



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:21 PM UTC

PDB ID : 4WFD / pdb_00004wfd
Title : Structure of the Rrp6-Rrp47-Mtr4 interaction
Authors : Schuch, B.; Conti, E.
Deposited on : 2014-09-14
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
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

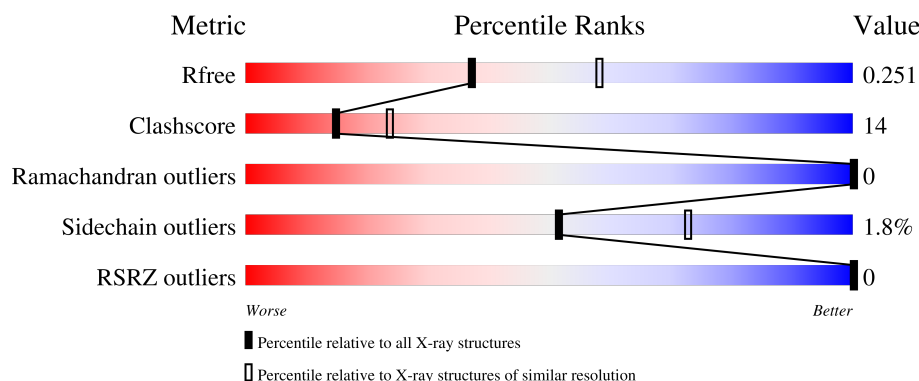
1 Overall quality at a glance

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)

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	115	
1	D	115	
1	G	115	
2	B	103	
2	E	103	

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Mol	Chain	Length	Quality of chain
2	H	103	61%28%10%
3	C	20	55%10%5%30%
3	F	20	35%5%60%
3	I	20	40%60%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4711 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 Exosome complex exonuclease RRP6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	91	Total	C	N	O	S	0	0	0
			702	435	116	147	4			
1	D	85	Total	C	N	O	S	0	0	0
			643	396	110	133	4			
1	G	82	Total	C	N	O	S	0	0	0
			617	382	108	123	4			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP Q12149
A	-2	PRO	-	expression tag	UNP Q12149
A	-1	ASP	-	expression tag	UNP Q12149
A	0	SER	-	expression tag	UNP Q12149
D	-3	GLY	-	expression tag	UNP Q12149
D	-2	PRO	-	expression tag	UNP Q12149
D	-1	ASP	-	expression tag	UNP Q12149
D	0	SER	-	expression tag	UNP Q12149
G	-3	GLY	-	expression tag	UNP Q12149
G	-2	PRO	-	expression tag	UNP Q12149
G	-1	ASP	-	expression tag	UNP Q12149
G	0	SER	-	expression tag	UNP Q12149

- Molecule 2 is a protein called Exosome complex protein LRP1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	100	Total	C	N	O	S	0	0	0
			803	512	133	154	4			
2	E	101	Total	C	N	O	S	0	0	0
			810	516	134	156	4			
2	H	93	Total	C	N	O	S	0	0	0
			733	469	120	140	4			

- Molecule 3 is a protein called ATP-dependent RNA helicase DOB1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	14	Total	C	N	O	0	0	0
			114	75	14	25			
3	F	8	Total	C	N	O	0	0	0
			71	47	8	16			
3	I	8	Total	C	N	O	0	0	0
			67	45	8	14			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	20	LYS	-	expression tag	UNP P47047
F	20	LYS	-	expression tag	UNP P47047
I	20	LYS	-	expression tag	UNP P47047

- Molecule 4 is YTTRIUM (III) ION (CCD ID: YT3) (formula: Y).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	6	Total	Y	0	0
			6	6		
4	B	1	Total	Y	0	0
			1	1		
4	C	1	Total	Y	0	0
			1	1		
4	D	5	Total	Y	0	0
			5	5		
4	E	1	Total	Y	0	0
			1	1		
4	F	1	Total	Y	0	0
			1	1		
4	G	6	Total	Y	0	0
			6	6		
4	I	1	Total	Y	0	0
			1	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	31	Total	O	0	0
			31	31		
5	B	20	Total	O	0	0
			20	20		

