



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:53 AM UTC

PDB ID : 1CN3 / pdb_00001cn3
Title : INTERACTION OF POLYOMAVIRUS INTERNAL PROTEIN VP2 WITH
MAJOR CAPSID PROTEIN VP1 AND IMPLICATIONS FOR PARTICIPA-
TION OF VP2 IN VIRAL ENTRY
Authors : Chen, X.; Stehle, T.; Harrison, S.C.
Deposited on : 1999-05-24
Resolution : 2.20 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

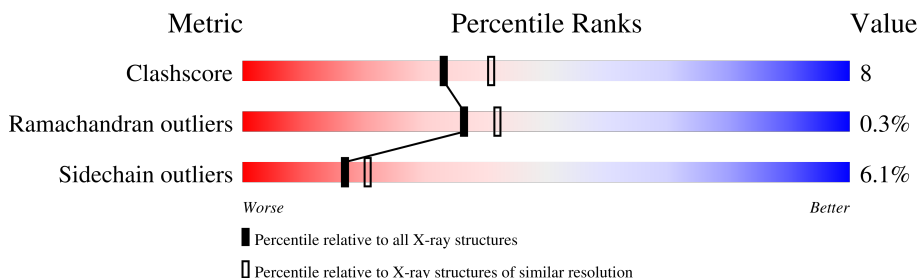
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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))
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	283	77% 20% .
1	B	283	76% 20% . .
1	C	283	76% 21% .
1	D	283	75% 21% .
1	E	283	74% 23% .
2	F	29	52% 34% 10% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11261 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 COAT PROTEIN VP1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	283	Total	C	N	O	S	0	0	0
			2213	1400	371	431	11			
1	B	283	Total	C	N	O	S	0	0	0
			2213	1400	371	431	11			
1	C	283	Total	C	N	O	S	0	0	0
			2213	1400	371	431	11			
1	D	283	Total	C	N	O	S	0	0	0
			2213	1400	371	431	11			
1	E	283	Total	C	N	O	S	0	0	0
			2213	1400	371	431	11			

- Molecule 2 is a protein called FRAGMENT OF COAT PROTEIN VP2.

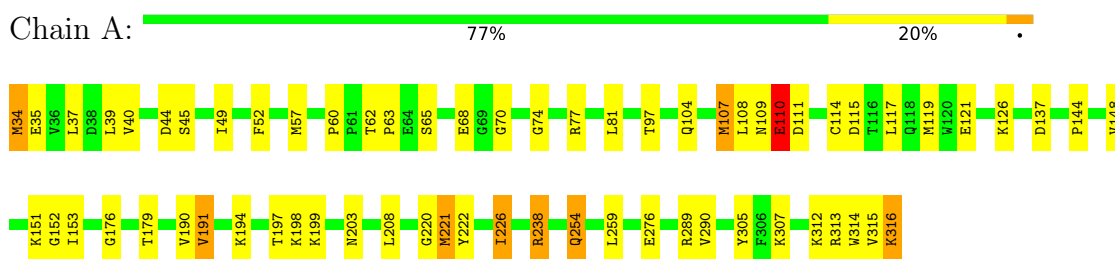
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	F	29	Total	C	N	O	S	0	0	0
			196	124	36	35	1			

3 Residue-property plots

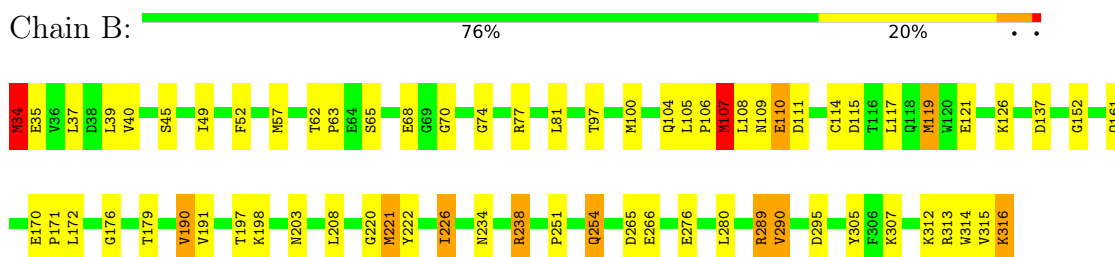
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS was not executed.

