



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 10:09 AM UTC

PDB ID : 9DEP / pdb_00009dep
Title : USP7 in complex with macrocycle MC09
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Deposited on : 2024-08-29
Resolution : 2.57 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

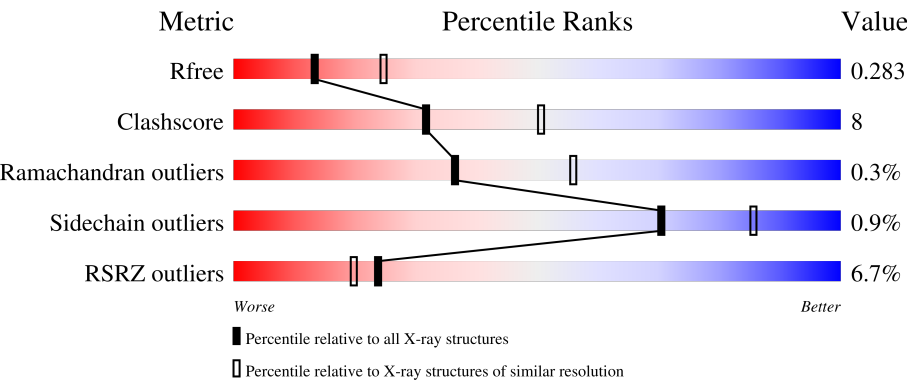
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.57 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	4770 (2.60-2.56)
Clashscore	190562	5124 (2.60-2.56)
Ramachandran outliers	187476	5046 (2.60-2.56)
Sidechain outliers	187428	5046 (2.60-2.56)
RSRZ outliers	180081	4770 (2.60-2.56)

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	368	2% 75% 15% • 9%
1	B	368	% 77% 15% 8%
1	E	368	14% 68% 19% 12%
2	C	15	7% 100%
2	D	15	87% 7% 7%

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Mol	Chain	Length	Quality of chain
2	F	15	20% 67% 27% 7%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8006 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 Ubiquitin carboxyl-terminal hydrolase 7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	334	Total	C	N	O	S	0	0	0
			2582	1634	429	503	16			
1	B	340	Total	C	N	O	S	0	0	0
			2675	1689	454	516	16			
1	E	323	Total	C	N	O	S	0	0	0
			2380	1507	390	468	15			

There are 63 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	187	MET	-	initiating methionine	UNP Q93009
A	188	GLY	-	expression tag	UNP Q93009
A	189	SER	-	expression tag	UNP Q93009
A	190	SER	-	expression tag	UNP Q93009
A	191	HIS	-	expression tag	UNP Q93009
A	192	HIS	-	expression tag	UNP Q93009
A	193	HIS	-	expression tag	UNP Q93009
A	194	HIS	-	expression tag	UNP Q93009
A	195	HIS	-	expression tag	UNP Q93009
A	196	HIS	-	expression tag	UNP Q93009
A	197	SER	-	expression tag	UNP Q93009
A	198	SER	-	expression tag	UNP Q93009
A	199	GLY	-	expression tag	UNP Q93009
A	200	LEU	-	expression tag	UNP Q93009
A	201	VAL	-	expression tag	UNP Q93009
A	202	PRO	-	expression tag	UNP Q93009
A	203	ARG	-	expression tag	UNP Q93009
A	204	GLY	-	expression tag	UNP Q93009
A	205	SER	-	expression tag	UNP Q93009
A	206	HIS	-	expression tag	UNP Q93009
A	207	MET	-	expression tag	UNP Q93009
B	187	MET	-	initiating methionine	UNP Q93009
B	188	GLY	-	expression tag	UNP Q93009

