



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:58 AM UTC

PDB ID : 7SA1 / pdb_00007sa1
Title : LRR-F-Box plant ubiquitin ligase
Authors : Palayam, M.; Shabek, N.
Deposited on : 2021-09-21
Resolution : 3.21 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

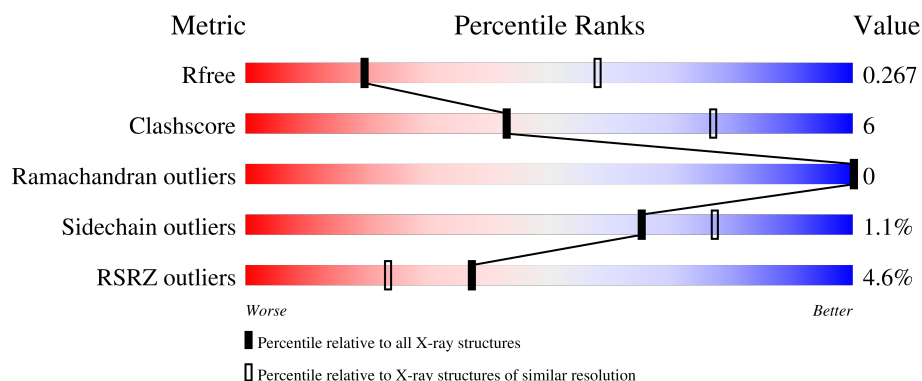
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.21 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	1768 (3.24-3.20)
Clashscore	190562	1879 (3.24-3.20)
Ramachandran outliers	187476	1844 (3.24-3.20)
Sidechain outliers	187428	1843 (3.24-3.20)
RSRZ outliers	180081	1768 (3.24-3.20)

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	160	
1	C	160	
2	B	688	
2	D	688	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11728 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 SKP1-like protein 1A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	122	Total	C	N	O	S	0	0	0
			978	621	159	194	4			
1	C	124	Total	C	N	O	S	0	0	0
			992	629	161	198	4			

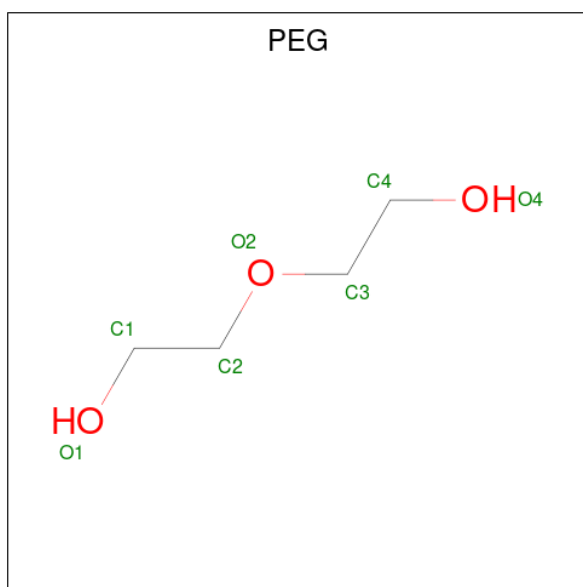
- Molecule 2 is a protein called F-box/LRR-repeat MAX2 homolog.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	624	Total	C	N	O	S	0	0	0
			4843	3081	849	883	30			
2	D	624	Total	C	N	O	S	0	0	0
			4843	3081	849	883	30			

There are 2 discrepancies between the modelled and reference sequences:

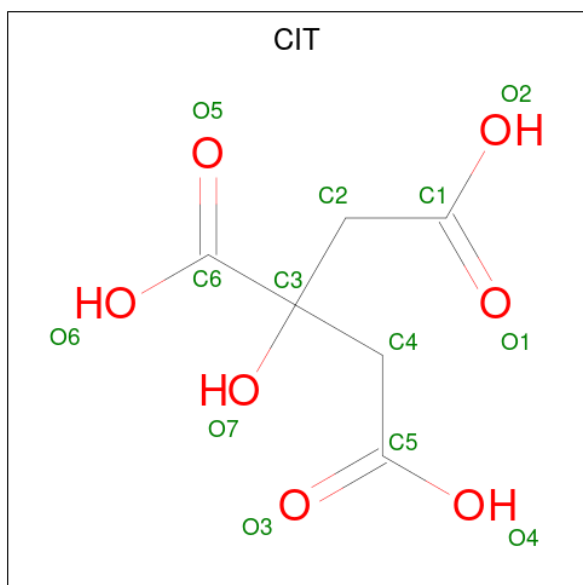
Chain	Residue	Modelled	Actual	Comment	Reference
B	720	LYS	ASP	engineered mutation	UNP Q5VMP0
D	720	LYS	ASP	engineered mutation	UNP Q5VMP0