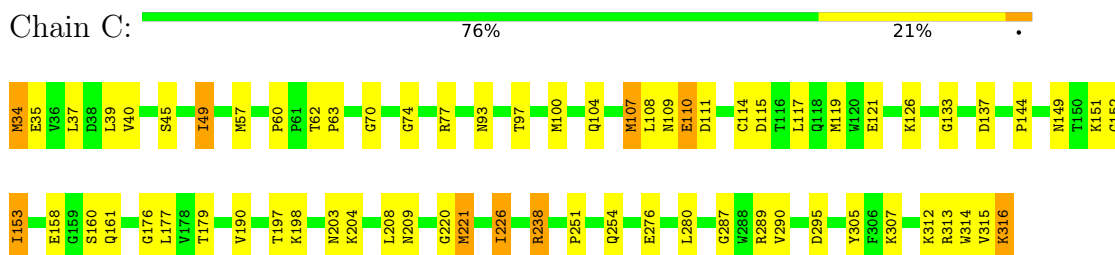
• Molecule 1: COAT PROTEIN VP1



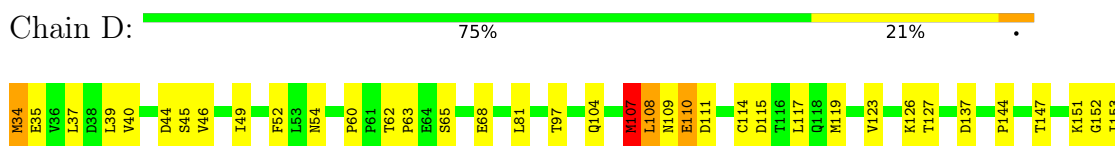
• Molecule 1: COAT PROTEIN VP1

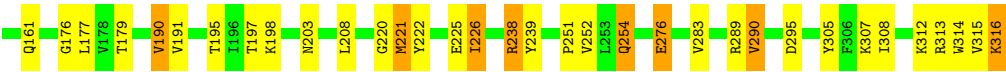


• Molecule 1: COAT PROTEIN VP1

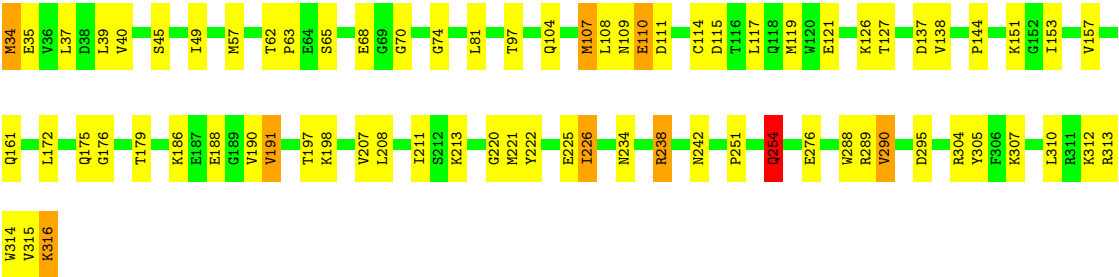


• Molecule 1: COAT PROTEIN VP1





• Molecule 1: COAT PROTEIN VP1



• Molecule 2: FRAGMENT OF COAT PROTEIN VP2



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	219.00 Å 219.00 Å 99.00 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.20	Depositor
% Data completeness (in resolution range)	95.0 (20.00-2.20)	Depositor
R_{merge}	0.09	Depositor
R_{sym}	0.15	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.230 , 0.260	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	11261	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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	1.26	5/2266 (0.2%)	1.28	12/3088 (0.4%)
1	B	1.26	7/2266 (0.3%)	1.26	11/3088 (0.4%)
1	C	1.26	7/2266 (0.3%)	1.28	10/3088 (0.3%)
1	D	1.33	10/2266 (0.4%)	1.25	12/3088 (0.4%)
1	E	1.29	8/2266 (0.4%)	1.26	9/3088 (0.3%)
2	F	0.80	0/201	1.33	3/271 (1.1%)
All	All	1.27	37/11531 (0.3%)	1.26	57/15711 (0.4%)