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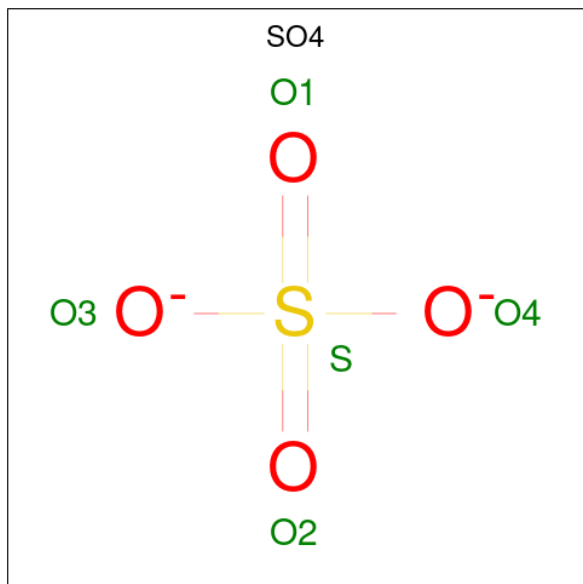
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Chain	Residue	Modelled	Actual	Comment	Reference
B	189	SER	-	expression tag	UNP Q93009
B	190	SER	-	expression tag	UNP Q93009
B	191	HIS	-	expression tag	UNP Q93009
B	192	HIS	-	expression tag	UNP Q93009
B	193	HIS	-	expression tag	UNP Q93009
B	194	HIS	-	expression tag	UNP Q93009
B	195	HIS	-	expression tag	UNP Q93009
B	196	HIS	-	expression tag	UNP Q93009
B	197	SER	-	expression tag	UNP Q93009
B	198	SER	-	expression tag	UNP Q93009
B	199	GLY	-	expression tag	UNP Q93009
B	200	LEU	-	expression tag	UNP Q93009
B	201	VAL	-	expression tag	UNP Q93009
B	202	PRO	-	expression tag	UNP Q93009
B	203	ARG	-	expression tag	UNP Q93009
B	204	GLY	-	expression tag	UNP Q93009
B	205	SER	-	expression tag	UNP Q93009
B	206	HIS	-	expression tag	UNP Q93009
B	207	MET	-	expression tag	UNP Q93009
E	187	MET	-	initiating methionine	UNP Q93009
E	188	GLY	-	expression tag	UNP Q93009
E	189	SER	-	expression tag	UNP Q93009
E	190	SER	-	expression tag	UNP Q93009
E	191	HIS	-	expression tag	UNP Q93009
E	192	HIS	-	expression tag	UNP Q93009
E	193	HIS	-	expression tag	UNP Q93009
E	194	HIS	-	expression tag	UNP Q93009
E	195	HIS	-	expression tag	UNP Q93009
E	196	HIS	-	expression tag	UNP Q93009
E	197	SER	-	expression tag	UNP Q93009
E	198	SER	-	expression tag	UNP Q93009
E	199	GLY	-	expression tag	UNP Q93009
E	200	LEU	-	expression tag	UNP Q93009
E	201	VAL	-	expression tag	UNP Q93009
E	202	PRO	-	expression tag	UNP Q93009
E	203	ARG	-	expression tag	UNP Q93009
E	204	GLY	-	expression tag	UNP Q93009
E	205	SER	-	expression tag	UNP Q93009
E	206	HIS	-	expression tag	UNP Q93009
E	207	MET	-	expression tag	UNP Q93009

- Molecule 2 is a protein called Macrocycle peptide MC09.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	15	Total	C	N	O	S	0	0	0
			121	81	21	18	1			
2	D	14	Total	C	N	O	S	0	0	0
			122	81	23	17	1			
2	F	14	Total	C	N	O	S	0	0	0
			104	66	20	17	1			