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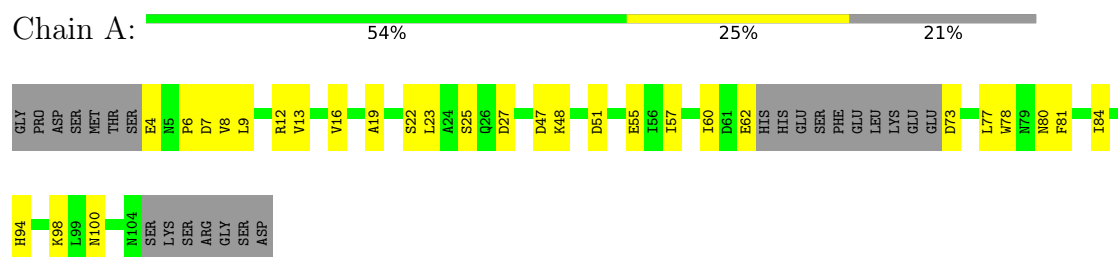
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	4	Total 4	O 4	0	0
5	D	18	Total 18	O 18	0	0
5	E	21	Total 21	O 21	0	0
5	G	20	Total 20	O 20	0	0
5	H	12	Total 12	O 12	0	0
5	I	3	Total 3	O 3	0	0

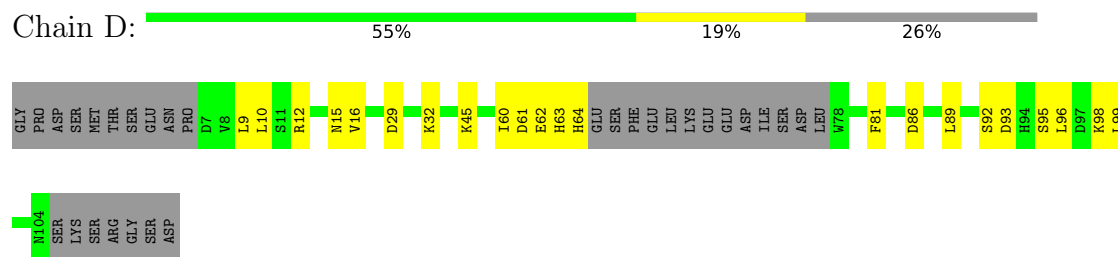
3 Residue-property plots [i](#)

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.

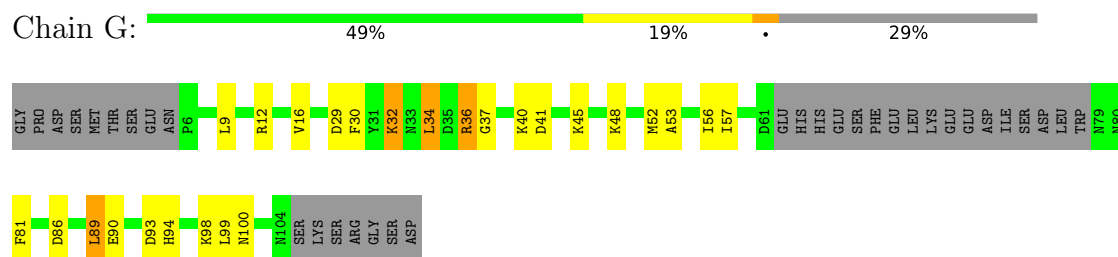
- Molecule 1: Exosome complex exonuclease RRP6