All (37) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	221	MET	SD-CE	14.02	2.14	1.79
1	D	221	MET	SD-CE	12.29	2.10	1.79
1	B	221	MET	SD-CE	11.75	2.08	1.79
1	B	107	MET	SD-CE	10.86	2.06	1.79
1	C	221	MET	SD-CE	10.31	2.05	1.79
1	A	57	MET	SD-CE	8.58	2.00	1.79
1	E	221	MET	SD-CE	8.43	2.00	1.79
1	D	107	MET	SD-CE	7.95	1.99	1.79
1	E	57	MET	SD-CE	7.84	1.99	1.79
1	E	107	MET	SD-CE	7.49	1.98	1.79
1	E	290	VAL	CA-CB	7.46	1.64	1.54
1	C	107	MET	SD-CE	7.40	1.98	1.79
1	E	238	ARG	NE-CZ	7.00	1.40	1.33
1	A	107	MET	SD-CE	6.95	1.97	1.79
1	B	57	MET	SD-CE	6.89	1.96	1.79
1	D	238	ARG	NE-CZ	6.70	1.40	1.33
1	D	123	VAL	CA-CB	6.36	1.62	1.54
1	B	254	GLN	CA-C	-6.32	1.45	1.52
1	C	153	ILE	CA-CB	6.23	1.64	1.54
1	D	276	GLU	C-N	6.20	1.36	1.33
1	D	254	GLN	CA-C	-6.13	1.45	1.52
1	C	238	ARG	CZ-NH2	6.05	1.41	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	238	ARG	NE-CZ	6.01	1.39	1.33
1	A	238	ARG	NE-CZ	5.95	1.39	1.33
1	C	158	GLU	CA-C	-5.87	1.45	1.52
1	C	238	ARG	NE-CZ	5.77	1.39	1.33
1	A	191	VAL	CA-CB	-5.73	1.47	1.54
1	E	254	GLN	CA-C	-5.64	1.46	1.52
1	D	46	VAL	CA-CB	-5.55	1.47	1.54
1	D	290	VAL	CA-CB	5.54	1.61	1.54
1	B	290	VAL	CA-CB	5.38	1.61	1.54
1	D	239	TYR	CA-C	-5.26	1.46	1.52
1	B	34	MET	SD-CE	5.23	1.92	1.79
1	E	191	VAL	CA-CB	-5.23	1.48	1.54
1	E	138	VAL	CA-CB	5.20	1.61	1.54
1	C	57	MET	SD-CE	5.12	1.92	1.79
1	D	308	ILE	CA-CB	5.08	1.60	1.54