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

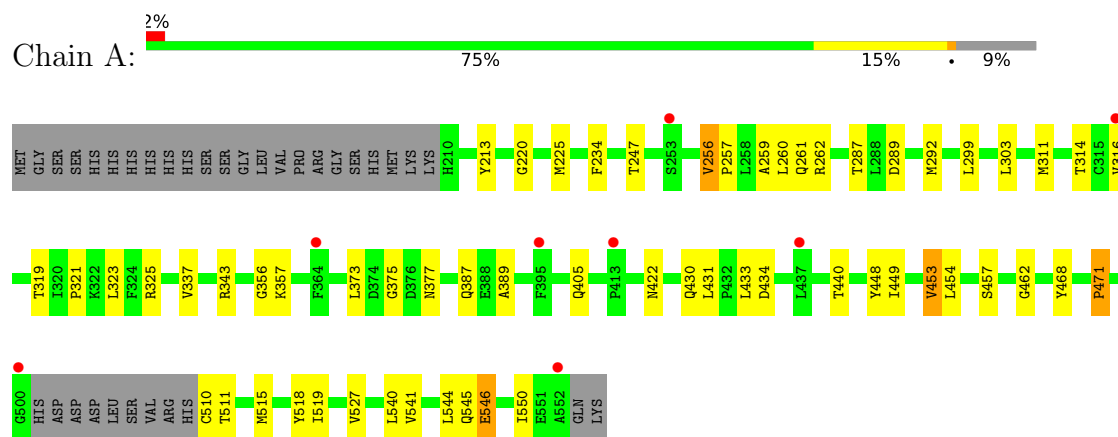
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	5	Total	O	0	0
			5	5		
4	B	8	Total	O	0	0
			8	8		
4	C	1	Total	O	0	0
			1	1		
4	E	3	Total	O	0	0
			3	3		

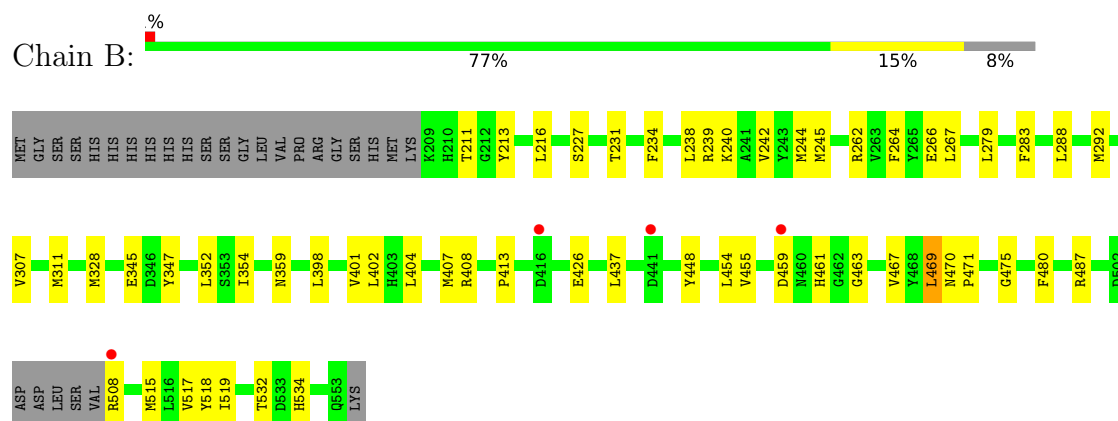
3 Residue-property plots [i](#)

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.

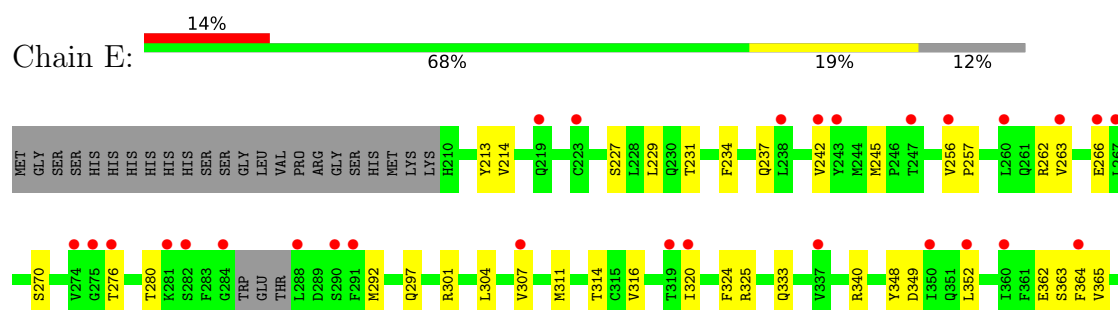
- Molecule 1: Ubiquitin carboxyl-terminal hydrolase 7

