



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:56 AM GMT

PDB ID : 2JEA
Title : STRUCTURE OF A 9-SUBUNIT ARCHAEAL EXOSOME BOUND TO
RNA
Authors : Lorentzen, E.; Conti, E.
Deposited on : 2007-01-16
Resolution : 2.33 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

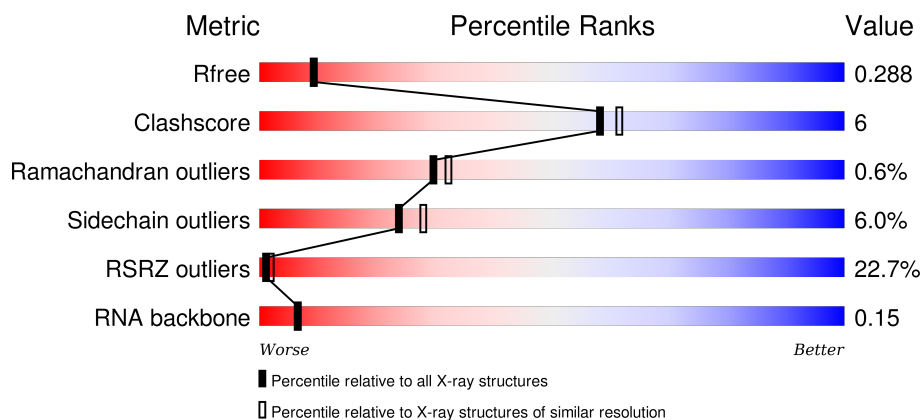
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.33 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}	91344	1406 (2.36-2.32)
Clashscore	102246	1509 (2.36-2.32)
Ramachandran outliers	100387	1490 (2.36-2.32)
Sidechain outliers	100360	1491 (2.36-2.32)
RSRZ outliers	91569	1412 (2.36-2.32)
RNA backbone	2183	1032 (2.88-1.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	277	<div> <div>15%</div> <div>85%</div> <div>11%</div> <div>• •</div> </div>
2	B	250	<div> <div>14%</div> <div>77%</div> <div>16%</div> <div>• 6%</div> </div>
3	C	35	<div> <div>17%</div> <div>9%</div> <div>6%</div> <div>6%</div> <div>80%</div> </div>
4	I	251	<div> <div>32%</div> <div>69%</div> <div>11%</div> <div>•</div> <div>18%</div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 5627 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 EXOSOME COMPLEX EXONUCLEASE 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	274	Total	C	N	O	S	0	9	0
			2091	1340	338	406	7			

- Molecule 2 is a protein called EXOSOME COMPLEX EXONUCLEASE 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	234	Total	C	N	O	S	0	7	0
			1809	1152	309	337	11			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	182	ALA	ASP	ENGINEERED MUTATION	UNP Q9UXC2

- Molecule 3 is a RNA chain called RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	7	Total	C	N	O	P	0	0	2
			112	50	25	32	5			

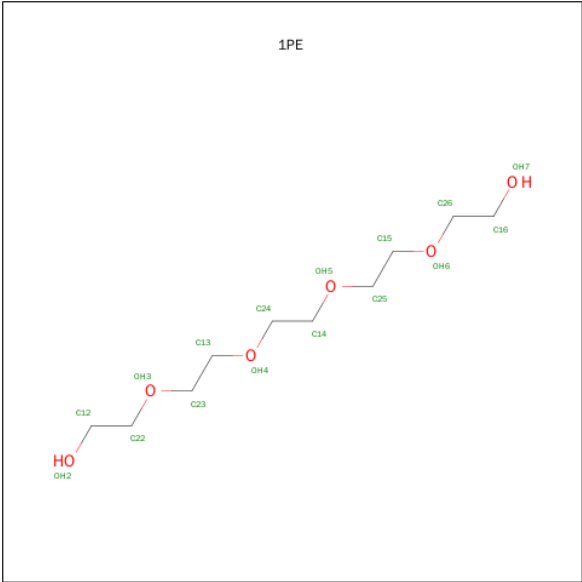
- Molecule 4 is a protein called EXOSOME COMPLEX RNA-BINDING PROTEIN 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	I	207	Total	C	N	O	S	0	0	0
			1477	963	238	274	2			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
I	8	GLU	LYS	CONFLICT	UNP Q9UXC4