- Molecule 3 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			7	4	3		
3	B	1	Total	C	O	0	0
			7	4	3		
3	B	1	Total	C	O	0	0
			7	4	3		
3	C	1	Total	C	O	0	0
			7	4	3		

- Molecule 4 is CITRIC ACID (CCD ID: CIT) (formula: $C_6H_8O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			13	6	7		
4	D	1	Total	C	O	0	0
			13	6	7		

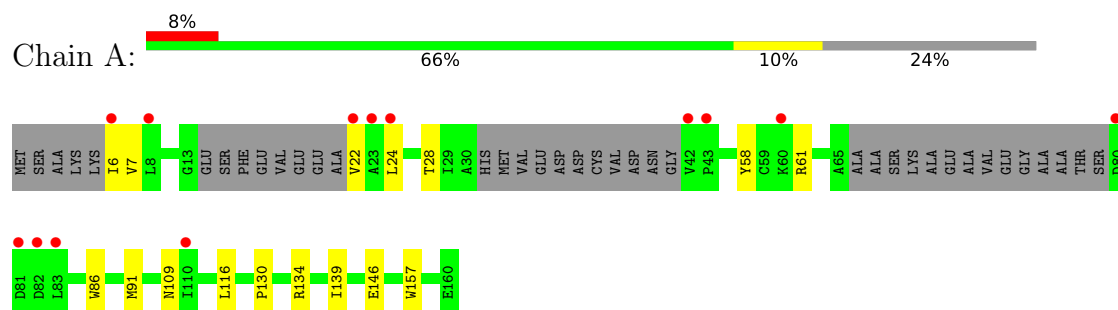
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	11	Total	O	0	0
			11	11		
5	C	1	Total	O	0	0
			1	1		
5	D	6	Total	O	0	0
			6	6		

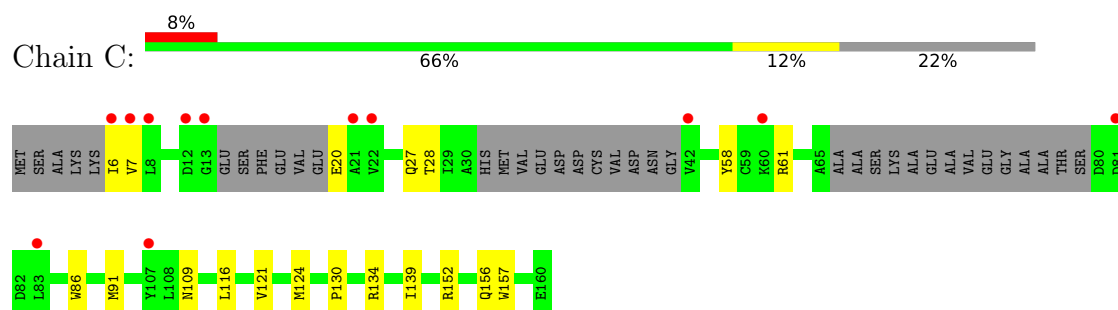
3 Residue-property plots [i](#)

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.