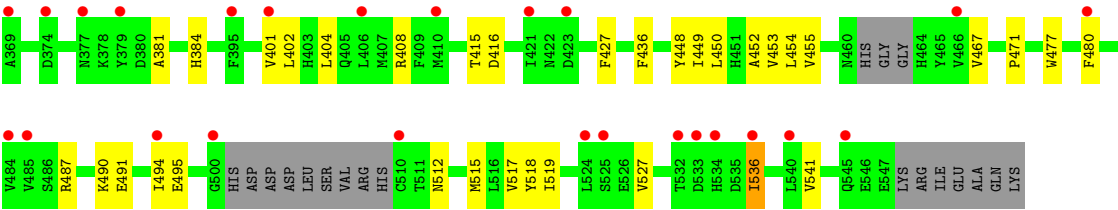


- Molecule 1: Ubiquitin carboxyl-terminal hydrolase 7



- Molecule 1: Ubiquitin carboxyl-terminal hydrolase 7

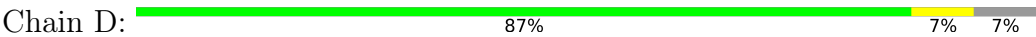




● Molecule 2: Macrocycle peptide MC09



● Molecule 2: Macrocycle peptide MC09



● Molecule 2: Macrocycle peptide MC09



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	245.00Å 68.85Å 76.05Å 90.00° 105.21° 90.00°	Depositor
Resolution (Å)	73.39 – 2.57 73.39 – 2.57	Depositor EDS
% Data completeness (in resolution range)	69.2 (73.39-2.57) 69.2 (73.39-2.57)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.40 (at 2.58Å)	Xtriage
Refinement program	PHENIX 1.21.2_5419	Depositor
R, R_{free}	0.247 , 0.284 0.249 , 0.283	Depositor DCC
R_{free} test set	1371 reflections (3.50%)	wwPDB-VP
Wilson B-factor (Å ²)	68.2	Xtriage
Anisotropy	0.038	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 55.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.005 for -h-2*1,-k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8006	wwPDB-VP
Average B, all atoms (Å ²)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 12.72% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ACE, 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	A	0.12	0/2638	0.32	0/3589
1	B	0.13	0/2733	0.29	0/3707
1	E	0.14	0/2429	0.32	0/3320
2	C	0.14	0/122	0.23	0/164
2	D	0.15	0/123	0.33	0/164
2	F	0.15	0/103	0.36	0/139
All	All	0.13	0/8148	0.31	0/11083

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	2582	0	2378	41	0
1	B	2675	0	2504	36	0
1	E	2380	0	2063	49	0
2	C	121	0	111	0	0
2	D	122	0	120	2	0
2	F	104	0	95	1	0
3	B	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	5	0	0	0	0
4	B	8	0	0	0	0
4	C	1	0	0	0	0
4	E	3	0	0	3	0
All	All	8006	0	7271	123	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 8.

