



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 03:48 PM GMT

PDB ID : 3SAG
Title : Crystal structure of the human RRP6 catalytic domain with D313N mutation
in the active site
Authors : Januszyk, K.; Liu, Q.; Lima, C.D.
Deposited on : 2011-06-02
Resolution : 2.70 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

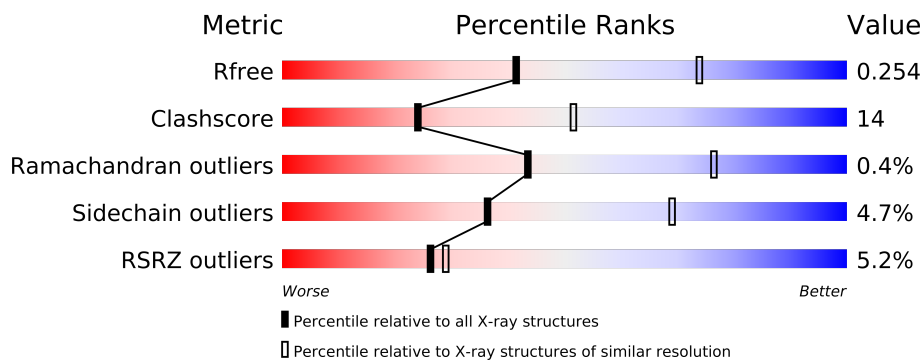
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.70 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}	66092	1557 (2.70-2.70)
Clashscore	79885	1939 (2.70-2.70)
Ramachandran outliers	78287	1905 (2.70-2.70)
Sidechain outliers	78261	1905 (2.70-2.70)
RSRZ outliers	66119	1559 (2.70-2.70)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	428	
1	B	428	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6537 atoms, of which 0 are hydrogen and 0 are deuterium.

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 component 10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	381	Total	C	N	O	S	0	0	0
			3168	2029	548	573	18			
1	B	380	Total	C	N	O	S	0	0	0
			3137	2010	537	572	18			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	179	SER	-	EXPRESSION TAG	UNP Q01780
A	313	ASN	ASP	ENGINEERED MUTATION	UNP Q01780
B	179	SER	-	EXPRESSION TAG	UNP Q01780
B	313	ASN	ASP	ENGINEERED MUTATION	UNP Q01780

- Molecule 2 is YTTRIUM (III) ION (three-letter code: YT3) (formula: Y).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Y	0	0
			1	1		
2	A	2	Total	Y	0	0
			2	2		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Mg	0	0
			1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	114	Total	O	0	0
			114	114		

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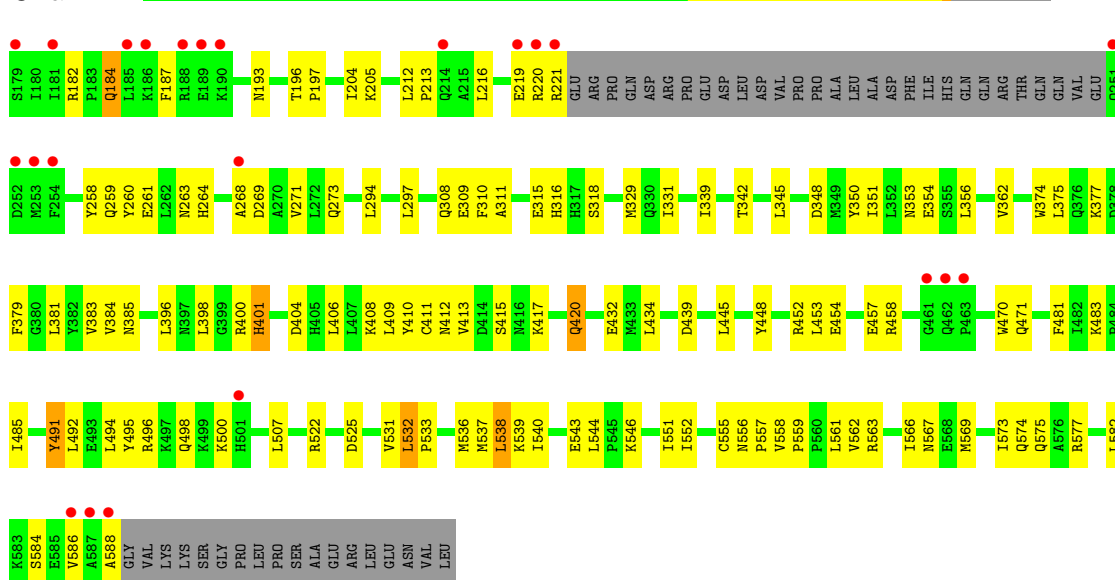
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	114	Total 114	O 114	0	0