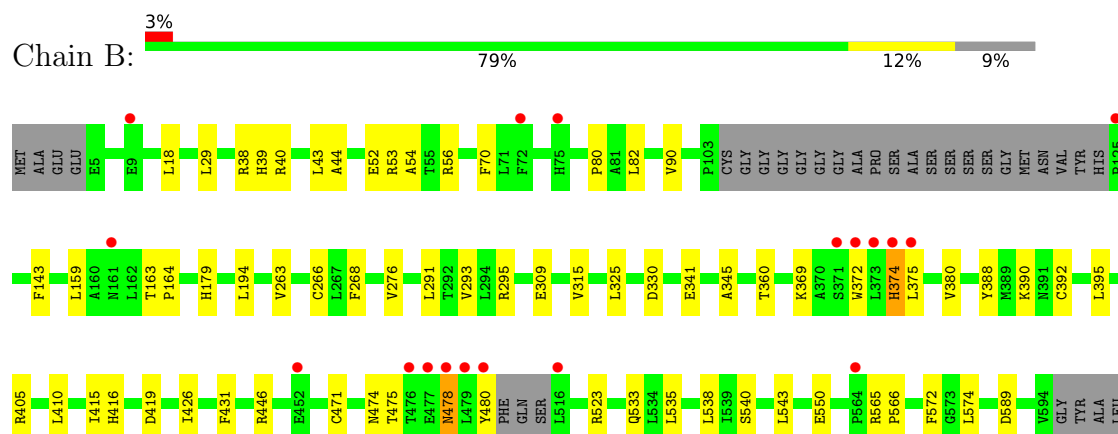
• Molecule 1: SKP1-like protein 1A

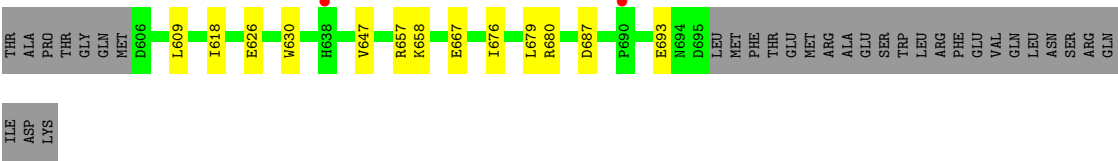


• Molecule 1: SKP1-like protein 1A

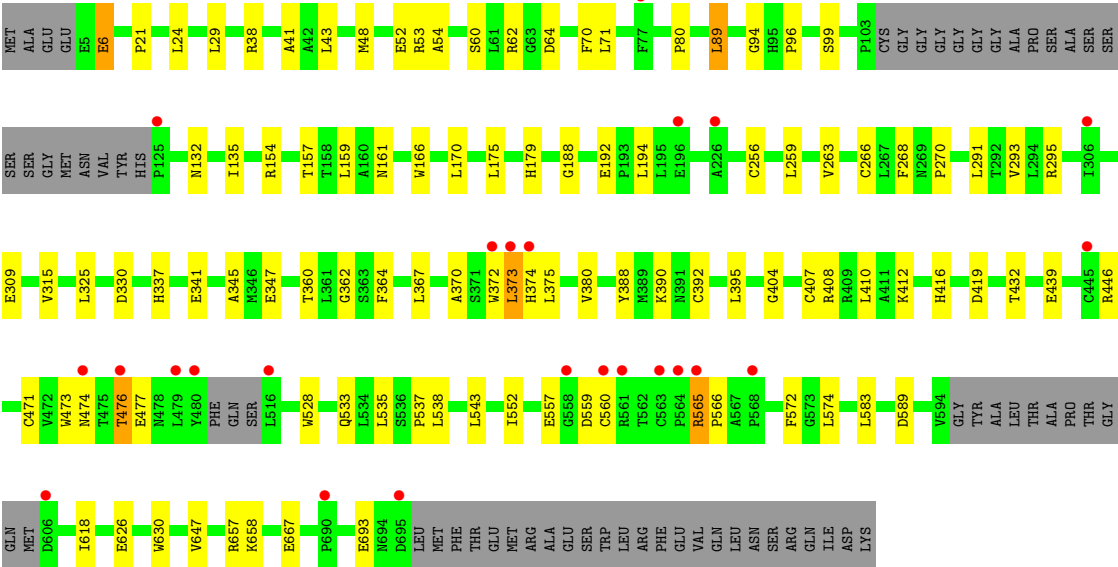
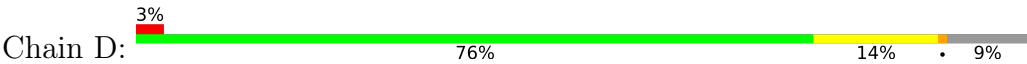


• Molecule 2: F-box/LRR-repeat MAX2 homolog





● Molecule 2: F-box/LRR-repeat MAX2 homolog



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	231.93Å 79.41Å 150.59Å 90.00° 127.98° 90.00°	Depositor
Resolution (Å)	48.90 – 3.21 48.90 – 3.21	Depositor EDS
% Data completeness (in resolution range)	97.6 (48.90-3.21) 87.6 (48.90-3.21)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.06 (at 3.19Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.208 , 0.264 0.210 , 0.267	Depositor DCC
R_{free} test set	1986 reflections (5.73%)	wwPDB-VP
Wilson B-factor (Å ²)	33.8	Xtriage
Anisotropy	0.812	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 39.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.136 for -h-2*k,l	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	11728	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 3.68% 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: CIT, PEG