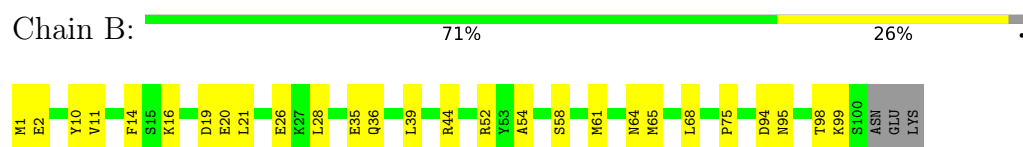
- Molecule 1: Exosome complex exonuclease RRP6



- Molecule 1: Exosome complex exonuclease RRP6



- Molecule 2: Exosome complex protein LRP1



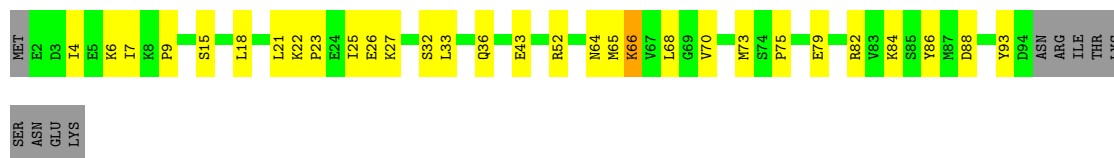
- Molecule 2: Exosome complex protein LRP1

Chain E:  67% 31% .



- Molecule 2: Exosome complex protein LRP1

Chain H:  61% 28% . 10%



- Molecule 3: ATP-dependent RNA helicase DOB1

Chain C:  55% 10% 5% 30%



- Molecule 3: ATP-dependent RNA helicase DOB1

Chain F:  35% 5% 60%



- Molecule 3: ATP-dependent RNA helicase DOB1

Chain I:  40% 60%



4 Data and refinement statistics

Property	Value	Source
Space group	P 3 1 2	Depositor
Cell constants a, b, c, α , β , γ	142.75Å 142.75Å 63.45Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	47.42 – 2.40 47.42 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.5 (47.42-2.40) 99.9 (47.42-2.40)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.55 (at 2.39Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.201 , 0.243 0.197 , 0.251	Depositor DCC
R_{free} test set	1470 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	56.3	Xtriage
Anisotropy	0.332	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 51.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.489 for -h,-k,l	Xtriage
Reported twinning fraction	0.500 for -h,-k,l	Depositor
Outliers	3 of 28898 reflections (0.010%)	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	4711	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.12% 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: YT3

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	0.62	0/708	1.05	0/956
1	D	0.64	0/647	1.05	1/871 (0.1%)
1	G	0.56	0/621	1.08	1/834 (0.1%)
2	B	0.61	0/812	0.98	0/1086
2	E	0.64	0/819	0.98	0/1097
2	H	0.57	0/741	1.01	1/993 (0.1%)
3	C	0.43	0/116	0.94	2/158 (1.3%)
3	F	0.50	0/72	0.74	0/96
3	I	0.48	0/68	0.73	0/91
All	All	0.60	0/4604	1.01	5/6182 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	D	60	ILE	N-CA-C	6.60	118.58	111.58
2	H	33	LEU	N-CA-C	6.31	117.82	111.07
1	G	34	LEU	N-CA-C	-5.90	104.77	111.14
3	C	13	THR	CA-C-N	5.75	126.07	119.92
3	C	13	THR	C-N-CA	5.75	126.07	119.92

There are no chirality outliers.

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	702	0	674	27	1
1	D	643	0	612	26	0
1	G	617	0	595	29	0
2	B	803	0	838	24	0
2	E	810	0	842	30	0
2	H	733	0	753	28	0
3	C	114	0	99	3	0
3	F	71	0	57	1	0
3	I	67	0	53	0	0
4	A	6	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	5	0	0	0	0
4	E	1	0	0	0	0
4	F	1	0	0	0	0
4	G	6	0	0	0	0
4	I	1	0	0	0	0
5	A	31	0	0	2	0
5	B	20	0	0	1	1
5	C	4	0	0	0	0
5	D	18	0	0	2	0
5	E	21	0	0	2	0
5	G	20	0	0	3	0
5	H	12	0	0	0	0
5	I	3	0	0	0	0
All	All	4711	0	4523	126	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 14.