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

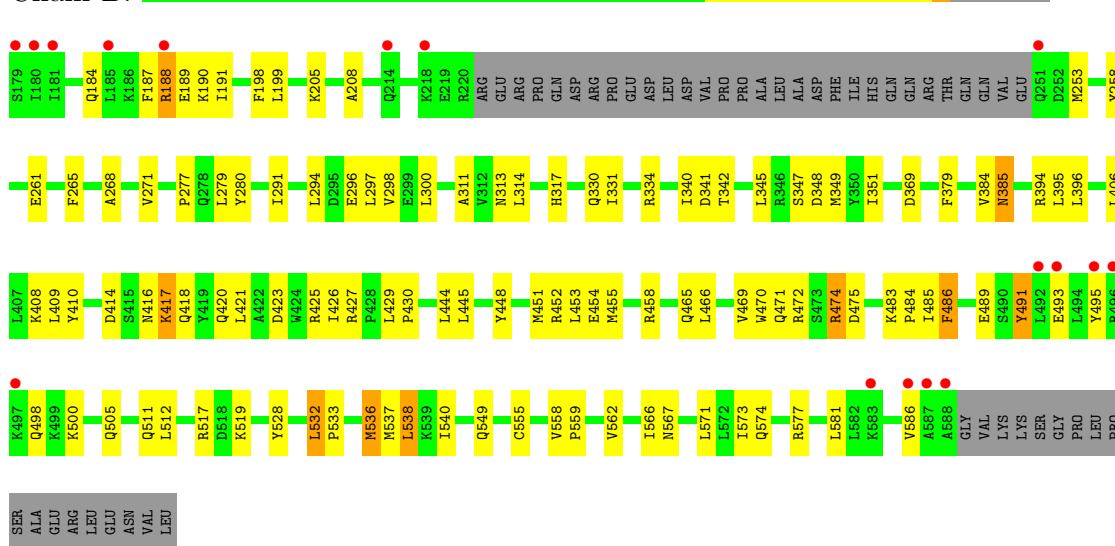
• Molecule 1: Exosome component 10

Chain A:



• Molecule 1: Exosome component 10

Chain B:



4 Data and refinement statistics

Property	Value	Source
Space group	P 4	Depositor
Cell constants a, b, c, α , β , γ	138.80Å 138.80Å 60.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.70 – 2.70 34.70 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.8 (34.70-2.70) 99.8 (34.70-2.70)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.42 (at 2.68Å)	Xtriage
Refinement program	CNS 1.3	Depositor
R, R_{free}	0.208 , 0.262 0.205 , 0.254	Depositor DCC
R_{free} test set	1588 reflections (5.24%)	DCC
Wilson B-factor (Å ²)	53.4	Xtriage
Anisotropy	0.580	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 31.1	EDS
Estimated twinning fraction	0.033 for h,-k,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	1 of 31952 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6537	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.68% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, 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	0.36	0/3245	0.58	0/4394
1	B	0.38	0/3213	0.59	0/4355
All	All	0.37	0/6458	0.58	0/8749

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3168	0	3162	88	0
1	B	3137	0	3110	90	0
2	A	2	0	0	0	0
2	B	1	0	0	0	0
3	B	1	0	0	0	0
4	A	114	0	0	4	0
4	B	114	0	0	2	0
All	All	6537	0	6272	178	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including

hydrogens) of the entry. The overall clashscore for this entry is 14.

