



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:56 PM UTC

PDB ID : 7JZJ / pdb_00007jzj
Title : Crystal structure demonstrating CTD-CTD interactions of Zaire Ebola virus VP40 dimer
Authors : Norris, M.J.; Bornholdt, Z.A.; Saphire, E.O.
Deposited on : 2020-09-02
Resolution : 2.46 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

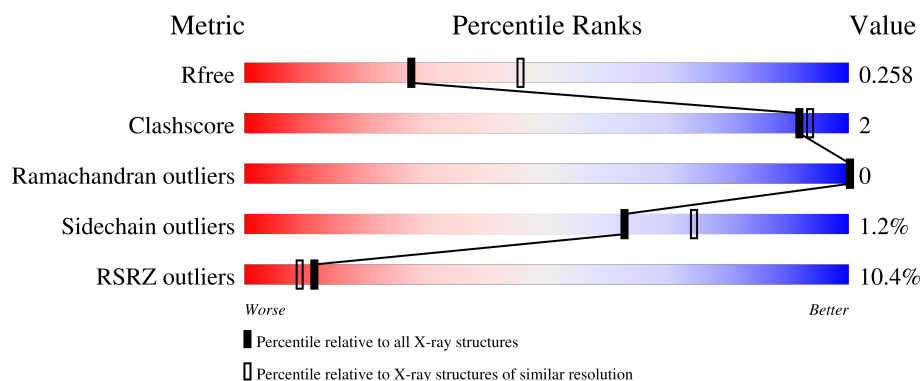
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.46 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	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

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	297	9% 77% 6% 16%
1	B	297	6% 85% 5% 10%
1	C	297	12% 73% 7% 19%
1	D	297	8% 78% 5% 17%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 14511 atoms, of which 7064 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 Matrix protein VP40.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	249	Total	C	H	N	O	S	0	0	0
			3611	1180	1777	308	339	7			
1	B	267	Total	C	H	N	O	S	0	0	0
			3943	1275	1961	329	370	8			
1	C	240	Total	C	H	N	O	S	0	0	0
			3300	1098	1587	292	316	7			
1	D	246	Total	C	H	N	O	S	0	0	0
			3540	1162	1739	301	330	8			

There are 52 discrepancies between the modelled and reference sequences:

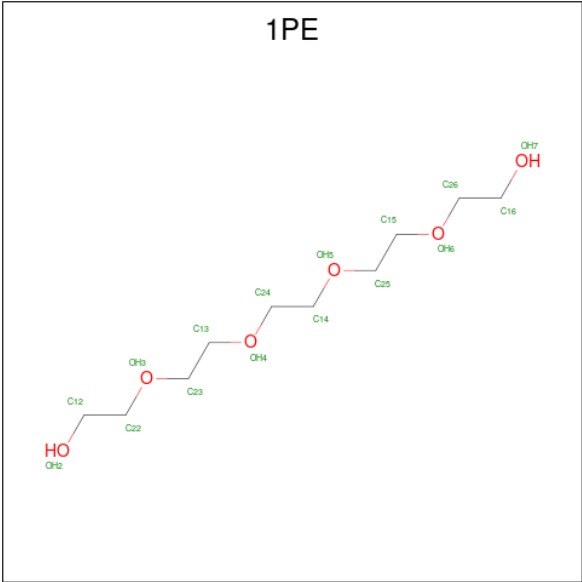
Chain	Residue	Modelled	Actual	Comment	Reference
A	30	MET	-	initiating methionine	UNP Q05128
A	31	ALA	-	expression tag	UNP Q05128
A	32	HIS	-	expression tag	UNP Q05128
A	33	HIS	-	expression tag	UNP Q05128
A	34	HIS	-	expression tag	UNP Q05128
A	35	HIS	-	expression tag	UNP Q05128
A	36	HIS	-	expression tag	UNP Q05128
A	37	HIS	-	expression tag	UNP Q05128
A	38	VAL	-	expression tag	UNP Q05128
A	39	ASP	-	expression tag	UNP Q05128
A	40	ASP	-	expression tag	UNP Q05128
A	41	ASP	-	expression tag	UNP Q05128
A	42	ASP	-	expression tag	UNP Q05128
B	30	MET	-	initiating methionine	UNP Q05128
B	31	ALA	-	expression tag	UNP Q05128
B	32	HIS	-	expression tag	UNP Q05128
B	33	HIS	-	expression tag	UNP Q05128
B	34	HIS	-	expression tag	UNP Q05128
B	35	HIS	-	expression tag	UNP Q05128
B	36	HIS	-	expression tag	UNP Q05128
B	37	HIS	-	expression tag	UNP Q05128