All (126) 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:22:SER:HA	3:C:14:PRO:HB3	1.55	0.87
1:A:19:ALA:HB1	2:B:14:PHE:HA	1.67	0.77
1:A:27:ASP:OD1	5:A:301:HOH:O	2.04	0.74
2:H:64:ASN:O	2:H:68:LEU:HG	1.90	0.71
2:B:64:ASN:O	2:B:68:LEU:HG	1.91	0.70
1:G:29:ASP:OD1	1:G:32:LYS:NZ	2.24	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:78:TRP:HZ3	2:B:61:MET:HE2	1.57	0.70
2:H:65:MET:HB3	2:H:73:MET:HE1	1.74	0.69
1:G:9:LEU:HD23	2:H:52:ARG:HH11	1.57	0.69
1:D:62:GLU:O	1:D:64:HIS:ND1	2.26	0.68
1:A:8:VAL:O	1:A:12:ARG:HG2	1.95	0.67
1:D:99:LEU:HD11	2:E:44:ARG:HA	1.76	0.67
1:D:12:ARG:NH1	2:E:20:GLU:OE1	2.28	0.66
1:A:7:ASP:OD1	2:B:52:ARG:NH1	2.30	0.65
1:A:78:TRP:CZ3	2:B:61:MET:HE2	2.31	0.64
2:B:54:ALA:O	2:B:58:SER:OG	2.14	0.64
1:D:96:LEU:HD21	2:E:89:LYS:HB3	1.80	0.64
1:G:89:LEU:HD23	2:H:82:ARG:CZ	2.28	0.63
1:D:81:PHE:HE2	2:E:79:GLU:HG3	1.63	0.63
2:E:43:GLU:OE2	2:E:46:LYS:HD2	2.00	0.62
1:G:56:ILE:HD11	2:H:25:ILE:HD12	1.82	0.61
1:G:29:ASP:OD2	5:G:301:HOH:O	2.16	0.61
2:B:16:LYS:O	2:B:20:GLU:HG3	2.01	0.60
2:E:12:ARG:NH1	5:E:303:HOH:O	2.33	0.59
2:H:32:SER:O	2:H:36:GLN:HG3	2.03	0.58
2:B:94:ASP:O	2:B:98:THR:HG23	2.03	0.58
1:A:4:GLU:HG3	1:A:6:PRO:HD3	1.84	0.58
1:D:29:ASP:O	1:D:32:LYS:HG2	2.03	0.58
1:D:12:ARG:HD2	2:E:20:GLU:HB3	1.87	0.57
1:D:89:LEU:HD22	2:E:86:TYR:CD2	2.40	0.56
2:H:84:LYS:NZ	2:H:88:ASP:OD2	2.38	0.56
1:G:16:VAL:HG21	2:H:21:LEU:HD22	1.89	0.55
1:D:81:PHE:CE2	2:E:79:GLU:HG3	2.42	0.55
2:B:35:GLU:O	2:B:39:LEU:HG	2.06	0.55
1:G:99:LEU:HD21	2:H:43:GLU:HB3	1.87	0.55
2:E:71:LYS:NZ	5:E:305:HOH:O	2.39	0.55
1:G:32:LYS:NZ	5:G:306:HOH:O	2.32	0.55
2:B:28:LEU:O	2:B:36:GLN:NE2	2.40	0.54
2:E:74:SER:N	2:E:75:PRO:HD2	2.23	0.54
1:A:16:VAL:HG21	2:B:21:LEU:HD22	1.89	0.54
2:B:10:TYR:OH	3:C:17:LEU:HD13	2.07	0.54
1:G:36:ARG:HH11	1:G:40:LYS:HD3	1.74	0.53
1:G:36:ARG:HG3	1:G:37:GLY:N	2.23	0.53
2:B:65:MET:HE1	2:B:75:PRO:HG2	1.92	0.52
1:G:12:ARG:O	1:G:16:VAL:HG23	2.10	0.51
1:A:25:SER:HB3	3:C:14:PRO:HG2	1.92	0.51
2:B:1:MET:O	2:B:2:GLU:HB2	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:93:ASP:HB2	2:H:86:TYR:HE1	1.75	0.51
2:B:95:ASN:O	2:B:99:LYS:HG2	2.09	0.51
2:E:94:ASP:O	2:E:98:THR:HG23	2.11	0.51
1:A:48:LYS:HE3	2:B:19:ASP:OD1	2.11	0.51
2:B:39:LEU:HB2	5:B:302:HOH:O	2.11	0.51
2:E:37:LEU:HD23	2:E:49:LEU:HD23	1.93	0.51
2:H:6:LYS:O	2:H:9:PRO:HD2	2.11	0.51
1:A:60:ILE:HG22	1:A:62:GLU:HG2	1.93	0.50
2:E:6:LYS:O	2:E:9:PRO:HD2	2.11	0.50
1:G:93:ASP:HB2	2:H:86:TYR:CE1	2.47	0.50
1:A:81:PHE:O	1:A:84:ILE:HG22	2.12	0.49
2:E:46:LYS:O	2:E:50:ILE:HG13	2.12	0.49
1:G:30:PHE:O	1:G:34:LEU:HB2	2.13	0.49
2:H:4:ILE:O	2:H:7:ILE:HG22	2.10	0.49
2:H:65:MET:HE1	2:H:75:PRO:HG2	1.95	0.49
1:G:48:LYS:HD2	2:H:18:LEU:HB3	1.95	0.49