- Molecule 5 is PENTAETHYLENE GLYCOL (three-letter code: 1PE) (formula: C₁₀H₂₂O₆).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	C	O	0
			16	10	6	

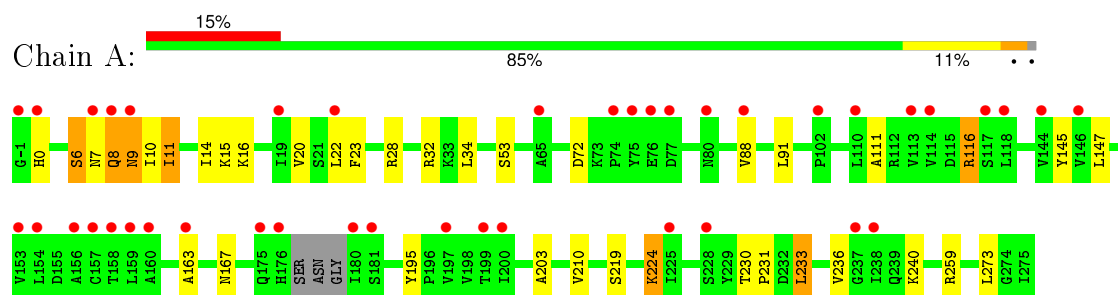
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	62	Total	O	0	0
			62	62		
6	B	51	Total	O	0	0
			51	51		
6	C	3	Total	O	0	0
			3	3		
6	I	6	Total	O	0	0
			6	6		

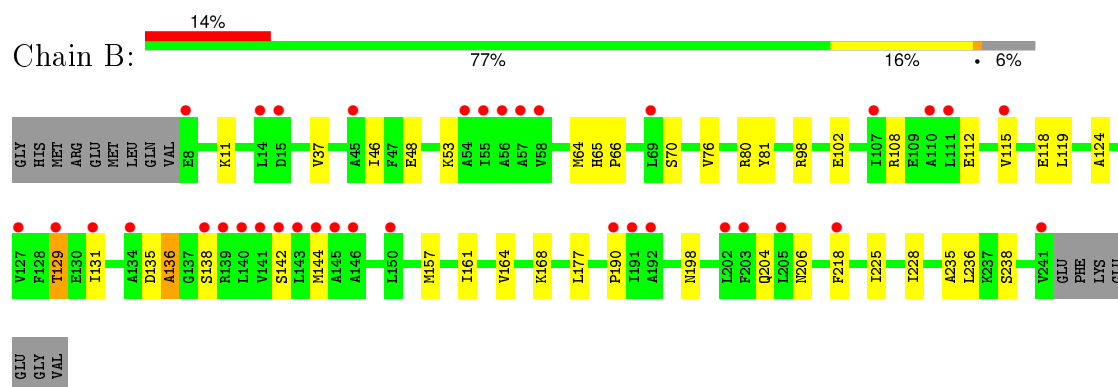
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

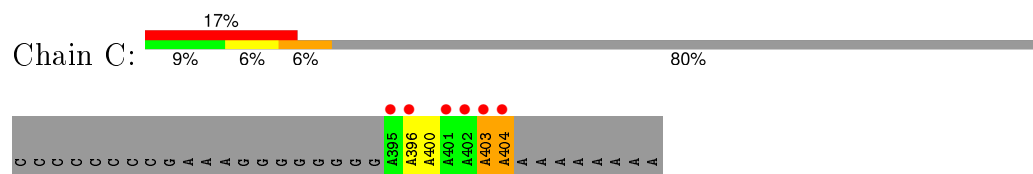
• Molecule 1: EXOSOME COMPLEX EXONUCLEASE 2



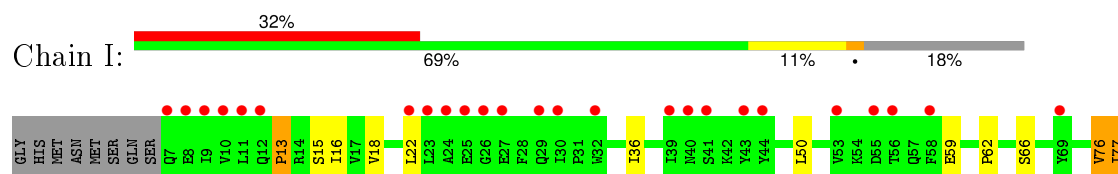
• Molecule 2: EXOSOME COMPLEX EXONUCLEASE 1