All (178) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:416:ASN:HB2	1:B:418:GLN:HE21	1.36	0.88
1:B:394:ARG:HH22	1:B:472:ARG:HE	1.18	0.88
1:B:571:LEU:HA	1:B:574:GLN:HE21	1.43	0.83
1:A:494:LEU:HG	1:A:498:GLN:HE21	1.44	0.81
1:B:394:ARG:NH2	1:B:472:ARG:HE	1.82	0.78
1:B:396:LEU:HD21	1:B:455:MET:HG2	1.66	0.78
1:A:258:TYR:HD1	1:A:261:GLU:HG3	1.50	0.76
1:B:416:ASN:HB2	1:B:418:GLN:NE2	2.02	0.74
1:A:362:VAL:HG22	1:A:385:ASN:HD21	1.52	0.74
1:B:297:LEU:HD22	1:B:345:LEU:HD13	1.70	0.73
1:A:485:ILE:O	1:A:485:ILE:HD12	1.88	0.73
1:B:258:TYR:HD1	1:B:261:GLU:HG3	1.53	0.73
1:A:294:LEU:HD13	1:A:348:ASP:HB3	1.71	0.72
1:B:205:LYS:HG2	1:B:208:ALA:HB2	1.72	0.71
1:A:533:PRO:HG2	1:A:536:MET:HG3	1.72	0.71
1:B:505:GLN:HE21	1:B:581:LEU:HG	1.57	0.69
1:A:556:ASN:HA	4:A:127:HOH:O	1.93	0.68
1:B:417:LYS:HD2	1:B:417:LYS:H	1.58	0.68
1:A:552:ILE:HG23	1:A:558:VAL:HG21	1.75	0.68
1:A:182:ARG:HB3	1:A:182:ARG:NH1	2.10	0.67
1:A:362:VAL:CG2	1:A:385:ASN:HD21	2.07	0.67
1:B:491:TYR:HD1	1:B:491:TYR:H	1.45	0.64
1:B:191:ILE:HD12	1:B:191:ILE:N	2.12	0.64
1:A:538:LEU:HD22	1:A:538:LEU:O	1.97	0.64
1:B:187:PHE:HD2	1:B:191:ILE:HD11	1.62	0.63
1:A:259:GLN:HE21	1:A:263:ASN:HD21	1.45	0.63
1:A:448:TYR:CE1	1:A:452:ARG:HD2	2.34	0.62
1:A:269:ASP:O	1:A:273:GLN:HG3	2.00	0.61
1:A:532:LEU:HD13	1:A:537:MET:HB2	1.81	0.61
1:A:184:GLN:HA	1:A:187:PHE:CD2	2.37	0.59
1:B:296:GLU:HA	1:B:296:GLU:OE1	2.01	0.59
1:B:294:LEU:HD13	1:B:348:ASP:HB3	1.85	0.59
1:A:213:PRO:HG2	1:A:216:LEU:HD12	1.83	0.59
1:B:573:ILE:O	1:B:577:ARG:HG3	2.02	0.59
1:A:362:VAL:HG22	1:A:385:ASN:ND2	2.16	0.59
1:B:495:TYR:HB3	1:B:500:LYS:CB	2.33	0.59
1:B:279:LEU:HD13	1:B:280:TYR:O	2.03	0.59
1:B:340:ILE:HG23	1:B:345:LEU:HD12	1.85	0.58
1:A:353:ASN:ND2	1:A:381:LEU:HD23	2.19	0.58
1:A:557:PRO:HA	4:A:127:HOH:O	2.04	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:398:LEU:C	1:A:400:ARG:H	2.07	0.57
1:B:559:PRO:HG2	1:B:562:VAL:CG2	2.34	0.57
1:A:552:ILE:HD13	1:A:563:ARG:NH2	2.21	0.56