1:D:99:LEU:HD21	2:E:44:ARG:N	2.28	0.48
1:D:62:GLU:OE1	2:E:32:SER:HB2	2.14	0.48
2:E:37:LEU:CD2	2:E:49:LEU:HD23	2.44	0.48
1:D:45:LYS:NZ	5:D:302:HOH:O	2.27	0.48
2:H:73:MET:HA	2:H:73:MET:HE2	1.96	0.47
1:D:89:LEU:O	1:D:93:ASP:N	2.38	0.47
2:E:86:TYR:HA	2:E:89:LYS:HD3	1.97	0.47
2:B:26:GLU:HA	2:B:26:GLU:OE1	2.14	0.47
1:G:52:MET:CE	2:H:25:ILE:HG13	2.45	0.47
1:D:12:ARG:O	1:D:16:VAL:HG23	2.15	0.47
1:G:41:ASP:OD1	1:G:45:LYS:HE2	2.15	0.46
1:A:94:HIS:CD2	1:A:98:LYS:HE2	2.50	0.46
1:G:45:LYS:HD2	2:H:15:SER:HB2	1.97	0.46
2:B:95:ASN:O	2:B:98:THR:OG1	2.24	0.46
2:H:66:LYS:HA	2:H:66:LYS:HD2	1.77	0.46
1:A:19:ALA:HB1	2:B:14:PHE:CA	2.41	0.46
1:D:12:ARG:HD3	1:D:12:ARG:HA	1.64	0.46
1:A:9:LEU:O	1:A:13:VAL:HG23	2.15	0.46
2:H:65:MET:HA	2:H:68:LEU:HD12	1.97	0.45
1:A:100:ASN:HB3	5:A:307:HOH:O	2.16	0.45
1:D:95:SER:OG	2:E:47:LEU:HB2	2.17	0.45
1:G:36:ARG:HG3	1:G:37:GLY:H	1.81	0.45
1:A:9:LEU:HD23	2:B:52:ARG:HH11	1.81	0.45
1:D:61:ASP:N	1:D:61:ASP:OD1	2.50	0.45
1:A:19:ALA:CB	2:B:14:PHE:HA	2.42	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:77:LEU:O	1:A:80:ASN:HB2	2.16	0.45
1:D:86:ASP:OD1	2:E:82:ARG:NH2	2.48	0.44
2:H:68:LEU:HB2	2:H:70:VAL:HG23	1.99	0.44
1:D:62:GLU:O	1:D:64:HIS:CE1	2.71	0.44
2:E:61:MET:HE3	2:E:65:MET:SD	2.58	0.44
1:A:57:ILE:HA	1:A:60:ILE:HD12	1.99	0.44
1:D:98:LYS:NZ	2:E:43:GLU:OE1	2.47	0.44
2:H:23:PRO:O	2:H:27:LYS:HG3	2.18	0.44
1:A:47:ASP:O	1:A:51:ASP:N	2.44	0.44
2:B:44:ARG:O	2:B:44:ARG:NH1	2.47	0.44
1:G:36:ARG:NH1	1:G:40:LYS:HD3	2.31	0.44
1:D:10:LEU:HD13	3:F:6:LEU:HD22	2.00	0.43
1:D:9:LEU:HD21	2:E:56:VAL:HG21	1.99	0.43
1:D:15:ASN:OD1	5:D:301:HOH:O	2.21	0.43
1:G:52:MET:HE3	2:H:22:LYS:HA	2.00	0.43
1:G:86:ASP:O	1:G:90:GLU:N	2.46	0.43
1:A:60:ILE:CD1	1:A:84:ILE:HD11	2.49	0.43
1:D:89:LEU:HA	1:D:92:SER:HB3	2.00	0.42
1:G:98:LYS:HD2	2:H:43:GLU:OE1	2.17	0.42
1:A:60:ILE:HD13	1:A:84:ILE:HD11	2.00	0.42
1:A:73:ASP:O	1:A:77:LEU:N	2.52	0.42
1:A:23:LEU:HD12	1:A:23:LEU:O	2.20	0.42
1:D:99:LEU:CD2	2:E:43:GLU:HB3	2.49	0.42
1:G:100:ASN:ND2	5:G:308:HOH:O	2.49	0.42
2:B:11:VAL:O	2:B:14:PHE:HB3	2.19	0.42
1:G:53:ALA:O	1:G:57:ILE:HG13	2.20	0.42
2:E:21:LEU:HG	2:E:25:ILE:HG12	2.01	0.41
1:G:99:LEU:CD2	2:H:43:GLU:HB3	2.50	0.41
1:G:81:PHE:HE2	2:H:79:GLU:HG3	1.84	0.41
2:E:65:MET:HG2	2:E:70:VAL:HG21	2.03	0.41
1:A:77:LEU:C	1:A:77:LEU:HD13	2.45	0.41
2:H:26:GLU:OE1	2:H:26:GLU:HA	2.19	0.41
2:E:52:ARG:O	2:E:56:VAL:HG23	2.21	0.41
1:G:99:LEU:HB3	2:H:93:TYR:CE1	2.57	0.40
1:G:94:HIS:CE1	1:G:98:LYS:HZ2	2.39	0.40
1:D:62:GLU:HG2	1:D:63:HIS:N	2.37	0.40
2:E:1:MET:O	2:E:2:GLU:HB2	2.22	0.40
2:E:16:LYS:NZ	2:E:20:GLU:OE2	2.40	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)
1:A:55:GLU:OE1	5:B:301:HOH:O[5_557]	2.18	0.02