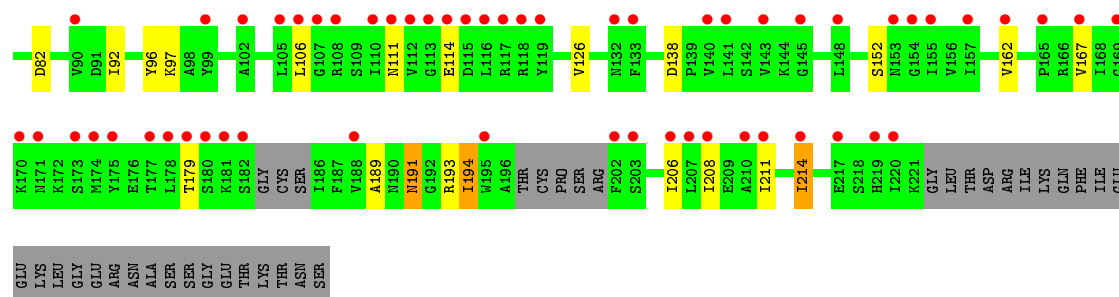


• Molecule 3: RNA



• Molecule 4: EXOSOME COMPLEX RNA-BINDING PROTEIN 1





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants a, b, c, α , β , γ	135.85Å 135.85Å 135.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.95 – 2.33 36.31 – 2.33	Depositor EDS
% Data completeness (in resolution range)	100.0 (42.95-2.33) 99.7 (36.31-2.33)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.46 (at 2.34Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.205 , 0.243 0.273 , 0.288	Depositor DCC
R_{free} test set	1788 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	43.4	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 61.8	EDS
Estimated twinning fraction	0.038 for l,-k,h	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 35804 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	5627	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.45% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: 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	0.61	1/2147 (0.0%)	0.77	2/2922 (0.1%)
2	B	0.60	0/1858	0.76	0/2514
3	C	1.22	1/125 (0.8%)	1.58	2/193 (1.0%)
4	I	0.49	0/1505	0.64	0/2064
All	All	0.60	2/5635 (0.0%)	0.76	4/7693 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	400	A	O3'-P	-10.52	1.48	1.61
1	A	224	LYS	CE-NZ	5.47	1.62	1.49

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	116	ARG	NE-CZ-NH2	-10.72	114.94	120.30
1	A	116	ARG	NE-CZ-NH1	7.54	124.07	120.30
3	C	403	A	P-O3'-C3'	-7.43	110.78	119.70
3	C	396	A	O4'-C1'-N9	5.43	112.54	108.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	6	SER	Peptide