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	0.28	0/990	0.36	0/1337
1	C	0.07	0/1004	0.22	0/1356
2	B	0.27	0/4956	0.44	3/6745 (0.0%)
2	D	0.28	1/4956 (0.0%)	0.44	1/6745 (0.0%)
All	All	0.27	1/11906 (0.0%)	0.42	4/16183 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	96	PRO	C-O	-5.12	1.17	1.23

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	370	ALA	N-CA-C	8.50	120.86	110.91
2	B	478	ASN	CB-CA-C	-7.82	104.62	114.40
2	B	372	TRP	N-CA-C	-6.24	103.77	112.25
2	B	478	ASN	N-CA-C	-5.11	102.08	109.29

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	978	0	974	13	0
1	C	992	0	985	20	0
2	B	4843	0	4853	53	0
2	D	4843	0	4853	64	0
3	A	7	0	10	0	0
3	B	14	0	20	0	0
3	C	7	0	10	0	0
4	B	13	0	5	3	0
4	D	13	0	5	5	0
5	B	11	0	0	0	0
5	C	1	0	0	0	0
5	D	6	0	0	0	0
All	All	11728	0	11715	144	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 6.

All (144) 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)
2:B:375:LEU:O	2:B:380:VAL:HG23	1.75	0.87
2:B:179:HIS:HD2	2:B:693:GLU:HG3	1.50	0.77
2:D:657:ARG:HH12	4:D:801:CIT:H22	1.49	0.76
2:D:372:TRP:H	2:D:372:TRP:HE3	1.34	0.73
2:B:375:LEU:O	2:B:380:VAL:CG2	2.36	0.72
2:B:419:ASP:HB3	2:B:446:ARG:HE	1.52	0.72
1:C:28:THR:HG23	1:C:109:ASN:HB2	1.74	0.70
1:C:6:ILE:HG13	1:C:7:VAL:N	2.06	0.69
2:D:179:HIS:HD2	2:D:693:GLU:HG3	1.59	0.67
2:B:565:ARG:HH21	2:B:609:LEU:HB3	1.62	0.63
1:C:6:ILE:HG13	1:C:7:VAL:H	1.65	0.61
1:C:6:ILE:HG23	1:C:20:GLU:HA	1.81	0.61
2:B:478:ASN:HB3	2:B:540:SER:HB2	1.81	0.61
2:D:419:ASP:HB3	2:D:446:ARG:HE	1.66	0.61
2:D:404:GLY:O	2:D:408:ARG:NH1	2.34	0.60
2:B:375:LEU:HD21	2:B:405:ARG:HB2	1.84	0.59
2:D:565:ARG:CZ	2:D:566:PRO:HD2	2.33	0.58
2:D:157:THR:O	2:D:161:ASN:ND2	2.35	0.58
2:B:266:CYS:HB3	2:B:276:VAL:HG21	1.87	0.57
2:B:538:LEU:HD11	2:B:543:LEU:HD12	1.87	0.56
2:B:475:THR:HG22	2:B:475:THR:O	2.05	0.56
1:A:130:PRO:HD3	2:B:40:ARG:HG2	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:21:PRO:HD2	2:D:24:LEU:HD12	1.88	0.55
2:D:159:LEU:HD13	2:D:194:LEU:HG	1.90	0.54
2:B:431:PHE:HE2	2:D:6:GLU:O	1.91	0.54
2:D:626:GLU:HG3	2:D:658:LYS:HB2	1.89	0.54
2:B:263:VAL:HG22	2:B:295:ARG:HB2	1.89	0.54
2:B:159:LEU:HD13	2:B:194:LEU:HG	1.89	0.54
2:D:263:VAL:HG22	2:D:295:ARG:HB2	1.89	0.53
2:D:538:LEU:HD11	2:D:543:LEU:HD12	1.91	0.53
2:D:471:CYS:O	2:D:533:GLN:NE2	2.34	0.53
2:D:291:LEU:HB2	2:D:325:LEU:HD22	1.90	0.53
1:A:134:ARG:HG3	1:A:139:ILE:HB	1.91	0.53
2:B:471:CYS:O	2:B:533:GLN:NE2	2.29	0.53
2:D:159:LEU:HD21	2:D:175:LEU:HD21	1.89	0.52
2:D:528:TRP:HE3	2:D:557:GLU:OE1	1.92	0.52
2:D:574:LEU:HD23	2:D:618:ILE:HB	1.91	0.52