1:B:294:LEU:O	1:B:298:VAL:HG23	2.06	0.56
1:A:533:PRO:HD2	1:A:536:MET:HE2	1.88	0.56
1:A:182:ARG:HH11	1:A:182:ARG:HB3	1.70	0.56
1:B:532:LEU:HD21	1:B:536:MET:HB3	1.88	0.56
1:A:184:GLN:HE21	1:A:184:GLN:H	1.53	0.55
1:A:268:ALA:O	1:A:271:VAL:HG22	2.06	0.55
1:B:474:ARG:HD2	1:B:474:ARG:C	2.27	0.55
1:B:394:ARG:HH22	1:B:472:ARG:NE	1.96	0.55
1:A:351:ILE:O	1:A:354:GLU:HG3	2.06	0.54
1:A:495:TYR:O	1:A:500:LYS:HB2	2.07	0.54
1:B:258:TYR:CD1	1:B:261:GLU:HG3	2.39	0.54
1:A:491:TYR:H	1:A:491:TYR:HD1	1.55	0.54
1:A:453:LEU:O	1:A:457:GLU:HG3	2.08	0.54
1:A:398:LEU:C	1:A:400:ARG:N	2.61	0.53
1:B:420:GLN:O	1:B:421:LEU:HB2	2.07	0.53
1:A:398:LEU:HD22	1:A:401:HIS:NE2	2.24	0.53
1:B:559:PRO:HG2	1:B:562:VAL:HB	1.91	0.53
1:B:188:ARG:HG3	1:B:347:SER:HB3	1.91	0.53
1:A:409:LEU:HD12	1:A:409:LEU:O	2.09	0.52
1:A:540:ILE:HD13	1:A:551:ILE:HG23	1.91	0.52
1:A:212:LEU:HD12	1:A:213:PRO:HD2	1.92	0.52
1:B:533:PRO:HD2	1:B:536:MET:HE2	1.91	0.52
1:A:311:ALA:O	1:A:331:ILE:HA	2.09	0.52
1:B:472:ARG:O	1:B:475:ASP:HB2	2.10	0.52
1:B:505:GLN:HE21	1:B:581:LEU:CG	2.22	0.52
1:B:317:HIS:CD2	1:B:425:ARG:HD3	2.45	0.52
1:B:484:PRO:HB2	1:B:486:PHE:CZ	2.45	0.51
1:B:445:LEU:O	1:B:448:TYR:HB3	2.10	0.51
1:A:182:ARG:HH11	1:A:182:ARG:CB	2.23	0.51
1:B:455:MET:SD	1:B:466:LEU:HD12	2.51	0.51
1:B:489:GLU:O	1:B:493:GLU:HG3	2.10	0.51
1:B:517:ARG:NE	1:B:528:TYR:HD2	2.08	0.51
1:B:369:ASP:HB2	4:B:162:HOH:O	2.10	0.51
1:A:494:LEU:HG	1:A:498:GLN:NE2	2.21	0.50
1:B:491:TYR:CD1	1:B:491:TYR:N	2.75	0.50
1:A:193:ASN:O	1:A:522:ARG:HD3	2.11	0.50
1:A:555:CYS:HB2	1:A:558:VAL:HG22	1.93	0.50
1:B:536:MET:HE3	1:B:559:PRO:CD	2.42	0.50
1:B:536:MET:CE	1:B:559:PRO:HD3	2.41	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:532:LEU:CD2	1:B:536:MET:HB3	2.41	0.50
1:B:495:TYR:HA	1:B:498:GLN:HB3	1.93	0.50
1:A:260:TYR:CE1	1:A:264:HIS:CE1	3.00	0.49
1:A:377:LYS:HG3	1:A:481:PHE:CE2	2.48	0.49
1:A:182:ARG:NH2	1:A:525:ASP:O	2.46	0.49
1:A:310:PHE:HD1	1:A:331:ILE:HG23	1.77	0.49