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	2091	0	2128	21	0
2	B	1809	0	1859	26	0
3	C	112	0	57	2	0
4	I	1477	0	1373	18	0
5	A	16	0	22	0	0
6	A	62	0	0	1	0
6	B	51	0	0	4	0
6	C	3	0	0	0	0
6	I	6	0	0	0	0
All	All	5627	0	5439	65	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 (65) 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)
4:I:167:VAL:HG22	4:I:214:ILE:HD11	1.49	0.92
1:A:53:SER:H	1:A:167:ASN:HD22	1.27	0.82
2:B:81:TYR:HA	2:B:129:THR:HG22	1.65	0.78
4:I:208:ILE:HA	4:I:211:ILE:HG22	1.67	0.74
1:A:14:ILE:H	1:A:14:ILE:HD12	1.54	0.71
2:B:206:ASN:ND2	6:B:2039:HOH:O	2.27	0.68
2:B:129:THR:HG21	2:B:142:SER:OG	1.94	0.68
4:I:76:VAL:HG22	4:I:191:ASN:CG	2.13	0.68
1:A:8:GLN:O	1:A:10:ILE:N	2.30	0.64
2:B:235:ALA:O	2:B:238:SER:OG	2.15	0.63
2:B:118:GLU:HG3	6:B:2017:HOH:O	1.99	0.61
1:A:91:LEU:HD12	1:A:147:LEU:HD23	1.85	0.59
2:B:198:ASN:O	2:B:198:ASN:ND2	2.36	0.59
1:A:116:ARG:NH2	2:B:102:GLU:OE1	2.35	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:I:206:ILE:O	4:I:206:ILE:HG23	2.04	0.58
4:I:13:PRO:HD2	4:I:15:SER:OG	2.04	0.58
2:B:168:LYS:NZ	6:B:2028:HOH:O	2.27	0.57
1:A:11:ILE:HG21	1:A:219[A]:SER:OG	2.07	0.53
1:A:91:LEU:HD11	1:A:145:TYR:HD2	1.74	0.53
4:I:16:ILE:HD13	4:I:50:LEU:HD12	1.92	0.51
2:B:164[A]:VAL:CG1	2:B:228:ILE:CD1	2.88	0.51
2:B:198:ASN:CG	2:B:198:ASN:O	2.49	0.51
2:B:135:ASP:O	2:B:136:ALA:HB3	2.09	0.51
1:A:116:ARG:HD2	6:B:2037:HOH:O	2.11	0.50
4:I:167:VAL:CG2	4:I:214:ILE:HD11	2.33	0.50
1:A:240:LYS:NZ	6:A:2055:HOH:O	2.42	0.50
2:B:64:MET:HE1	2:B:76:VAL:CG2	2.42	0.50
2:B:37:VAL:HG11	2:B:46:ILE:HD11	1.93	0.49
1:A:28:ARG:CZ	1:A:210[B]:VAL:HG22	2.43	0.49
1:A:20:VAL:HG13	1:A:23:PHE:CD2	2.49	0.48
4:I:194:ILE:O	4:I:194:ILE:HG13	2.13	0.47
4:I:16:ILE:HD13	4:I:50:LEU:CD1	2.45	0.47
4:I:62:PRO:HG2	4:I:162:VAL:HG22	1.96	0.47
1:A:28:ARG:CZ	1:A:210[A]:VAL:HG13	2.44	0.47
1:A:28:ARG:HD2	1:A:34:LEU:HA	1.97	0.47
2:B:108:ARG:O	2:B:112:GLU:HG2	2.15	0.46
2:B:48:GLU:HG2	2:B:53:LYS:HG3	1.98	0.46
1:A:116:ARG:HH22	2:B:102:GLU:CD	2.17	0.46
4:I:36:ILE:HD11	4:I:59:GLU:HA	1.97	0.46
4:I:96:TYR:HB3	4:I:138:ASP:OD1	2.17	0.45
2:B:64:MET:CE	2:B:70:SER:HA	2.46	0.45
2:B:135:ASP:O	2:B:136:ALA:CB	2.65	0.44
4:I:62:PRO:O	4:I:162:VAL:HG13	2.18	0.44
2:B:115:VAL:HA	2:B:157:MET:HG2	2.00	0.44
1:A:163:ALA:HA	1:A:273:LEU:HD21	2.00	0.44
4:I:191:ASN:HA	4:I:191:ASN:HD22	1.57	0.44
2:B:65:HIS:N	2:B:66:PRO:CD	2.81	0.44
2:B:131:ILE:HD12	2:B:138:SER:HB2	2.00	0.43
1:A:230:THR:HB	1:A:231:PRO:CD	2.49	0.43
2:B:164[A]:VAL:HG13	2:B:228:ILE:CD1	2.49	0.42
3:C:404:A:H5'	3:C:404:A:C8	2.54	0.42
2:B:190:PRO:HG2	2:B:204:GLN:HB2	2.02	0.42
4:I:189:ALA:C	4:I:191:ASN:H	2.21	0.42
3:C:404:A:C8	3:C:404:A:C5'	3.02	0.42
4:I:77:ILE:HD11	4:I:193:ARG:HB3	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:144:MET:HG2	2:B:161:ILE:O	2.21	0.41
1:A:7:ASN:O	1:A:9:ASN:N	2.54	0.41
1:A:88:VAL:HG23	1:A:111:ALA:HA	2.02	0.41
2:B:164[A]:VAL:HG22	2:B:225:ILE:HG13	2.03	0.41
4:I:92:ILE:O	4:I:191:ASN:ND2	2.51	0.41
1:A:203:ALA:HB3	1:A:210[B]:VAL:HG12	2.03	0.41
4:I:111:ASN:H	4:I:114:GLU:HG3	1.86	0.40
1:A:11:ILE:HD11	1:A:16:LYS:HG3	2.03	0.40
2:B:76:VAL:O	2:B:124:ALA:HA	2.22	0.40
1:A:233:LEU:HG	1:A:259:ARG:HG3	2.03	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	279/277 (101%)	269 (96%)	8 (3%)	2 (1%)	26	28
2	B	239/250 (96%)	232 (97%)	6 (2%)	1 (0%)	39	45
4	I	201/251 (80%)	183 (91%)	17 (8%)	1 (0%)	34	37
All	All	719/778 (92%)	684 (95%)	31 (4%)	4 (1%)	30	32