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	109	ASN	N-CA-C	8.69	122.68	108.52
1	A	109	ASN	N-CA-C	8.62	122.58	108.52
1	D	109	ASN	N-CA-C	8.29	122.04	108.52
1	B	109	ASN	N-CA-C	8.04	121.62	108.52
1	E	109	ASN	N-CA-C	7.84	121.30	108.52
2	F	11	SER	N-CA-C	7.53	126.84	110.80
1	B	289	ARG	N-CA-C	-7.26	98.40	109.95
1	A	289	ARG	N-CA-C	-7.04	98.76	109.95
1	C	108	LEU	N-CA-C	6.98	122.47	113.18
1	A	108	LEU	N-CA-C	6.92	122.39	113.18
1	D	289	ARG	N-CA-C	-6.71	99.27	109.95
1	C	226	ILE	CB-CA-C	-6.71	104.34	110.91
1	D	226	ILE	CB-CA-C	-6.64	104.40	110.91
1	B	108	LEU	N-CA-C	6.63	122.00	113.18
1	D	108	LEU	N-CA-C	6.62	121.98	113.18
1	E	289	ARG	N-CA-C	-6.57	99.50	109.95
1	E	108	LEU	N-CA-C	6.48	121.80	113.18
1	E	226	ILE	CB-CA-C	-6.45	105.10	111.30
1	C	289	ARG	N-CA-C	-6.45	99.79	110.17
1	B	126	LYS	N-CA-C	-6.44	98.02	108.52
1	B	226	ILE	CB-CA-C	-6.39	104.64	110.91
1	D	126	LYS	N-CA-C	-6.37	98.14	108.52
1	E	126	LYS	N-CA-C	-6.25	98.34	108.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	126	LYS	N-CA-C	-6.21	98.39	108.52
1	A	226	ILE	CB-CA-C	-6.13	105.42	111.30
1	D	60	PRO	CA-C-N	6.09	125.98	119.28
1	D	60	PRO	C-N-CA	6.09	125.98	119.28
2	F	12	HIS	N-CA-C	6.05	119.39	109.76
1	D	45	SER	N-CA-C	-6.02	105.77	113.23
1	C	60	PRO	CA-C-N	5.98	125.86	119.28
1	C	60	PRO	C-N-CA	5.98	125.86	119.28
1	C	74	GLY	N-CA-C	-5.93	106.34	114.64
1	A	60	PRO	CA-C-N	5.89	125.76	119.28
1	A	60	PRO	C-N-CA	5.89	125.76	119.28
1	A	126	LYS	N-CA-C	-5.56	99.67	108.73
1	A	45	SER	N-CA-C	-5.52	106.38	113.23
1	B	52	PHE	N-CA-C	-5.44	100.53	109.40
1	D	52	PHE	N-CA-C	-5.42	100.57	109.40
1	D	177	LEU	N-CA-C	-5.31	99.48	110.80
1	C	45	SER	N-CA-C	-5.29	106.67	113.23
1	A	74	GLY	N-CA-C	-5.28	107.25	114.64
1	B	45	SER	N-CA-C	-5.22	106.75	113.23
2	F	8	GLY	N-CA-C	5.18	125.47	113.18
1	C	177	LEU	N-CA-C	-5.18	99.76	110.80
1	E	45	SER	N-CA-C	-5.15	106.84	113.23
1	E	225	GLU	N-CA-C	-5.13	106.79	113.16
1	A	254	GLN	CA-CB-CG	-5.12	103.86	114.10
1	A	52	PHE	N-CA-C	-5.10	100.72	109.24
1	E	254	GLN	CA-CB-CG	-5.10	103.89	114.10
1	A	77	ARG	N-CA-C	-5.07	102.50	110.36
1	B	190	VAL	CB-CA-C	-5.06	102.08	111.18
1	B	74	GLY	N-CA-C	-5.03	107.60	114.64
1	D	195	THR	N-CA-C	-5.02	105.81	111.28
1	D	190	VAL	CB-CA-C	-5.01	102.15	111.18
1	B	77	ARG	N-CA-C	-5.01	104.06	110.53
1	B	254	GLN	CA-CB-CG	-5.01	104.08	114.10
1	E	74	GLY	N-CA-C	-5.01	107.63	114.64

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	2213	0	2181	36	1
1	B	2213	0	2181	40	0
1	C	2213	0	2181	37	1
1	D	2213	0	2181	36	1
1	E	2213	0	2181	40	3
2	F	196	0	192	16	0
All	All	11261	0	11097	175	3