1:B:454:GLU:O	1:B:458:ARG:HG2	2.12	0.49
1:B:265:PHE:CZ	1:B:385:ASN:HA	2.48	0.48
1:B:191:ILE:CD1	1:B:191:ILE:N	2.76	0.48
1:A:339:ILE:HD12	1:A:434:LEU:HD22	1.95	0.48
1:A:398:LEU:O	1:A:400:ARG:N	2.46	0.48
1:B:485:ILE:O	1:B:485:ILE:HG23	2.14	0.48
1:B:511:GLN:HG3	1:B:512:LEU:N	2.28	0.48
1:A:259:GLN:HE21	1:A:263:ASN:ND2	2.11	0.47
1:B:395:LEU:HD12	1:B:469:VAL:HG22	1.96	0.47
1:A:362:VAL:HA	1:A:385:ASN:ND2	2.29	0.47
1:A:445:LEU:O	1:A:448:TYR:HB3	2.14	0.47
1:B:555:CYS:HB2	1:B:558:VAL:HG22	1.95	0.47
1:A:408:LYS:HE2	1:A:412:ASN:OD1	2.15	0.47
1:B:448:TYR:CE1	1:B:452:ARG:HD2	2.49	0.47
1:B:190:LYS:HG2	1:B:191:ILE:N	2.29	0.47
1:A:491:TYR:CD1	1:A:491:TYR:N	2.82	0.47
1:B:313:ASN:C	1:B:314:LEU:HD12	2.35	0.47
1:B:409:LEU:HD12	1:B:409:LEU:O	2.13	0.47
1:A:582:LEU:O	1:A:586:VAL:HG23	2.15	0.47
1:A:559:PRO:HG2	1:A:562:VAL:CG2	2.44	0.47
1:B:349:MET:HB2	1:B:379:PHE:HB3	1.97	0.46
1:B:334:ARG:HD3	1:B:445:LEU:HD12	1.97	0.46
1:B:536:MET:O	1:B:540:ILE:HG13	2.15	0.46
1:A:411:CYS:HB2	1:A:413:VAL:HG23	1.98	0.46
1:B:384:VAL:O	1:B:385:ASN:HB3	2.16	0.45
1:A:404:ASP:OD1	1:A:415:SER:HB3	2.16	0.45
1:B:408:LYS:HE3	1:B:414:ASP:OD1	2.16	0.45
1:A:546:LYS:HG3	1:A:577:ARG:NH1	2.32	0.45
1:B:341:ASP:OD1	1:B:427:ARG:NH1	2.49	0.45
1:B:291:ILE:HD12	1:B:300:LEU:HD22	1.98	0.45
1:A:353:ASN:HD21	1:A:381:LEU:HA	1.82	0.45
1:B:423:ASP:O	1:B:426:ILE:HG22	2.17	0.45
1:A:329:MET:HE3	1:A:375:LEU:HD21	1.98	0.45
1:B:342:THR:HG22	1:B:349:MET:SD	2.57	0.45
1:A:182:ARG:HH12	1:A:184:GLN:HB2	1.82	0.44
1:A:384:VAL:O	1:A:385:ASN:HB3	2.17	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:187:PHE:CD2	1:B:191:ILE:HD11	2.49	0.44
1:B:536:MET:HE3	1:B:559:PRO:HD3	1.99	0.44
1:B:268:ALA:O	1:B:271:VAL:HG22	2.18	0.44
1:B:451:MET:O	1:B:455:MET:HG3	2.18	0.44
1:B:536:MET:HG2	1:B:555:CYS:SG	2.58	0.44
1:B:406:LEU:HD22	1:B:410:TYR:HE1	1.82	0.44
1:B:416:ASN:O	1:B:418:GLN:HG2	2.18	0.44
1:B:486:PHE:CD1	1:B:486:PHE:N	2.85	0.44
1:A:196:THR:O	1:A:483:LYS:HD2	2.18	0.43