All (123) 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:E:452:ALA:HB3	1:E:517:VAL:HB	1.71	0.71
1:E:450:LEU:HD22	1:E:477:TRP:HH2	1.57	0.69
1:B:413:PRO:HD3	2:D:4:ILE:HD12	1.78	0.66
1:A:457:SER:HB2	1:A:511:THR:HG23	1.76	0.66
1:A:337:VAL:HG22	1:A:387:GLN:HG2	1.79	0.64
1:E:450:LEU:HD21	1:E:453:VAL:HG23	1.79	0.64
1:E:494:ILE:HB	4:E:601:HOH:O	2.02	0.60
1:A:247:THR:HG21	1:A:261:GLN:OE1	2.02	0.59
1:E:467:VAL:HG13	1:E:480:PHE:HB2	1.84	0.59
1:A:259:ALA:HA	1:A:262:ARG:HH12	1.67	0.58
1:B:459:ASP:HB2	1:B:463:GLY:H	1.69	0.58
1:E:237:GLN:HB2	1:E:527:VAL:HA	1.84	0.58
1:A:448:TYR:HB3	1:A:518:TYR:HB3	1.85	0.57
1:E:454:LEU:HB2	1:E:515:MET:HB3	1.86	0.57
1:E:242:VAL:HA	1:E:245:MET:HE2	1.85	0.57
1:A:343:ARG:NH1	1:E:495:GLU:HG2	2.19	0.56
1:E:415:THR:HG22	1:E:416:ASP:H	1.71	0.56
1:A:356:GLY:C	1:A:357:LYS:HE2	2.31	0.56
1:E:477:TRP:N	1:E:487:ARG:HH21	2.04	0.55
1:E:229:LEU:HD23	1:E:263:VAL:HG11	1.87	0.55
1:A:292:MET:HE1	1:B:288:LEU:HD22	1.88	0.55
1:E:490:LYS:HD3	4:E:601:HOH:O	2.06	0.54
1:B:328:MET:HE3	1:B:345:GLU:HG3	1.90	0.54
1:E:262:ARG:O	1:E:266:GLU:HG2	2.08	0.54
1:E:491:GLU:HA	4:E:601:HOH:O	2.06	0.54
1:A:259:ALA:HA	1:A:262:ARG:NH1	2.24	0.53
1:E:448:TYR:HB3	1:E:518:TYR:HB3	1.91	0.53
1:E:519:ILE:HG21	1:E:527:VAL:HG11	1.91	0.53
1:B:454:LEU:HB2	1:B:515:MET:HB3	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:245:MET:SD	1:E:307:VAL:HG23	2.49	0.52
1:A:319:THR:O	1:A:323:LEU:HD12	2.10	0.52
1:B:234:PHE:HD2	1:B:469:LEU:HD13	1.75	0.52
1:A:422:ASN:ND2	1:A:510:CYS:HB3	2.25	0.51
1:B:402:LEU:HG	1:B:404:LEU:HD21	1.91	0.51
1:E:427:PHE:CE1	1:E:494:ILE:HG23	2.45	0.51
1:B:467:VAL:HG13	1:B:480:PHE:HB2	1.91	0.51
1:B:242:VAL:HG11	1:B:264:PHE:HD2	1.76	0.51
1:A:319:THR:HG22	1:A:323:LEU:HD11	1.94	0.50
1:E:314:THR:HG23	1:E:316:VAL:H	1.75	0.50
1:B:242:VAL:HG11	1:B:264:PHE:CD2	2.47	0.50
1:A:540:LEU:O	1:A:544:LEU:HG	2.12	0.50
1:A:287:THR:HG22	1:A:289:ASP:H	1.77	0.50
1:E:352:LEU:HB3	1:E:363:SER:OG	2.12	0.50