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	8	GLN
1	A	9	ASN
4	I	13	PRO
2	B	136	ALA

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	231/243 (95%)	220 (95%)	11 (5%)	31	39
2	B	193/208 (93%)	184 (95%)	9 (5%)	32	40
4	I	140/223 (63%)	126 (90%)	14 (10%)	9	8
All	All	564/674 (84%)	530 (94%)	34 (6%)	24	28

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	0	HIS
1	A	6	SER
1	A	11	ILE
1	A	15	LYS
1	A	22	LEU
1	A	32	ARG
1	A	72	ASP
1	A	195	TYR
1	A	224	LYS
1	A	233	LEU
1	A	236	VAL
2	B	11	LYS
2	B	80	ARG
2	B	98	ARG
2	B	119	LEU
2	B	129	THR
2	B	177	LEU
2	B	218[A]	PHE
2	B	218[B]	PHE
2	B	236	LEU
4	I	18	VAL
4	I	22	LEU
4	I	66	SER
4	I	76	VAL
4	I	77	ILE
4	I	82	ASP

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Mol	Chain	Res	Type
4	I	97	LYS
4	I	106	LEU
4	I	126	VAL
4	I	152	SER
4	I	179	THR
4	I	191	ASN
4	I	194	ILE
4	I	214	ILE

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

Mol	Chain	Res	Type
1	A	167	ASN
2	B	198	ASN
4	I	191	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
3	C	3/35 (8%)	2 (66%)	0

All (2) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
3	C	403	A
3	C	404	A

There are no RNA pucker outliers to report.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

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	$\# Z > 2$
5	1PE	A	1276	-	15,15,15	0.56	0	14,14,14	0.25	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
5	1PE	A	1276	-	-	0/13/13/13	0/0/0/0

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	274/277 (98%)	0.90	41 (14%) 3 5	47, 54, 68, 73	1 (0%)
2	B	234/250 (93%)	0.91	36 (15%) 3 5	47, 54, 63, 72	0
3	C	7/35 (20%)	7.92	6 (85%) 0 0	31, 36, 41, 43	7 (100%)
4	I	207/251 (82%)	1.79	81 (39%) 0 0	46, 53, 60, 62	0
All	All	722/813 (88%)	1.23	164 (22%) 1 2	31, 53, 65, 73	8 (1%)

All (164) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	404	A	15.5
3	C	403	A	13.1
4	I	113	GLY	9.1
3	C	395	A	8.6
3	C	402	A	8.2
1	A	-1	GLY	7.6
4	I	9	ILE	6.2
3	C	401	A	6.1
1	A	75	TYR	5.7
4	I	112	VAL	5.7
4	I	8	GLU	5.2
4	I	110	ILE	5.1
4	I	26	GLY	5.1
4	I	106	LEU	5.1
4	I	169	GLY	5.1
4	I	41	SER	4.8
1	A	9	ASN	4.8
4	I	170	LYS	4.8
4	I	220	ILE	4.7
4	I	208	ILE	4.7
4	I	203	SER	4.6

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Mol	Chain	Res	Type	RSRZ
4	I	180	SER	4.6
4	I	107	GLY	4.6
4	I	105	LEU	4.6
4	I	58	PHE	4.5
4	I	119	TYR	4.5
1	A	176	HIS	4.2
4	I	24	ALA	4.1
4	I	182	SER	4.1
4	I	10	VAL	4.0
4	I	206	ILE	4.0
4	I	211	ILE	4.0
2	B	143	LEU	3.9
4	I	56	THR	3.9
1	A	114	VAL	3.9
4	I	178	LEU	3.9
1	A	102	PRO	3.8
4	I	102	ALA	3.8
1	A	157	CYS	3.8
4	I	171	ASN	3.8
4	I	174	MET	3.7
1	A	154	LEU	3.7
4	I	115	ASP	3.7
4	I	165	PRO	3.7
4	I	27	GLU	3.6
4	I	179	THR	3.6
2	B	140	LEU	3.6
4	I	40	ASN	3.5
4	I	43	TYR	3.5
2	B	107	ILE	3.5
4	I	207	LEU	3.5
2	B	110	ALA	3.5
4	I	181	LYS	3.5
1	A	19	ILE	3.5
1	A	238	ILE	3.4
2	B	241	VAL	3.4
2	B	8	GLU	3.4
4	I	11	LEU	3.4
4	I	202	PHE	3.4
4	I	148	LEU	3.3
4	I	32	TRP	3.3
4	I	219	HIS	3.3
4	I	108	ARG	3.2