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 (175) 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:C:221:MET:SD	1:C:221:MET:CE	2.05	1.44
1:B:107:MET:SD	1:B:107:MET:CE	2.06	1.43
1:B:221:MET:SD	1:B:221:MET:CE	2.08	1.40
1:D:221:MET:CE	1:D:221:MET:SD	2.10	1.38
1:A:221:MET:CE	1:A:221:MET:SD	2.14	1.35
1:A:119:MET:HE3	1:A:315:VAL:HG21	1.26	1.18
1:D:119:MET:HE3	1:D:315:VAL:HG21	1.22	1.12
1:B:119:MET:HE3	1:B:315:VAL:HG21	1.23	1.10
1:C:119:MET:HE3	1:C:315:VAL:HG21	1.27	1.09
1:E:119:MET:HE3	1:E:315:VAL:HG21	1.28	1.07
1:E:254:GLN:O	2:F:10:ALA:HB3	1.58	1.03
1:A:259:LEU:HD13	2:F:15:VAL:HG21	1.56	0.87
1:A:34:MET:HA	1:A:34:MET:CE	2.09	0.82
1:C:34:MET:CE	1:C:34:MET:HA	2.08	0.82
1:A:259:LEU:CD1	2:F:15:VAL:HG21	2.10	0.82
1:D:34:MET:HA	1:D:34:MET:CE	2.10	0.82
1:E:34:MET:HA	1:E:34:MET:CE	2.12	0.79
1:B:34:MET:HA	1:B:34:MET:CE	2.13	0.78
1:C:34:MET:HA	1:C:34:MET:HE2	1.67	0.75
1:C:111:ASP:HB3	1:C:114:CYS:HB2	1.73	0.71
1:B:266:GLU:HA	2:F:18:ASP:O	1.91	0.70
1:B:34:MET:HA	1:B:34:MET:HE2	1.73	0.69
1:B:111:ASP:HB3	1:B:114:CYS:HB2	1.75	0.68
1:A:34:MET:HA	1:A:34:MET:HE2	1.77	0.67
1:D:34:MET:HA	1:D:34:MET:HE3	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:111:ASP:HB3	1:D:114:CYS:HB2	1.78	0.65
1:A:34:MET:HA	1:A:34:MET:HE3	1.79	0.65
1:E:111:ASP:HB3	1:E:114:CYS:HB2	1.81	0.63
1:C:197:THR:O	1:C:198:LYS:HB2	1.99	0.63
1:E:34:MET:HA	1:E:34:MET:HE3	1.80	0.62
1:D:34:MET:HA	1:D:34:MET:HE2	1.81	0.62
1:E:242:ASN:ND2	2:F:7:GLY:C	2.59	0.61
1:E:34:MET:HA	1:E:34:MET:HE2	1.80	0.60
1:A:111:ASP:HB3	1:A:114:CYS:HB2	1.81	0.60
1:B:197:THR:O	1:B:198:LYS:HB2	2.02	0.60
1:B:110:GLU:OE2	1:B:110:GLU:HA	2.05	0.57
1:A:259:LEU:HD13	2:F:15:VAL:HG11	1.87	0.56
1:E:242:ASN:HD22	2:F:7:GLY:C	2.12	0.56
1:E:197:THR:O	1:E:198:LYS:HB2	2.04	0.55
1:B:39:LEU:HD23	1:B:313:ARG:HD2	1.89	0.55
1:E:254:GLN:O	2:F:10:ALA:CB	2.45	0.55
1:C:110:GLU:HA	1:C:110:GLU:OE2	2.06	0.55
1:C:153:ILE:HA	1:D:81:LEU:HD21	1.88	0.55
1:D:197:THR:O	1:D:198:LYS:HB2	2.06	0.55
1:C:39:LEU:HD23	1:C:313:ARG:HD2	1.89	0.55
1:C:305:TYR:OH	1:C:307:LYS:HE2	2.07	0.54
1:B:37:LEU:HD11	1:B:316:LYS:HB2	1.90	0.54
1:E:305:TYR:OH	1:E:307:LYS:HE2	2.08	0.54
1:D:40:VAL:HB	1:D:312:LYS:HB2	1.89	0.53
1:A:37:LEU:HD11	1:A:316:LYS:HB2	1.90	0.53
1:C:37:LEU:HD11	1:C:316:LYS:HB2	1.90	0.53
1:C:40:VAL:HB	1:C:312:LYS:HB2	1.89	0.53
1:E:40:VAL:HB	1:E:312:LYS:HB2	1.91	0.53
1:A:259:LEU:HD22	2:F:15:VAL:CG2	2.39	0.53
1:A:110:GLU:OE2	1:A:110:GLU:HA	2.09	0.53
1:B:70:GLY:O	1:C:203:ASN:HB2	2.09	0.53
1:E:304:ARG:NH2	2:F:8:GLY:O	2.36	0.53