1:B:532:THR:HG23	1:B:534:HIS:H	1.76	0.50
1:A:321:PRO:O	1:A:325:ARG:HB3	2.13	0.49
1:E:297:GLN:O	1:E:301:ARG:HD3	2.12	0.49
1:E:408:ARG:HD3	1:E:512:ASN:ND2	2.27	0.49
1:E:364:PHE:HB3	1:E:436:PHE:CE2	2.47	0.49
1:E:311:MET:HE1	1:E:320:ILE:HD13	1.94	0.49
1:E:333:GLN:HB2	1:E:340:ARG:NH1	2.28	0.49
1:A:462:GLY:HA3	1:B:508:ARG:NH2	2.28	0.48
1:B:448:TYR:HB3	1:B:518:TYR:HB3	1.95	0.48
1:E:234:PHE:CD1	1:E:471:PRO:HB3	2.47	0.48
1:A:373:LEU:HD12	1:A:389:ALA:HB3	1.96	0.48
1:B:262:ARG:O	1:B:266:GLU:HG2	2.14	0.48
1:B:413:PRO:CD	2:D:4:ILE:HD12	2.44	0.48
1:E:402:LEU:HG	1:E:404:LEU:HD13	1.95	0.48
1:E:325:ARG:HB2	1:E:348:TYR:HE1	1.79	0.47
1:E:231:THR:HG21	1:E:517:VAL:HG21	1.96	0.47
1:E:304:LEU:HA	1:E:307:VAL:HG12	1.97	0.47
1:B:470:ASN:HD22	1:B:475:GLY:H	1.62	0.47
1:E:401:VAL:HG22	1:E:519:ILE:HG12	1.97	0.47
1:B:216:LEU:HD21	1:B:267:LEU:HD21	1.96	0.47
1:A:453:VAL:HG13	1:A:468:TYR:HB2	1.96	0.47
1:A:220:GLY:HA2	1:B:292:MET:HE1	1.96	0.47
1:B:359:ASN:HB3	1:B:426:GLU:HB2	1.96	0.47
1:A:314:THR:HG23	1:A:316:VAL:H	1.80	0.46
1:B:231:THR:HG21	1:B:517:VAL:HG21	1.97	0.46
1:B:242:VAL:HG23	1:B:245:MET:HE2	1.98	0.45
1:B:279:LEU:HG	1:B:283:PHE:HE2	1.80	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:546:GLU:O	1:A:550:ILE:HD12	2.16	0.45
1:E:292:MET:HE3	1:E:292:MET:HB2	1.80	0.45
1:B:211:THR:HG22	1:B:487:ARG:HG2	1.99	0.45
1:B:213:TYR:CZ	1:B:471:PRO:HG2	2.52	0.45
1:B:238:LEU:O	1:B:242:VAL:HG12	2.17	0.45
1:B:459:ASP:HB3	1:B:461:HIS:H	1.82	0.45
1:E:229:LEU:CD2	1:E:263:VAL:HG11	2.47	0.45
1:E:408:ARG:NH1	1:E:512:ASN:HD22	2.15	0.45
1:A:405:GLN:HG3	1:A:515:MET:HG3	1.99	0.45
1:E:227:SER:HB3	1:E:467:VAL:HB	1.98	0.45
1:A:213:TYR:CZ	1:A:471:PRO:HG2	2.52	0.44
1:A:260:LEU:HD22	1:A:303:LEU:HD23	1.99	0.44
1:E:213:TYR:CZ	1:E:471:PRO:HG2	2.52	0.44
1:A:541:VAL:O	1:A:545:GLN:HG3	2.16	0.44
1:A:434:ASP:OD2	1:A:440:THR:HB	2.17	0.44
1:B:307:VAL:O	1:B:311:MET:HG3	2.17	0.44
1:A:213:TYR:CE2	1:A:471:PRO:HG2	2.53	0.44