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Chain	Residue	Modelled	Actual	Comment	Reference
B	38	VAL	-	expression tag	UNP Q05128
B	39	ASP	-	expression tag	UNP Q05128
B	40	ASP	-	expression tag	UNP Q05128
B	41	ASP	-	expression tag	UNP Q05128
B	42	ASP	-	expression tag	UNP Q05128
C	30	MET	-	initiating methionine	UNP Q05128
C	31	ALA	-	expression tag	UNP Q05128
C	32	HIS	-	expression tag	UNP Q05128
C	33	HIS	-	expression tag	UNP Q05128
C	34	HIS	-	expression tag	UNP Q05128
C	35	HIS	-	expression tag	UNP Q05128
C	36	HIS	-	expression tag	UNP Q05128
C	37	HIS	-	expression tag	UNP Q05128
C	38	VAL	-	expression tag	UNP Q05128
C	39	ASP	-	expression tag	UNP Q05128
C	40	ASP	-	expression tag	UNP Q05128
C	41	ASP	-	expression tag	UNP Q05128
C	42	ASP	-	expression tag	UNP Q05128
D	30	MET	-	initiating methionine	UNP Q05128
D	31	ALA	-	expression tag	UNP Q05128
D	32	HIS	-	expression tag	UNP Q05128
D	33	HIS	-	expression tag	UNP Q05128
D	34	HIS	-	expression tag	UNP Q05128
D	35	HIS	-	expression tag	UNP Q05128
D	36	HIS	-	expression tag	UNP Q05128
D	37	HIS	-	expression tag	UNP Q05128
D	38	VAL	-	expression tag	UNP Q05128
D	39	ASP	-	expression tag	UNP Q05128
D	40	ASP	-	expression tag	UNP Q05128
D	41	ASP	-	expression tag	UNP Q05128
D	42	ASP	-	expression tag	UNP Q05128

- Molecule 2 is PENTAETHYLENE GLYCOL (CCD ID: 1PE) (formula: C₁₀H₂₂O₆).



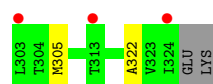
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			16	10	6		
2	B	1	Total	C	O	0	0
			16	10	6		
2	D	1	Total	C	O	0	0
			16	10	6		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	15	Total	O	0	0
			15	15		
3	B	21	Total	O	0	0
			21	21		
3	C	17	Total	O	0	0
			17	17		
3	D	16	Total	O	0	0
			16	16		

- Molecule 1: Matrix protein VP40





4 Data and refinement statistics

Property	Value	Source
Space group	P 62	Depositor
Cell constants a, b, c, α , β , γ	159.94Å 159.94Å 89.76Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	46.17 – 2.46 46.17 – 2.46	Depositor EDS
% Data completeness (in resolution range)	91.7 (46.17-2.46) 91.9 (46.17-2.46)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.59 (at 2.45Å)	Xtriage
Refinement program	PHENIX 1.17.1-3660	Depositor
R, R_{free}	0.249 , 0.258 0.252 , 0.258	Depositor DCC
R_{free} test set	2191 reflections (4.60%)	wwPDB-VP
Wilson B-factor (Å ²)	69.3	Xtriage
Anisotropy	0.346	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 76.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.029 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	14511	wwPDB-VP
Average B, all atoms (Å ²)	99.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.78% 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: 1PE