1:A:197:THR:O	1:A:198:LYS:HB2	2.09	0.52
1:B:137:ASP:OD1	1:B:137:ASP:C	2.51	0.52
1:E:39:LEU:HD23	1:E:313:ARG:HD2	1.90	0.52
1:A:40:VAL:HB	1:A:312:LYS:HB2	1.91	0.52
1:D:37:LEU:HD11	1:D:316:LYS:HB2	1.91	0.52
1:E:137:ASP:OD1	1:E:137:ASP:C	2.53	0.52
1:A:259:LEU:HD13	2:F:15:VAL:CG2	2.35	0.51
1:B:305:TYR:OH	1:B:307:LYS:HE2	2.10	0.51
1:D:110:GLU:HA	1:D:110:GLU:OE2	2.09	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:97:THR:OG1	1:B:220:GLY:HA2	2.11	0.51
1:C:34:MET:HA	1:C:34:MET:HE3	1.88	0.51
1:A:81:LEU:HD21	1:E:153:ILE:HA	1.92	0.50
1:D:252:VAL:HG21	2:F:4:GLY:H	1.77	0.50
1:E:35:GLU:HB3	1:E:316:LYS:HB3	1.93	0.50
1:C:35:GLU:HB3	1:C:316:LYS:HB3	1.93	0.50
1:C:97:THR:OG1	1:C:220:GLY:HA2	2.11	0.50
1:D:39:LEU:HD23	1:D:313:ARG:HD2	1.94	0.50
1:C:40:VAL:HG22	1:C:314:TRP:CE2	2.47	0.50
1:C:221:MET:CE	1:C:221:MET:CG	2.90	0.50
1:B:34:MET:HA	1:B:34:MET:HE3	1.92	0.50
1:A:115:ASP:OD2	1:A:316:LYS:HD2	2.12	0.50
1:B:266:GLU:HA	2:F:19:TRP:HA	1.92	0.49
1:D:35:GLU:HB3	1:D:316:LYS:HB3	1.92	0.49
1:A:191:VAL:HG11	1:A:222:TYR:CE1	2.47	0.49
1:E:110:GLU:OE2	1:E:110:GLU:HA	2.11	0.49
1:A:203:ASN:HB2	1:E:70:GLY:O	2.12	0.49
1:B:35:GLU:HB3	1:B:316:LYS:HB3	1.94	0.49
1:D:144:PRO:HB2	1:D:151:LYS:O	2.13	0.49
1:A:62:THR:HA	1:A:63:PRO:C	2.37	0.49
1:E:37:LEU:HD11	1:E:316:LYS:HB2	1.94	0.49
1:A:35:GLU:HB3	1:A:316:LYS:HB3	1.93	0.49
1:B:40:VAL:HB	1:B:312:LYS:HB2	1.94	0.49
1:A:39:LEU:HD23	1:A:313:ARG:HD2	1.95	0.48
1:D:305:TYR:OH	1:D:307:LYS:HE2	2.12	0.48
1:D:97:THR:OG1	1:D:220:GLY:HA2	2.14	0.48
1:B:40:VAL:HG22	1:B:314:TRP:CE2	2.49	0.48
1:C:62:THR:HA	1:C:63:PRO:C	2.38	0.48
1:E:62:THR:HA	1:E:63:PRO:C	2.39	0.47
1:A:305:TYR:OH	1:A:307:LYS:HE2	2.15	0.47
1:B:121:GLU:OE2	1:B:313:ARG:HD3	2.15	0.47
1:A:97:THR:OG1	1:A:220:GLY:HA2	2.15	0.47
1:D:40:VAL:HG22	1:D:314:TRP:CE2	2.49	0.47
1:D:62:THR:HA	1:D:63:PRO:C	2.40	0.47
1:E:115:ASP:OD2	1:E:316:LYS:HD2	2.15	0.47
1:E:176:GLY:HA2	1:E:226:ILE:O	2.15	0.47
1:B:115:ASP:OD2	1:B:316:LYS:HD2	2.15	0.47
1:B:266:GLU:CA	2:F:18:ASP:O	2.62	0.46
1:B:62:THR:HA	1:B:63:PRO:C	2.39	0.46
1:B:100:MET:HA	1:B:280:LEU:O	2.15	0.46
1:B:265:ASP:O	2:F:19:TRP:HA	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:137:ASP:OD1	1:A:137:ASP:C	2.55	0.46
1:E:40:VAL:HG22	1:E:314:TRP:CE2	2.51	0.46
1:E:144:PRO:HB2	1:E:151:LYS:O	2.16	0.45
1:D:221:MET:CE	1:D:221:MET:CG	2.93	0.45
1:C:110:GLU:OE2	1:C:110:GLU:CA	2.65	0.45
1:C:70:GLY:O	1:D:203:ASN:HB2	2.16	0.45
1:B:191:VAL:HG11	1:B:222:TYR:CE1	2.52	0.45
1:C:176:GLY:HA2	1:C:226:ILE:O	2.17	0.45
1:D:127:THR:HA	1:D:305:TYR:O	2.17	0.45