1:E:453:VAL:HG12	1:E:455:VAL:HG23	2.00	0.44
1:A:319:THR:HG22	1:A:323:LEU:CD1	2.47	0.43
1:A:454:LEU:HB2	1:A:515:MET:HB3	2.00	0.43
1:E:436:PHE:CD1	1:E:436:PHE:N	2.84	0.43
1:B:352:LEU:O	1:B:407:MET:HG2	2.18	0.43
1:A:256:VAL:HG22	1:A:257:PRO:HD3	1.99	0.43
1:A:430:GLN:HG3	1:A:449:ILE:HD13	2.01	0.43
1:B:347:TYR:OH	1:B:398:LEU:HD23	2.19	0.43
1:B:227:SER:HB3	1:B:467:VAL:HB	2.01	0.42
1:A:431:LEU:HG	1:A:433:LEU:HG	2.00	0.42
1:A:225:MET:HG3	1:A:299:LEU:HD21	2.00	0.42
1:B:354:ILE:HD13	1:B:408:ARG:HG2	2.01	0.42
2:F:7:PHE:CG	2:F:11:VAL:HB	2.54	0.42
1:E:536:ILE:HD12	1:E:541:VAL:HG22	2.01	0.42
1:B:292:MET:HE2	1:B:292:MET:HB3	1.81	0.42
1:E:324:PHE:CD1	1:E:401:VAL:HB	2.55	0.42
1:A:311:MET:HG2	1:A:316:VAL:HG23	2.01	0.42
1:B:240:LYS:O	1:B:244:MET:HG2	2.20	0.42
1:B:401:VAL:HG22	1:B:519:ILE:HG12	2.02	0.42
1:E:450:LEU:HD22	1:E:477:TRP:CH2	2.47	0.42
1:B:239:ARG:HA	1:B:242:VAL:HG12	2.01	0.42
1:E:449:ILE:HD13	1:E:449:ILE:HA	1.86	0.42
1:A:453:VAL:CG1	1:A:468:TYR:HB2	2.50	0.41
1:E:362:GLU:O	1:E:365:VAL:HG12	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:375:GLY:C	1:A:377:ASN:H	2.28	0.41
1:E:276:THR:O	1:E:280:THR:HG23	2.20	0.41
1:A:311:MET:O	1:A:314:THR:HG22	2.19	0.41
1:A:234:PHE:CE1	1:A:471:PRO:HB3	2.55	0.41
1:E:381:ALA:HB3	1:E:384:HIS:HB2	2.02	0.41
1:A:256:VAL:HA	1:A:259:ALA:HB3	2.02	0.41
1:A:292:MET:HE3	1:A:292:MET:HB2	1.94	0.41
1:E:256:VAL:HG13	1:E:257:PRO:HD3	2.02	0.41
1:B:437:LEU:HD23	1:B:437:LEU:HA	1.87	0.41
1:E:214:VAL:HG11	1:E:270:SER:O	2.20	0.41
1:A:519:ILE:HD13	1:A:527:VAL:HG11	2.02	0.40
1:E:415:THR:HG22	1:E:416:ASP:N	2.36	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	330/368 (90%)	310 (94%)	18 (6%)	2 (1%)	21	39
1	B	336/368 (91%)	324 (96%)	12 (4%)	0	100	100
1	E	315/368 (86%)	298 (95%)	17 (5%)	0	100	100
2	C	13/15 (87%)	12 (92%)	1 (8%)	0	100	100
2	D	12/15 (80%)	11 (92%)	1 (8%)	0	100	100
2	F	12/15 (80%)	10 (83%)	1 (8%)	1 (8%)	0	0
All	All	1018/1149 (89%)	965 (95%)	50 (5%)	3 (0%)	36	56