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	1.02	4/1880 (0.2%)	1.43	7/2582 (0.3%)
1	B	1.01	1/2034 (0.0%)	1.45	5/2796 (0.2%)
1	C	1.07	3/1755 (0.2%)	1.45	9/2414 (0.4%)
1	D	1.00	2/1847 (0.1%)	1.38	4/2541 (0.2%)
All	All	1.03	10/7516 (0.1%)	1.43	25/10333 (0.2%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	108	PHE	C-O	-8.48	1.14	1.24
1	A	148	ARG	C-O	-6.38	1.15	1.24
1	A	108	PHE	C-O	-6.25	1.16	1.24
1	D	283	PRO	C-O	-5.99	1.16	1.23
1	A	164	PRO	C-O	-5.75	1.20	1.25
1	C	283	PRO	C-O	-5.59	1.16	1.23
1	A	70	SER	C-O	-5.37	1.17	1.24
1	D	97	PRO	C-O	-5.30	1.19	1.24
1	C	181	LEU	C-O	-5.26	1.17	1.24
1	C	143	PRO	C-O	-5.24	1.17	1.23

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	216	ILE	N-CA-C	-6.73	102.61	110.21
1	C	53	PRO	N-CA-C	6.05	120.49	111.41
1	B	108	PHE	CA-C-O	-6.05	114.13	120.55
1	C	89	MET	CA-C-O	-5.87	114.72	121.47
1	C	307	ILE	N-CA-C	-5.79	99.19	107.77
1	C	107	SER	CA-C-O	-5.70	115.26	121.87
1	D	53	PRO	N-CA-C	5.68	120.20	111.57

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	306	VAL	CA-C-O	-5.64	116.04	121.63
1	C	255	THR	N-CA-C	-5.56	101.62	109.96
1	B	53	PRO	N-CA-C	5.54	119.99	111.57
1	C	89	MET	N-CA-C	-5.50	102.33	110.48
1	A	70	SER	CA-C-O	-5.49	114.77	120.70
1	C	73	PHE	CA-C-O	-5.48	114.78	120.70
1	B	128	ALA	N-CA-C	-5.42	105.38	111.28
1	A	108	PHE	CA-C-O	-5.40	114.82	120.55
1	C	252	ILE	N-CA-C	-5.38	106.56	111.67
1	B	199	SER	N-CA-C	-5.37	105.42	111.28
1	A	268	VAL	N-CA-C	-5.36	105.16	110.62
1	D	283	PRO	CA-C-O	-5.29	115.57	121.23
1	D	238	GLN	N-CA-C	-5.25	105.56	111.28
1	C	321	PRO	N-CA-C	5.21	120.38	113.40
1	B	306	VAL	CA-C-O	-5.20	116.21	121.67
1	A	53	PRO	N-CA-C	5.12	119.35	111.57
1	D	237	ILE	N-CA-C	-5.08	106.11	113.07
1	A	216	ILE	CA-C-O	-5.04	116.53	122.13

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	1834	1777	1777	6	0
1	B	1982	1961	1961	5	5
1	C	1713	1587	1589	6	0
1	D	1801	1739	1746	6	5
2	A	16	0	22	1	0
2	B	16	0	22	2	0
2	D	16	0	22	2	0
3	A	15	0	0	0	0
3	B	21	0	0	0	0
3	C	17	0	0	0	0
3	D	16	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	7447	7064	7139	23	5

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 2.

All (23) 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:B:95:TRP:CZ3	2:B:401:1PE:H261	2.39	0.58
1:C:240:ILE:HG22	1:C:305:MET:HE2	1.84	0.58
1:C:249:ILE:HD11	1:C:303:LEU:HD21	1.85	0.57
1:D:240:ILE:HG22	1:D:305:MET:HE2	1.87	0.55
1:B:291:LYS:HG3	2:B:401:1PE:H142	1.89	0.55
1:A:219:PRO:O	1:A:220:ASN:C	2.50	0.54
1:D:257:ASN:HB2	2:D:401:1PE:H131	1.90	0.54
1:A:94:ILE:HD12	1:A:163:LEU:CD2	2.39	0.53
1:C:75:LEU:HG	1:C:98:LEU:HD11	1.92	0.52
1:D:241:MET:HG2	1:D:305:MET:HE1	1.94	0.50
1:D:145:HIS:CE1	1:D:150:LEU:HD12	2.48	0.49
1:A:145:HIS:CE1	1:A:150:LEU:HD12	2.48	0.49
1:C:81:VAL:HB	1:C:89:MET:CB	2.44	0.47
1:C:241:MET:CG	1:C:305:MET:HE1	2.46	0.46
1:B:94:ILE:HD12	1:B:163:LEU:CD2	2.47	0.45
1:B:145:HIS:CE1	1:B:150:LEU:HD12	2.53	0.44
1:D:74:ILE:HG21	1:D:95:TRP:CE2	2.53	0.43
1:A:93:PRO:C	1:A:94:ILE:HD13	2.45	0.42
1:B:94:ILE:HD12	1:B:163:LEU:HD23	2.01	0.41
1:D:95:TRP:CZ3	2:D:401:1PE:H222	2.56	0.41
1:A:108:PHE:HE1	1:A:183:THR:HG21	1.86	0.40
1:C:213:LEU:HD11	1:C:293:ILE:HD13	2.04	0.40
1:A:95:TRP:CE3	2:A:401:1PE:H222	2.56	0.40