1:A:121:GLU:OE2	1:A:313:ARG:HD3	2.16	0.45
1:D:153:ILE:HA	1:E:81:LEU:HD21	1.99	0.45
1:A:40:VAL:HG22	1:A:314:TRP:CE2	2.52	0.45
1:B:110:GLU:OE2	1:B:110:GLU:CA	2.65	0.44
1:B:176:GLY:HA2	1:B:226:ILE:O	2.17	0.44
1:A:194:LYS:HG3	1:A:199:LYS:O	2.18	0.44
1:C:144:PRO:HB2	1:C:151:LYS:O	2.18	0.44
1:A:153:ILE:HA	1:B:81:LEU:HD21	1.99	0.44
1:D:54:ASN:OD1	1:E:207:VAL:HB	2.18	0.44
1:D:115:ASP:OD2	1:D:316:LYS:HD2	2.17	0.44
1:B:65:SER:HB3	1:B:68:GLU:HB2	1.99	0.44
1:C:137:ASP:C	1:C:137:ASP:OD1	2.60	0.44
1:D:65:SER:HB3	1:D:68:GLU:HB2	1.99	0.44
1:B:197:THR:O	1:B:198:LYS:CB	2.66	0.44
1:D:161:GLN:CD	1:D:251:PRO:HA	2.43	0.44
1:B:221:MET:CE	1:B:221:MET:CG	2.96	0.43
1:C:160:SER:O	1:C:287:GLY:HA2	2.18	0.43
1:E:191:VAL:HG11	1:E:222:TYR:CE1	2.53	0.43
1:C:121:GLU:OE2	1:C:313:ARG:HD3	2.17	0.43
1:B:161:GLN:CD	1:B:251:PRO:HA	2.44	0.43
1:A:144:PRO:HB2	1:A:151:LYS:O	2.19	0.43
1:E:121:GLU:OE2	1:E:313:ARG:HD3	2.18	0.43
1:C:115:ASP:OD2	1:C:316:LYS:HD2	2.19	0.43
1:E:127:THR:HA	1:E:305:TYR:O	2.19	0.43
1:A:70:GLY:O	1:B:203:ASN:HB2	2.18	0.43
1:D:137:ASP:OD1	1:D:137:ASP:C	2.61	0.43
1:D:191:VAL:HG11	1:D:222:TYR:CE1	2.54	0.43
1:A:176:GLY:HA2	1:A:226:ILE:O	2.19	0.42
1:A:152:GLY:HA2	1:B:295:ASP:O	2.19	0.42
1:C:152:GLY:HA2	1:D:295:ASP:O	2.20	0.42
1:B:152:GLY:HA2	1:C:295:ASP:O	2.19	0.42
1:C:204:LYS:HB3	1:C:209:ASN:HB2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:100:MET:HA	1:C:280:LEU:O	2.20	0.41
1:A:110:GLU:OE2	1:A:110:GLU:CA	2.67	0.41
1:E:157:VAL:HA	1:E:288:TRP:O	2.20	0.41
1:A:44:ASP:O	1:A:312:LYS:HE3	2.21	0.41
1:E:97:THR:OG1	1:E:220:GLY:HA2	2.20	0.41
1:E:161:GLN:CD	1:E:251:PRO:HA	2.45	0.41
1:E:254:GLN:HE22	2:F:12:HIS:CG	2.39	0.41
1:C:77:ARG:HD2	1:C:77:ARG:HA	1.86	0.41
1:C:77:ARG:HG3	1:C:93:ASN:ND2	2.36	0.41
1:C:161:GLN:CD	1:C:251:PRO:HA	2.46	0.41
1:D:152:GLY:HA2	1:E:295:ASP:O	2.20	0.41
1:E:172:LEU:HD12	1:E:234:ASN:OD1	2.20	0.41
1:D:44:ASP:O	1:D:312:LYS:HE3	2.21	0.41
1:E:65:SER:HB3	1:E:68:GLU:HB2	2.03	0.41
1:E:121:GLU:O	1:E:310:LEU:HA	2.21	0.41
1:E:198:LYS:HD3	1:E:198:LYS:HA	1.72	0.41
1:D:107:MET:O	1:D:108:LEU:HD23	2.21	0.40
1:B:170:GLU:HB2	1:B:171:PRO:HD2	2.04	0.40
1:A:65:SER:HB3	1:A:68:GLU:HB2	2.02	0.40
1:C:49:ILE:HG23	1:C:49:ILE:HD12	1.74	0.40
1:C:198:LYS:HA	1:C:198:LYS:HD3	1.69	0.40
1:D:110:GLU:OE2	1:D:110:GLU:CA	2.69	0.40
1:D:176:GLY:HA2	1:D:226:ILE:O	2.20	0.40
1:B:105:LEU:HB3	1:B:106:PRO:HD2	2.03	0.40
1:B:172:LEU:HD12	1:B:234:ASN:OD1	2.21	0.40
1:C:133:GLY:O	1:D:225:GLU:HA	2.21	0.40
1:E:175:GLN:OE1	1:E:213:LYS:HE2	2.21	0.40

