1/1	0.98	0.06	-3.00	51,51,51,51	0
2	K	H	301	1/1	0.97	0.05	-3.72	60,60,60,60	0
2	K	B	301	1/1	0.99	0.04	-3.79	52,52,52,52	0

6.5 Other polymers [i](#)

There are no such residues in this entry.