All (5) 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:B:274:LYS:NZ	1:D:322:ALA:O[5_554]	1.38	0.82
1:B:274:LYS:CE	1:D:322:ALA:O[5_554]	1.72	0.48
1:B:274:LYS:HZ1	1:D:322:ALA:O[5_554]	1.23	0.37
1:B:274:LYS:HE3	1:D:322:ALA:O[5_554]	1.47	0.13

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:274:LYS:HZ2	1:D:322:ALA:O[5_554]	1.48	0.12

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	239/297 (80%)	237 (99%)	2 (1%)	0	100	100
1	B	261/297 (88%)	258 (99%)	3 (1%)	0	100	100
1	C	224/297 (75%)	219 (98%)	5 (2%)	0	100	100
1	D	236/297 (80%)	235 (100%)	1 (0%)	0	100	100
All	All	960/1188 (81%)	949 (99%)	11 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	195/259 (75%)	192 (98%)	3 (2%)	57	70
1	B	218/259 (84%)	214 (98%)	4 (2%)	51	65
1	C	171/259 (66%)	170 (99%)	1 (1%)	78	86
1	D	191/259 (74%)	190 (100%)	1 (0%)	81	87
All	All	775/1036 (75%)	766 (99%)	9 (1%)	63	74

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	152	ILE
1	A	243	SER
1	A	287	VAL
1	B	65	THR
1	B	261	ILE
1	B	296	ASP
1	B	313	THR
1	C	119	SER
1	D	123	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	170	GLN
1	C	103	GLN
1	D	103	GLN
1	D	124	HIS

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](#)

3 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	$\# Z > 2$
2	1PE	A	401	-	15,15,15	0.56	0	14,14,14	0.24	0
2	1PE	D	401	-	15,15,15	0.41	0	14,14,14	0.34	0
2	1PE	B	401	-	15,15,15	0.37	0	14,14,14	0.41	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
2	1PE	A	401	-	-	5/13/13/13	-
2	1PE	D	401	-	-	5/13/13/13	-
2	1PE	B	401	-	-	8/13/13/13	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	401	1PE	C13-C23-OH3-C22
2	D	401	1PE	C14-C24-OH4-C13
2	A	401	1PE	OH5-C14-C24-OH4
2	A	401	1PE	OH7-C16-C26-OH6
2	B	401	1PE	OH5-C14-C24-OH4
2	A	401	1PE	OH4-C13-C23-OH3
2	D	401	1PE	OH5-C14-C24-OH4
2	B	401	1PE	C14-C24-OH4-C13
2	A	401	1PE	OH6-C15-C25-OH5
2	A	401	1PE	C23-C13-OH4-C24
2	D	401	1PE	C12-C22-OH3-C23
2	B	401	1PE	C24-C14-OH5-C25
2	B	401	1PE	OH6-C15-C25-OH5
2	B	401	1PE	OH7-C16-C26-OH6
2	D	401	1PE	C16-C26-OH6-C15
2	B	401	1PE	C16-C26-OH6-C15
2	D	401	1PE	OH7-C16-C26-OH6
2	B	401	1PE	OH4-C13-C23-OH3

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	1PE	1	0
2	D	401	1PE	2	0
2	B	401	1PE	2	0