All (3) 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:C:149:ASN:OD1	1:E:186:LYS:NZ[3_564]	2.03	0.17
1:D:147:THR:CG2	1:E:188:GLU:OE2[3_564]	2.07	0.13
1:A:148:VAL:CG2	1:E:211:ILE:CD1[3_564]	2.15	0.05

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	281/283 (99%)	271 (96%)	9 (3%)	1 (0%)	30	34
1	B	281/283 (99%)	270 (96%)	11 (4%)	0	100	100
1	C	281/283 (99%)	272 (97%)	9 (3%)	0	100	100
1	D	281/283 (99%)	269 (96%)	12 (4%)	0	100	100
1	E	281/283 (99%)	270 (96%)	11 (4%)	0	100	100
2	F	27/29 (93%)	18 (67%)	5 (18%)	4 (15%)	0	0
All	All	1432/1444 (99%)	1370 (96%)	57 (4%)	5 (0%)	36	42

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	8	GLY
2	F	11	SER
2	F	10	ALA
2	F	23	LEU
1	A	110	GLU

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	249/249 (100%)	235 (94%)	14 (6%)	19	24
1	B	249/249 (100%)	233 (94%)	16 (6%)	16	19
1	C	249/249 (100%)	235 (94%)	14 (6%)	19	24

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	249/249 (100%)	234 (94%)	15 (6%)	17	21
1	E	249/249 (100%)	235 (94%)	14 (6%)	19	24
2	F	17/17 (100%)	13 (76%)	4 (24%)	1	0
All	All	1262/1262 (100%)	1185 (94%)	77 (6%)	17	20

All (77) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	34	MET
1	A	49	ILE
1	A	104	GLN
1	A	107	MET
1	A	110	GLU
1	A	117	LEU
1	A	179	THR
1	A	190	VAL
1	A	208	LEU
1	A	238	ARG
1	A	254	GLN
1	A	276	GLU
1	A	290	VAL
1	A	316	LYS
1	B	34	MET
1	B	49	ILE
1	B	104	GLN
1	B	107	MET
1	B	110	GLU
1	B	117	LEU
1	B	119	MET
1	B	179	THR
1	B	190	VAL
1	B	208	LEU
1	B	238	ARG
1	B	254	GLN
1	B	276	GLU
1	B	289	ARG
1	B	290	VAL
1	B	316	LYS
1	C	34	MET
1	C	49	ILE
1	C	104	GLN

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Mol	Chain	Res	Type
1	C	107	MET
1	C	110	GLU
1	C	117	LEU
1	C	179	THR
1	C	190	VAL
1	C	208	LEU
1	C	238	ARG
1	C	254	GLN
1	C	276	GLU
1	C	290	VAL
1	C	316	LYS
1	D	34	MET
1	D	49	ILE
1	D	104	GLN
1	D	107	MET
1	D	110	GLU
1	D	117	LEU
1	D	179	THR
1	D	190	VAL
1	D	208	LEU
1	D	238	ARG
1	D	254	GLN
1	D	276	GLU
1	D	283	VAL
1	D	290	VAL
1	D	316	LYS
1	E	34	MET
1	E	49	ILE
1	E	104	GLN
1	E	107	MET
1	E	110	GLU
1	E	117	LEU
1	E	179	THR
1	E	190	VAL
1	E	208	LEU
1	E	238	ARG
1	E	254	GLN
1	E	276	GLU
1	E	290	VAL
1	E	316	LYS
2	F	13	GLN
2	F	14	ARG

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Mol	Chain	Res	Type
2	F	25	LEU
2	F	27	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	142	ASN
1	E	242	ASN
1	E	254	GLN

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](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.