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	87/115 (76%)	86 (99%)	1 (1%)	0	100	100
1	D	81/115 (70%)	80 (99%)	1 (1%)	0	100	100
1	G	78/115 (68%)	78 (100%)	0	0	100	100
2	B	98/103 (95%)	96 (98%)	2 (2%)	0	100	100
2	E	99/103 (96%)	98 (99%)	1 (1%)	0	100	100
2	H	91/103 (88%)	91 (100%)	0	0	100	100
3	C	12/20 (60%)	11 (92%)	1 (8%)	0	100	100
3	F	6/20 (30%)	6 (100%)	0	0	100	100
3	I	6/20 (30%)	6 (100%)	0	0	100	100
All	All	558/714 (78%)	552 (99%)	6 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	D	71/104 (68%)	71 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	66/104 (64%)	63 (96%)	3 (4%)	24	42
2	E	91/96 (95%)	90 (99%)	1 (1%)	65	82
2	H	80/96 (83%)	79 (99%)	1 (1%)	61	80
3	C	12/20 (60%)	11 (92%)	1 (8%)	10	17
3	F	8/20 (40%)	8 (100%)	0	100	100
3	I	7/20 (35%)	7 (100%)	0	100	100
All	All	335/460 (73%)	329 (98%)	6 (2%)	51	73