2:B:309:GLU:OE1	2:B:309:GLU:N	2.42	0.52
1:C:91:MET:HE3	1:C:116:LEU:HG	1.93	0.51
4:D:801:CIT:O2	4:D:801:CIT:O7	2.27	0.51
2:D:589:ASP:HA	2:D:630:TRP:HB2	1.91	0.51
1:A:139:ILE:HG23	2:B:18:LEU:HD11	1.93	0.51
2:D:94:GLY:O	2:D:154:ARG:NH1	2.40	0.51
2:D:657:ARG:NH1	4:D:801:CIT:H22	2.23	0.50
2:B:29:LEU:O	2:B:38:ARG:NH1	2.44	0.50
2:D:54:ALA:HA	2:D:80:PRO:HD2	1.93	0.50
2:B:52:GLU:OE1	2:B:53:ARG:NH1	2.45	0.50
2:B:54:ALA:HA	2:B:80:PRO:HD2	1.94	0.50
2:D:407:CYS:HB3	2:D:410:LEU:HG	1.93	0.50
1:A:91:MET:HE3	1:A:116:LEU:HG	1.93	0.49
2:D:364:PHE:HB3	2:D:367:LEU:HD21	1.93	0.49
2:B:163:THR:N	2:B:164:PRO:HD2	2.27	0.49
2:D:647:VAL:HG11	2:D:667:GLU:HG3	1.94	0.49
1:C:6:ILE:CG1	1:C:7:VAL:H	2.21	0.49
2:B:574:LEU:HD23	2:B:618:ILE:HB	1.94	0.49
2:B:374:HIS:O	2:B:375:LEU:HG	2.13	0.48
2:D:60:SER:OG	2:D:89:LEU:CD1	2.61	0.48
1:C:6:ILE:CG2	1:C:20:GLU:HB2	2.44	0.48
2:D:657:ARG:HH22	4:D:801:CIT:C2	2.27	0.48
1:C:121:VAL:HA	1:C:124:MET:HE3	1.95	0.48
1:A:58:TYR:HA	1:A:86:TRP:HH2	1.78	0.48
2:B:676:ILE:HD12	2:B:679:LEU:HD22	1.95	0.48
2:B:266:CYS:HB2	2:B:268:PHE:CE2	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:392:CYS:SG	2:D:395:LEU:HB2	2.54	0.47
1:A:7:VAL:HG12	1:A:7:VAL:O	2.14	0.47
2:B:647:VAL:HG11	2:B:667:GLU:HG3	1.95	0.47
2:B:315:VAL:HG11	2:B:341:GLU:HG3	1.96	0.47
2:D:360:THR:HG23	2:D:388:TYR:HB3	1.95	0.47
2:B:565:ARG:NH1	2:B:566:PRO:O	2.47	0.47
1:C:6:ILE:HG23	1:C:20:GLU:CB	2.45	0.47
2:B:535:LEU:HB2	2:B:572:PHE:CZ	2.51	0.46
2:B:626:GLU:HG3	2:B:658:LYS:HB2	1.97	0.46
2:D:52:GLU:OE1	2:D:53:ARG:NH1	2.48	0.46
1:A:130:PRO:O	1:A:134:ARG:HD3	2.15	0.46
2:D:476:THR:HG23	2:D:477:GLU:HG3	1.98	0.46
1:A:61:ARG:HG2	1:A:86:TRP:CE3	2.51	0.46
2:D:99:SER:O	2:D:132:ASN:ND2	2.42	0.46
2:D:657:ARG:HH22	4:D:801:CIT:H22	1.80	0.46
2:B:392:CYS:SG	2:B:395:LEU:HB2	2.56	0.46
1:A:6:ILE:HG22	1:A:6:ILE:O	2.16	0.46
2:B:82:LEU:HD22	2:B:143:PHE:HB3	1.97	0.46
1:C:6:ILE:HG23	1:C:20:GLU:CA	2.45	0.45
1:C:156:GLN:HG3	2:D:62:ARG:HH12	1.81	0.45
2:B:657:ARG:NH2	2:B:680:ARG:HD2	2.30	0.45
2:B:40:ARG:HA	2:B:43:LEU:HD12	1.98	0.45
1:C:130:PRO:O	1:C:134:ARG:HD3	2.16	0.45
2:B:589:ASP:HA	2:B:630:TRP:HB2	1.97	0.45
1:C:157:TRP:CH2	2:D:70:PHE:HB2	2.51	0.45
2:B:291:LEU:HB2	2:B:325:LEU:HD22	1.98	0.45
2:D:374:HIS:CD2	2:D:375:LEU:HG	2.52	0.45
2:D:412:LYS:HG2	2:D:439:GLU:HB2	1.98	0.45
1:C:61:ARG:HG2	1:C:86:TRP:CE3	2.52	0.44
2:D:315:VAL:HG11	2:D:341:GLU:HG3	1.99	0.44
1:A:146:GLU:N	1:A:146:GLU:OE1	2.50	0.44
2:B:374:HIS:HB3	2:B:375:LEU:H	1.55	0.44
1:A:157:TRP:CH2	2:B:70:PHE:HB2	2.52	0.44
2:B:415:ILE:HD11	2:B:426:ILE:HD11	2.00	0.44