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

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	249/297 (83%)	0.76	27 (10%) 11 9	67, 99, 135, 148	0
1	B	267/297 (89%)	0.60	17 (6%) 25 23	59, 85, 133, 164	0
1	C	240/297 (80%)	0.89	35 (14%) 6 4	67, 103, 155, 184	0
1	D	246/297 (82%)	0.82	25 (10%) 12 10	66, 102, 150, 178	0
All	All	1002/1188 (84%)	0.76	104 (10%) 11 9	59, 98, 147, 184	0

All (104) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	220	ASN	5.0
1	D	122	ILE	4.9
1	C	87	VAL	4.9
1	B	231	LEU	4.6
1	C	322	ALA	4.5
1	D	249	ILE	4.2
1	A	138	LEU	4.0
1	B	194	ASP	3.9
1	A	216	ILE	3.8
1	A	280	ASN	3.8
1	C	122	ILE	3.5
1	A	231	LEU	3.4
1	D	324	ILE	3.4
1	C	302	ASP	3.3
1	B	209	PHE	3.3
1	C	280	ASN	3.3
1	D	280	ASN	3.2
1	C	128	ALA	3.2
1	A	316	SER	3.2
1	C	174	PHE	3.1
1	C	220	ASN	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	140	PRO	3.1
1	B	320	LEU	3.0
1	B	192	THR	3.0
1	C	195	THR	2.9
1	C	89	MET	2.9
1	A	232	THR	2.8
1	D	293	ILE	2.8
1	D	271	LEU	2.8
1	C	205	PRO	2.7
1	A	175	ASP	2.7
1	C	240	ILE	2.7
1	C	117	LEU	2.7
1	C	132	LEU	2.7
1	C	191	TRP	2.7
1	D	185	PRO	2.6
1	D	303	LEU	2.6
1	C	81	VAL	2.6
1	D	191	TRP	2.6
1	A	44	GLY	2.6
1	B	191	TRP	2.5
1	B	198	GLY	2.5
1	C	175	ASP	2.5
1	C	82	ILE	2.5
1	D	207	ILE	2.5
1	D	259	MET	2.5
1	B	321	PRO	2.5
1	A	141	GLY	2.5
1	C	281	GLY	2.5
1	D	138	LEU	2.5
1	C	233	SER	2.5
1	C	86	LYS	2.5
1	D	247	PHE	2.5
1	A	284	ILE	2.4
1	C	207	ILE	2.4
1	A	45	ASP	2.4
1	B	142	ILE	2.4
1	C	170	GLN	2.4
1	A	111	THR	2.4
1	B	173	THR	2.4
1	B	197	THR	2.4
1	D	88	LEU	2.4
1	A	209	PHE	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	279	LYS	2.3
1	D	236	LYS	2.3
1	B	200	ASN	2.3
1	A	250	VAL	2.3
1	A	272	THR	2.3
1	D	313	THR	2.3
1	D	267	LEU	2.3
1	D	149	LEU	2.3
1	C	151	ARG	2.3
1	C	293	ILE	2.2
1	B	276	VAL	2.2
1	A	136	ASN	2.2
1	C	255	THR	2.2
1	A	281	GLY	2.2
1	D	54	ILE	2.2
1	B	280	ASN	2.2
1	C	189	ALA	2.2
1	B	272	THR	2.2
1	C	172	PHE	2.2
1	A	314	CYS	2.1
1	D	57	ASP	2.1
1	C	146	PRO	2.1
1	D	211	PRO	2.1
1	A	114	ALA	2.1
1	A	132	LEU	2.1
1	C	158	LEU	2.1
1	C	313	THR	2.1
1	C	284	ILE	2.1
1	A	140	PRO	2.1
1	C	282	GLN	2.1
1	D	209	PHE	2.1
1	A	303	LEU	2.1
1	A	152	ILE	2.1
1	D	184	GLN	2.0
1	A	51	LEU	2.0
1	A	129	THR	2.0
1	C	261	ILE	2.0
1	D	263	VAL	2.0
1	D	128	ALA	2.0
1	A	122	ILE	2.0
1	A	277	THR	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
2	1PE	D	401	16/16	0.84	0.16	114,117,119,119	0
2	1PE	A	401	16/16	0.85	0.14	102,103,105,105	0
2	1PE	B	401	16/16	0.95	0.09	67,71,73,74	16

6.5 Other polymers [i](#)

There are no such residues in this entry.