1:B:505:GLN:NE2	1:B:581:LEU:HG	2.30	0.43
1:B:533:PRO:HD2	1:B:536:MET:CE	2.48	0.43
1:A:377:LYS:HG2	1:A:377:LYS:O	2.17	0.43
1:B:465:GLN:O	1:B:469:VAL:HG23	2.18	0.43
1:A:297:LEU:HD22	1:A:345:LEU:HD13	2.00	0.43
1:A:350:TYR:CD1	1:A:351:ILE:N	2.87	0.43
1:A:316:HIS:HB2	1:A:374:TRP:CZ3	2.53	0.43
1:A:406:LEU:O	1:A:410:TYR:HB2	2.18	0.43
1:A:562:VAL:HG13	1:A:569:MET:SD	2.59	0.43
1:A:342:THR:HB	1:A:379:PHE:CZ	2.54	0.43
1:A:492:LEU:O	1:A:496:ARG:HB2	2.18	0.43
1:A:445:LEU:HG	4:A:31:HOH:O	2.17	0.42
1:A:454:GLU:O	1:A:458:ARG:HG3	2.19	0.42
1:A:204:ILE:HG23	1:A:204:ILE:O	2.19	0.42
1:A:318:SER:HB2	4:A:172:HOH:O	2.18	0.42
1:B:297:LEU:HG	1:B:351:ILE:HD11	2.01	0.42
1:B:423:ASP:OD1	1:B:425:ARG:HB2	2.19	0.42
1:B:199:LEU:HD13	1:B:253:MET:HE2	2.02	0.42
1:B:340:ILE:HD12	1:B:340:ILE:N	2.35	0.42
1:B:311:ALA:O	1:B:331:ILE:HA	2.19	0.42
1:B:429:LEU:HA	1:B:430:PRO:HD2	1.91	0.42
1:A:584:SER:O	1:A:588:ALA:HB3	2.20	0.42
1:B:261:GLU:CD	1:B:261:GLU:H	2.23	0.42
1:B:566:ILE:HG23	1:B:567:ASN:N	2.35	0.42
1:B:559:PRO:HG2	1:B:562:VAL:CB	2.49	0.42
1:A:566:ILE:HG23	1:A:567:ASN:N	2.35	0.42
1:B:538:LEU:O	1:B:538:LEU:HD23	2.19	0.42
1:B:198:PHE:HB2	1:B:483:LYS:HG2	2.02	0.41
1:A:573:ILE:O	1:A:577:ARG:HG3	2.20	0.41
1:A:546:LYS:HG3	1:A:577:ARG:HH12	1.85	0.41
1:A:539:LYS:HB2	1:A:539:LYS:HE3	1.81	0.41
1:A:196:THR:HG23	1:A:197:PRO:HD2	2.01	0.41
1:B:341:ASP:CG	1:B:427:ARG:NH1	2.74	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:531:VAL:HG22	1:A:561:LEU:CD1	2.51	0.41
1:A:219:GLU:C	1:A:221:ARG:H	2.24	0.41
1:B:532:LEU:HD22	1:B:537:MET:HB2	2.03	0.41
1:A:205:LYS:HZ1	1:A:261:GLU:CD	2.24	0.40
1:A:258:TYR:CD1	1:A:261:GLU:HG3	2.41	0.40
1:B:536:MET:HE1	1:B:559:PRO:HD3	2.03	0.40
1:A:417:LYS:HG2	1:A:420:GLN:HE22	1.85	0.40
1:A:543:GLU:O	1:A:544:LEU:C	2.60	0.40
1:A:315:GLU:OE1	1:A:315:GLU:HA	2.21	0.40
1:A:396:LEU:O	1:A:398:LEU:HG	2.21	0.40
1:B:334:ARG:HG2	4:B:19:HOH:O	2.22	0.40
1:A:356:LEU:O	1:A:383:VAL:HA	2.21	0.40