2:D:362:GLY:HA2	2:D:390:LYS:HB2	1.98	0.44
2:B:390:LYS:HG2	2:B:416:HIS:HB3	2.00	0.44
2:D:60:SER:OG	2:D:89:LEU:HD12	2.18	0.44
2:D:71:LEU:HD21	2:D:135:ILE:HG12	2.00	0.44
2:B:315:VAL:HG23	2:B:345:ALA:HB2	2.00	0.44
2:B:410:LEU:HD23	2:B:410:LEU:HA	1.83	0.44
2:D:408:ARG:NH2	2:D:432:THR:O	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:803:CIT:O1	4:B:803:CIT:O7	2.33	0.43
1:C:58:TYR:HA	1:C:86:TRP:HH2	1.83	0.43
2:B:374:HIS:C	2:B:375:LEU:HG	2.42	0.43
2:B:657:ARG:HH22	4:B:803:CIT:H42	1.81	0.43
2:D:270:PRO:HD2	2:D:337:HIS:CE1	2.54	0.43
1:C:152:ARG:HG2	2:D:43:LEU:HD21	1.99	0.43
2:D:170:LEU:HD12	2:D:170:LEU:HA	1.86	0.43
2:B:39:HIS:CD2	2:B:56:ARG:HH22	2.37	0.43
2:B:687:ASP:OD1	2:B:687:ASP:N	2.47	0.43
2:B:360:THR:HG23	2:B:388:TYR:HB3	2.00	0.43
4:B:803:CIT:O7	4:B:803:CIT:O4	2.35	0.43
2:D:373:LEU:HB3	2:D:374:HIS:H	1.63	0.43
2:B:293:VAL:HG22	2:B:330:ASP:HB3	2.00	0.43
1:A:28:THR:HG23	1:A:109:ASN:HB2	2.01	0.43
2:D:29:LEU:HD21	2:D:48:MET:HE3	2.01	0.43
2:D:41:ALA:HB1	2:D:48:MET:HE2	2.01	0.43
2:D:552:ILE:HG22	2:D:583:LEU:HD11	2.01	0.43
1:C:157:TRP:HZ3	2:D:64:ASP:H	1.67	0.42
2:D:256:CYS:HB2	2:D:259:LEU:HG	2.01	0.42
2:D:347:GLU:HG2	2:D:380:VAL:HG21	2.02	0.42
2:D:29:LEU:O	2:D:38:ARG:NH1	2.51	0.42
2:D:315:VAL:HG23	2:D:345:ALA:HB2	2.02	0.42
2:D:410:LEU:HD23	2:D:410:LEU:HA	1.83	0.42
1:C:134:ARG:HA	1:C:139:ILE:HB	2.01	0.42
2:D:535:LEU:HB2	2:D:572:PHE:CZ	2.54	0.42
2:B:478:ASN:HD22	2:B:480:TYR:HD2	1.68	0.42
1:C:61:ARG:HG2	1:C:86:TRP:CD2	2.55	0.42
2:D:309:GLU:OE1	2:D:309:GLU:N	2.49	0.41
2:D:528:TRP:CE3	2:D:557:GLU:OE1	2.72	0.41
1:C:27:GLN:HB2	1:C:109:ASN:HB3	2.02	0.41
2:D:293:VAL:HG22	2:D:330:ASP:HB3	2.03	0.41
2:B:179:HIS:CD2	2:B:693:GLU:HG3	2.41	0.41
2:D:473:TRP:CD2	2:D:537:PRO:HB2	2.56	0.41
1:A:134:ARG:HD2	2:B:44:ALA:HA	2.03	0.41
2:B:523:ARG:NH1	2:B:550:GLU:OE1	2.54	0.41
2:B:375:LEU:O	2:B:380:VAL:HG22	2.19	0.40
2:D:188:GLY:O	2:D:192:GLU:HG3	2.22	0.40
2:D:390:LYS:HG2	2:D:416:HIS:HB3	2.02	0.40
2:D:266:CYS:HB2	2:D:268:PHE:CE2	2.57	0.40
2:D:166:TRP:O	2:D:170:LEU:HB2	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	114/160 (71%)	108 (95%)	6 (5%)	0	100	100
1	C	116/160 (72%)	113 (97%)	3 (3%)	0	100	100
2	B	616/688 (90%)	576 (94%)	40 (6%)	0	100	100
2	D	616/688 (90%)	582 (94%)	34 (6%)	0	100	100
All	All	1462/1696 (86%)	1379 (94%)	83 (6%)	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	A	109/137 (80%)	107 (98%)	2 (2%)	51	72
1	C	110/137 (80%)	110 (100%)	0	100	100
2	B	526/575 (92%)	522 (99%)	4 (1%)	73	80
2	D	526/575 (92%)	518 (98%)	8 (2%)	57	74
All	All	1271/1424 (89%)	1257 (99%)	14 (1%)	65	77