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Mol	Chain	Res	Type	RSRZ
4	I	153	ASN	3.2
2	B	55	ILE	3.2
4	I	210	ALA	3.2
2	B	129	THR	3.2
4	I	30	ILE	3.1
4	I	55	ASP	3.1
2	B	54	ALA	3.1
2	B	131	ILE	3.1
2	B	192	ALA	3.1
2	B	15	ASP	3.1
2	B	203	PHE	3.1
4	I	53	VAL	3.0
4	I	175	TYR	3.0
4	I	141	LEU	3.0
4	I	29	GLN	3.0
1	A	180	ILE	3.0
4	I	25	GLU	3.0
2	B	45	ALA	3.0
1	A	8	GLN	3.0
4	I	195	TRP	2.9
2	B	138	SER	2.9
4	I	39	ILE	2.9
1	A	76	GLU	2.9
2	B	141	VAL	2.9
2	B	202	LEU	2.9
4	I	23	LEU	2.9
4	I	22	LEU	2.9
1	A	199	THR	2.9
4	I	214	ILE	2.8
4	I	111	ASN	2.8
4	I	167	VAL	2.8
4	I	114	GLU	2.8
4	I	177	THR	2.8
1	A	237	GLY	2.7
4	I	145	GLY	2.7
1	A	113	VAL	2.7
2	B	69	LEU	2.7
1	A	88	VAL	2.7
4	I	90	VAL	2.7
2	B	146	ALA	2.7
1	A	22	LEU	2.7
2	B	111	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
2	B	142	SER	2.7
1	A	175	GLN	2.6
2	B	190	PRO	2.6
2	B	115	VAL	2.6
1	A	200	ILE	2.6
2	B	14	LEU	2.6
1	A	77	ASP	2.6
4	I	162	VAL	2.6
1	A	225	ILE	2.6
1	A	181	SER	2.6
4	I	132	ASN	2.6
4	I	99	TYR	2.5
4	I	116	LEU	2.5
1	A	197	VAL	2.5
2	B	191	ILE	2.5
2	B	218[A]	PHE	2.5
3	C	396	A	2.5
4	I	154	GLY	2.5
4	I	155	ILE	2.5
1	A	110	LEU	2.5
1	A	144	VAL	2.5
4	I	140	VAL	2.5
4	I	69	TYR	2.4
4	I	143	VAL	2.4
1	A	117	SER	2.4
2	B	145	ALA	2.4
2	B	205[A]	LEU	2.4
4	I	133	PHE	2.4
1	A	0	HIS	2.4
1	A	80	ASN	2.4
4	I	7	GLN	2.4
1	A	153[A]	VAL	2.3
1	A	65	ALA	2.3
2	B	56	ALA	2.3
2	B	58	VAL	2.3
4	I	44	TYR	2.3
1	A	159	LEU	2.3
2	B	57	ALA	2.3
4	I	217	GLU	2.2
4	I	118	ARG	2.2
2	B	150	LEU	2.2
1	A	156	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	146	VAL	2.2
4	I	12	GLN	2.2
1	A	118	LEU	2.2
1	A	158	THR	2.2
2	B	134	ALA	2.2
2	B	144	MET	2.2
4	I	157	ILE	2.1
1	A	228	SER	2.1
1	A	7	ASN	2.1
4	I	173	SER	2.1
4	I	188	VAL	2.1
1	A	163	ALA	2.1
1	A	74	PRO	2.1
2	B	127	VAL	2.1
2	B	139	ARG	2.1
4	I	117	ARG	2.1
1	A	160	ALA	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	1PE	A	1276	16/16	0.85	0.18	0.58	56,62,77,77	0

6.5 Other polymers

There are no such residues in this entry.