There are no symmetry-related clashes.

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	377/428 (88%)	344 (91%)	32 (8%)	1 (0%)	50	82
1	B	376/428 (88%)	352 (94%)	22 (6%)	2 (0%)	38	70
All	All	753/856 (88%)	696 (92%)	54 (7%)	3 (0%)	43	76

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	220	ARG
1	B	385	ASN
1	B	277	PRO

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	353/395 (89%)	338 (96%)	15 (4%)	40	73
1	B	348/395 (88%)	330 (95%)	18 (5%)	32	63
All	All	701/790 (89%)	668 (95%)	33 (5%)	36	69

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	184	GLN
1	A	308	GLN
1	A	309	GLU
1	A	401	HIS
1	A	420	GLN
1	A	432	GLU
1	A	439	ASP
1	A	470	TRP
1	A	471	GLN
1	A	491	TYR
1	A	507	LEU
1	A	532	LEU
1	A	538	LEU
1	A	574	GLN
1	A	575	GLN
1	B	184	GLN
1	B	188	ARG
1	B	189	GLU
1	B	330	GLN
1	B	417	LYS
1	B	444	LEU
1	B	453	LEU
1	B	470	TRP
1	B	471	GLN
1	B	474	ARG
1	B	486	PHE
1	B	491	TYR
1	B	519	LYS
1	B	532	LEU
1	B	536	MET
1	B	538	LEU
1	B	549	GLN
1	B	586	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	184	GLN
1	A	259	GLN
1	A	385	ASN
1	A	465	GLN
1	A	467	GLN
1	A	471	GLN
1	A	498	GLN
1	A	505	GLN
1	A	575	GLN
1	B	184	GLN
1	B	259	GLN
1	B	278	GLN
1	B	313	ASN
1	B	418	GLN
1	B	420	GLN
1	B	462	GLN
1	B	505	GLN
1	B	574	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

Of 4 ligands modelled in this entry, 4 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.

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 ⓘ

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	381/428 (89%)	0.15	23 (6%) 21 23	42, 65, 97, 111	0
1	B	380/428 (88%)	0.14	17 (4%) 32 36	40, 61, 96, 125	0
All	All	761/856 (88%)	0.14	40 (5%) 26 28	40, 63, 97, 125	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	587	ALA	6.0
1	B	588	ALA	5.2
1	A	588	ALA	4.9
1	B	179	SER	4.8
1	B	181	ILE	4.5
1	B	496	ARG	3.9
1	A	220	ARG	3.6
1	A	587	ALA	3.6
1	B	218	LYS	3.5
1	A	179	SER	3.4
1	B	180	ILE	3.4
1	A	586	VAL	3.3
1	A	189	GLU	2.9
1	B	586	VAL	2.9
1	B	251	GLN	2.9
1	B	495	TYR	2.8
1	B	497	LYS	2.8
1	A	251	GLN	2.8
1	A	252	ASP	2.7
1	A	268	ALA	2.7
1	A	190	LYS	2.6
1	A	181	ILE	2.6
1	A	501	HIS	2.5
1	A	185	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	188	ARG	2.4
1	A	461	GLY	2.4
1	A	463	PRO	2.4
1	A	253	MET	2.4
1	A	221	ARG	2.3
1	B	214	GLN	2.2
1	B	185	LEU	2.2
1	B	583	LYS	2.2
1	A	186	LYS	2.1
1	A	219	GLU	2.1
1	B	492	LEU	2.1
1	A	462	GLN	2.0
1	B	493	GLU	2.0
1	A	254	PHE	2.0
1	A	214	GLN	2.0
1	B	188	ARG	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	MG	B	1	1/1	0.26	1.51	60,60,60,60	0
2	YT3	B	2	1/1	0.14	-0.50	65,65,65,65	0
2	YT3	A	1	1/1	0.10	-1.40	63,63,63,63	0
2	YT3	A	3	1/1	0.11	-4.19	71,71,71,71	0

6.5 Other polymers ⓘ

There are no such residues in this entry.