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	VAL
1	A	24	LEU

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Mol	Chain	Res	Type
2	B	90	VAL
2	B	369	LYS
2	B	374	HIS
2	B	474	ASN
2	D	6	GLU
2	D	89	LEU
2	D	373	LEU
2	D	474	ASN
2	D	476	THR
2	D	559	ASP
2	D	560	CYS
2	D	565	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	62	HIS
1	A	119	GLN
2	B	131	GLN
2	B	165	HIS
2	B	305	ASN
2	D	165	HIS
2	D	337	HIS
2	D	374	HIS

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

6 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
4	CIT	B	803	-	12,12,12	1.10	0	17,17,17	1.37	1 (5%)
4	CIT	D	801	-	12,12,12	1.15	1 (8%)	17,17,17	1.12	2 (11%)
3	PEG	A	801	-	6,6,6	0.51	0	5,5,5	0.29	0
3	PEG	B	801	-	6,6,6	0.50	0	5,5,5	0.28	0
3	PEG	B	802	-	6,6,6	0.49	0	5,5,5	0.33	0
3	PEG	C	201	-	6,6,6	0.50	0	5,5,5	0.26	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
4	CIT	B	803	-	-	8/16/16/16	-
4	CIT	D	801	-	-	9/16/16/16	-
3	PEG	A	801	-	-	2/4/4/4	-
3	PEG	B	801	-	-	3/4/4/4	-
3	PEG	B	802	-	-	2/4/4/4	-
3	PEG	C	201	-	-	1/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	801	CIT	C3-C6	2.06	1.55	1.53

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	803	CIT	O6-C6-C3	3.60	120.04	113.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	801	CIT	O5-C6-C3	-2.47	117.31	122.09
4	D	801	CIT	O6-C6-C3	2.19	117.34	113.14

There are no chirality outliers.