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	1	PHE

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Mol	Chain	Res	Type
1	A	546	GLU
1	A	471	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	270/332 (81%)	268 (99%)	2 (1%)	76	88
1	B	284/332 (86%)	282 (99%)	2 (1%)	76	88
1	E	229/332 (69%)	227 (99%)	2 (1%)	70	85
2	C	11/13 (85%)	11 (100%)	0	100	100
2	D	12/13 (92%)	12 (100%)	0	100	100
2	F	9/13 (69%)	8 (89%)	1 (11%)	6	11
All	All	815/1035 (79%)	808 (99%)	7 (1%)	70	85

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	256	VAL
1	A	453	VAL
1	B	455	VAL
1	B	469	LEU
1	E	349	ASP
1	E	536	ILE
2	F	13	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	306	ASN
1	A	377	ASN
1	A	403	HIS
1	A	422	ASN

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Mol	Chain	Res	Type
1	B	309	ASN
1	B	414	GLN
1	E	237	GLN
1	E	384	HIS
1	E	512	ASN
2	F	10	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is 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	$\# Z > 2$
3	SO4	B	601	-	4,4,4	0.67	0	6,6,6	0.04	0

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

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	334/368 (90%)	0.44	8 (2%) 59 55	41, 68, 101, 138	0
1	B	340/368 (92%)	0.24	4 (1%) 76 74	35, 58, 87, 142	0
1	E	323/368 (87%)	1.18	53 (16%) 4 3	67, 92, 137, 168	0
2	C	14/15 (93%)	0.88	1 (7%) 22 18	50, 64, 86, 112	0
2	D	13/15 (86%)	0.41	0 100 100	47, 59, 81, 89	0
2	F	13/15 (86%)	1.46	3 (23%) 2 2	90, 104, 133, 147	0
All	All	1037/1149 (90%)	0.62	69 (6%) 24 20	35, 73, 119, 168	0

All (69) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	364	PHE	4.6
1	E	219	GLN	4.2
1	A	395	PHE	4.2
1	E	256	VAL	4.1
1	E	288	LEU	4.1
1	E	290	SER	3.9
1	E	260	LEU	3.6
1	E	282	SER	3.4
2	C	2	PHE	3.4
1	E	377	ASN	3.4
1	E	350	ILE	3.4
1	E	540	LEU	3.3
1	E	480	PHE	3.2
1	E	243	TYR	3.1
1	E	275	GLY	3.0
1	B	508	ARG	3.0
1	E	500	GLY	3.0
1	E	494	ILE	2.9
1	B	416	ASP	2.9

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Mol	Chain	Res	Type	RSRZ
2	F	11	VAL	2.8
1	E	379	TYR	2.7
1	E	421	ILE	2.7
1	E	406	LEU	2.7
1	E	395	PHE	2.6
1	E	410	MET	2.6
1	E	352	LEU	2.6
1	E	484	VAL	2.6
1	E	284	GLY	2.5
1	E	319	THR	2.5
2	F	3	ARG	2.5
1	E	307	VAL	2.4
1	E	401	VAL	2.4
1	E	369	ALA	2.4
1	E	485	VAL	2.4
1	E	281	LYS	2.4
1	E	247	THR	2.3
1	E	267	LEU	2.3
1	B	459	ASP	2.3
1	A	316	VAL	2.3
1	E	266	GLU	2.3
1	E	274	VAL	2.3
1	E	466	VAL	2.3
1	A	500	GLY	2.3
1	E	510	CYS	2.3
1	E	532	THR	2.2
1	A	552	ALA	2.2
1	E	534	HIS	2.2
1	E	545	GLN	2.2
1	E	276	THR	2.2
1	A	437	LEU	2.2
1	A	364	PHE	2.2
1	E	374	ASP	2.2
1	E	291	PHE	2.1
1	E	536	ILE	2.1
2	F	7	PHE	2.1
1	E	337	VAL	2.1
1	E	524	LEU	2.1
1	E	533	ASP	2.1
1	A	413	PRO	2.1
1	E	242	VAL	2.1
1	E	320	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	E	360	ILE	2.1
1	E	525	SER	2.1
1	E	238	LEU	2.0
1	E	223	CYS	2.0
1	E	263	VAL	2.0
1	B	441	ASP	2.0
1	E	423	ASP	2.0
1	A	253	SER	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
3	SO4	B	601	5/5	0.63	0.13	101,109,113,118	0

6.5 Other polymers [i](#)

There are no such residues in this entry.