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	C	17	LEU
2	E	3	ASP
1	G	32	LYS
1	G	36	ARG
1	G	89	LEU
2	H	66	LYS

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

Mol	Chain	Res	Type
1	G	26	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 oligosaccharides in this entry.

5.6 Ligand geometry

Of 22 ligands modelled in this entry, 22 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

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 [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	91/115 (79%)	-1.17	0 100 100	43, 61, 88, 99	0
1	D	85/115 (73%)	-1.27	0 100 100	42, 61, 100, 116	0
1	G	82/115 (71%)	-0.84	0 100 100	61, 73, 101, 114	0
2	B	100/103 (97%)	-1.28	0 100 100	43, 65, 95, 101	0
2	E	101/103 (98%)	-1.34	0 100 100	44, 62, 95, 112	0
2	H	93/103 (90%)	-1.09	0 100 100	63, 76, 95, 110	0
3	C	14/20 (70%)	-0.61	0 100 100	55, 88, 100, 111	0
3	F	8/20 (40%)	-1.46	0 100 100	52, 56, 79, 88	0
3	I	8/20 (40%)	-1.47	0 100 100	75, 86, 92, 98	0
All	All	582/714 (81%)	-1.17	0 100 100	42, 69, 97, 116	0

There are no RSRZ outliers to report.

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	YT3	G	204	1/1	0.98	0.10	241,241,241,241	0
4	YT3	A	203	1/1	0.99	0.12	109,109,109,109	0
4	YT3	A	204	1/1	0.99	0.08	78,78,78,78	1
4	YT3	A	205	1/1	0.99	0.10	90,90,90,90	1
4	YT3	A	206	1/1	0.99	0.08	98,98,98,98	1
4	YT3	D	205	1/1	0.99	0.11	206,206,206,206	0
4	YT3	G	203	1/1	0.99	0.06	117,117,117,117	0
4	YT3	A	202	1/1	0.99	0.06	189,189,189,189	0
4	YT3	G	206	1/1	0.99	0.08	190,190,190,190	0
4	YT3	D	202	1/1	1.00	0.01	69,69,69,69	0
4	YT3	D	203	1/1	1.00	0.04	73,73,73,73	1
4	YT3	D	204	1/1	1.00	0.04	74,74,74,74	1
4	YT3	A	201	1/1	1.00	0.01	69,69,69,69	0
4	YT3	E	201	1/1	1.00	0.01	56,56,56,56	1
4	YT3	F	101	1/1	1.00	0.02	74,74,74,74	1
4	YT3	G	201	1/1	1.00	0.01	85,85,85,85	0
4	YT3	G	202	1/1	1.00	0.01	87,87,87,87	0
4	YT3	B	201	1/1	1.00	0.01	56,56,56,56	1
4	YT3	C	101	1/1	1.00	0.02	78,78,78,78	1
4	YT3	G	205	1/1	1.00	0.04	145,145,145,145	0
4	YT3	D	201	1/1	1.00	0.02	52,52,52,52	0
4	YT3	I	101	1/1	1.00	0.04	121,121,121,121	0

6.5 Other polymers ⓘ

There are no such residues in this entry.