All (25) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	803	CIT	O7-C3-C6-O5
4	B	803	CIT	O7-C3-C6-O6
4	B	803	CIT	C4-C3-C6-O5
4	B	803	CIT	C4-C3-C6-O6
4	D	801	CIT	O7-C3-C6-O5
4	D	801	CIT	O7-C3-C6-O6
4	D	801	CIT	C4-C3-C6-O5
4	D	801	CIT	C4-C3-C6-O6
4	D	801	CIT	C1-C2-C3-O7
4	D	801	CIT	C1-C2-C3-C6
3	A	801	PEG	O1-C1-C2-O2
4	D	801	CIT	C1-C2-C3-C4
3	B	801	PEG	O2-C3-C4-O4
3	A	801	PEG	C1-C2-O2-C3
3	B	802	PEG	O1-C1-C2-O2
3	B	801	PEG	C1-C2-O2-C3
3	B	802	PEG	C4-C3-O2-C2
4	B	803	CIT	C3-C4-C5-O4
4	D	801	CIT	O2-C1-C2-C3
3	B	801	PEG	C4-C3-O2-C2
4	D	801	CIT	O1-C1-C2-C3
4	B	803	CIT	C3-C4-C5-O3
4	B	803	CIT	C1-C2-C3-O7
4	B	803	CIT	C1-C2-C3-C6
3	C	201	PEG	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	803	CIT	3	0
4	D	801	CIT	5	0

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	122/160 (76%)	0.78	13 (10%) 11 7	21, 69, 111, 145	0
1	C	124/160 (77%)	0.70	12 (9%) 13 9	25, 72, 113, 152	0
2	B	624/688 (90%)	0.25	20 (3%) 50 32	12, 32, 83, 137	0
2	D	624/688 (90%)	0.33	24 (3%) 44 28	14, 36, 89, 137	0
All	All	1494/1696 (88%)	0.36	69 (4%) 37 23	12, 36, 101, 152	0

All (69) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	373	LEU	6.7
2	B	371	SER	5.3
1	A	22	VAL	4.9
2	B	374	HIS	4.6
2	B	125	PRO	4.5
1	A	23	ALA	4.2
2	D	373	LEU	4.2
2	B	372	TRP	4.1
2	D	125	PRO	4.1
1	C	21	ALA	3.8
2	D	374	HIS	3.6
2	D	558	GLY	3.5
1	A	81	ASP	3.3
2	B	9	GLU	3.2
1	A	43	PRO	3.1
2	D	480	TYR	3.1
1	A	24	LEU	3.0
2	B	72	PHE	3.0
2	D	563	CYS	3.0
1	C	7	VAL	3.0
1	A	83	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
2	B	479	LEU	2.9
1	C	8	LEU	2.9
1	C	6	ILE	2.8
2	D	564	PRO	2.8
2	B	452	GLU	2.7
1	C	22	VAL	2.7
2	D	695	ASP	2.7
2	D	372	TRP	2.6
2	B	564	PRO	2.6
2	D	474	ASN	2.6
2	B	75	HIS	2.6
1	C	13	GLY	2.6
1	A	8	LEU	2.6
1	C	107	TYR	2.6
2	B	476	THR	2.5
2	D	306	ILE	2.5
1	C	81	ASP	2.4
2	B	477	GLU	2.4
2	B	638	HIS	2.3
2	D	565	ARG	2.3
2	D	516	LEU	2.3
1	C	60	LYS	2.3
2	B	516	LEU	2.3
2	B	375	LEU	2.3
2	D	445	CYS	2.3
1	A	80	ASP	2.3
1	A	42	VAL	2.2
2	B	161	ASN	2.2
1	C	83	LEU	2.2
2	D	77	PHE	2.2
2	D	476	THR	2.2
2	D	560	CYS	2.2
1	C	42	VAL	2.2
1	A	6	ILE	2.2
2	D	479	LEU	2.2
2	B	478	ASN	2.2
2	D	196	GLU	2.2
2	D	568	PRO	2.2
2	D	690	PRO	2.1
1	A	82	ASP	2.1
2	D	606	ASP	2.1
1	A	60	LYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	110	ILE	2.1
1	C	12	ASP	2.1
2	D	561	ARG	2.1
2	D	226	ALA	2.1
2	B	690	PRO	2.1
2	B	480	TYR	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(Å ²)	Q<0.9
4	CIT	D	801	13/13	0.54	0.21	46,69,101,107	0
3	PEG	B	802	7/7	0.66	0.19	39,44,58,64	0
4	CIT	B	803	13/13	0.67	0.16	47,65,77,80	0
3	PEG	A	801	7/7	0.84	0.14	24,29,39,43	0
3	PEG	C	201	7/7	0.87	0.12	18,21,35,44	0
3	PEG	B	801	7/7	0.88	0.14	29,44,62,68	0

6.5 Other polymers [i](#)

There are no such residues in this entry.