



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:13 PM UTC

PDB ID : 6S8S / pdb_00006s8s
Title : Extended structure of the human DDX6 C-terminal domain in complex with an EDC3 FDF peptide
Authors : Peter, D.; Valkov, E.
Deposited on : 2019-07-10
Resolution : 2.21 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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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

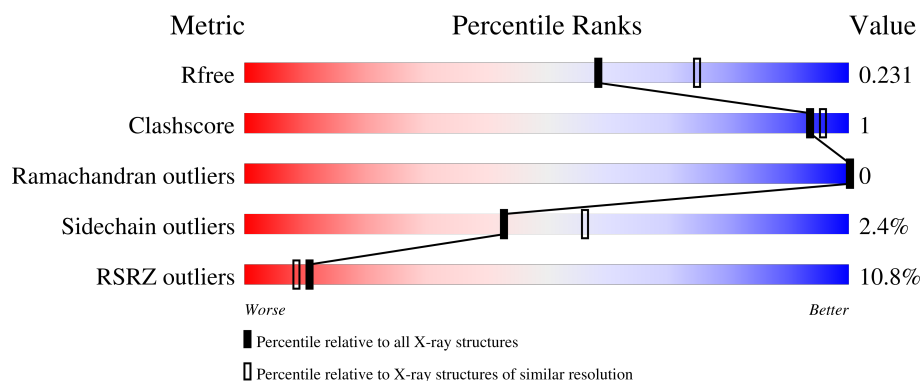
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.21 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	7682 (2.24-2.20)
Clashscore	190562	8402 (2.24-2.20)
Ramachandran outliers	187476	8303 (2.24-2.20)
Sidechain outliers	187428	8304 (2.24-2.20)
RSRZ outliers	180081	7683 (2.24-2.20)

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 $\leq 5\%$. 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	193	5% 82% 13%
1	C	193	7% 84% 10%
2	B	44	25% 82% 16%
2	D	44	25% 73% 5% 5% 16%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3775 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 Probable ATP-dependent RNA helicase DDX6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	168	Total	C	N	O	S	0	0	0
			1381	880	250	245	6			
1	C	173	Total	C	N	O	S	0	0	0
			1419	903	255	255	6			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	280	GLY	-	expression tag	UNP P26196
A	281	PRO	-	expression tag	UNP P26196
A	282	GLN	-	expression tag	UNP P26196
A	283	ASP	-	expression tag	UNP P26196
C	280	GLY	-	expression tag	UNP P26196
C	281	PRO	-	expression tag	UNP P26196
C	282	GLN	-	expression tag	UNP P26196
C	283	ASP	-	expression tag	UNP P26196

- Molecule 2 is a protein called Enhancer of mRNA-decapping protein 3.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	37	Total	C	N	O	0	0	0
			301	193	39	69			
2	D	37	Total	C	N	O	0	0	0
			301	193	39	69			

There are 14 discrepancies between the modelled and reference sequences:

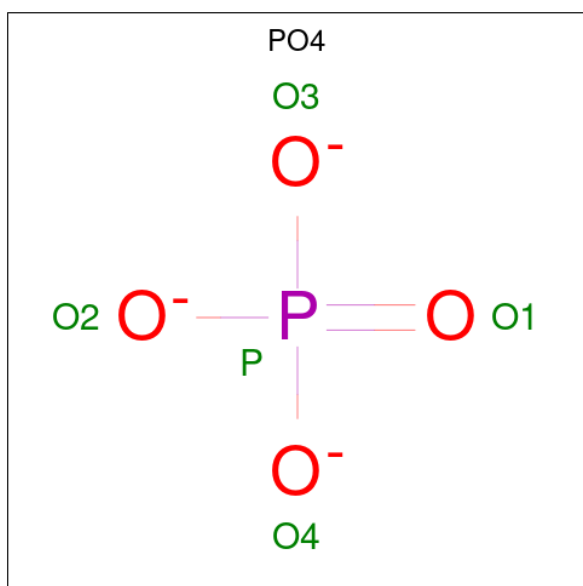
Chain	Residue	Modelled	Actual	Comment	Reference
B	185	GLY	-	expression tag	UNP Q96F86
B	186	PRO	-	expression tag	UNP Q96F86
B	187	HIS	-	expression tag	UNP Q96F86
B	188	MET	-	expression tag	UNP Q96F86

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Chain	Residue	Modelled	Actual	Comment	Reference
B	189	ALA	-	expression tag	UNP Q96F86
B	190	ASP	-	expression tag	UNP Q96F86
B	191	LEU	-	expression tag	UNP Q96F86
D	185	GLY	-	expression tag	UNP Q96F86
D	186	PRO	-	expression tag	UNP Q96F86
D	187	HIS	-	expression tag	UNP Q96F86
D	188	MET	-	expression tag	UNP Q96F86
D	189	ALA	-	expression tag	UNP Q96F86
D	190	ASP	-	expression tag	UNP Q96F86
D	191	LEU	-	expression tag	UNP Q96F86

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



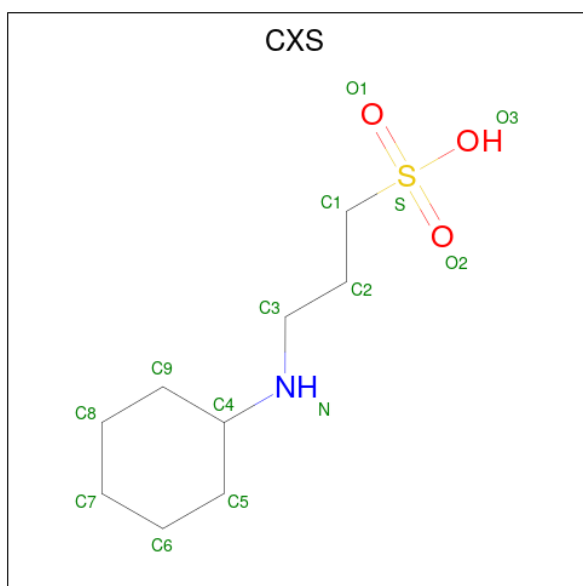
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O P 5 4 1	0	0
3	A	1	Total O P 5 4 1	0	0
3	A	1	Total O P 5 4 1	0	0
3	A	1	Total O P 5 4 1	0	0
3	A	1	Total O P 5 4 1	0	0
3	C	1	Total O P 5 4 1	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	C	1	Total	O	P	0	0
			5	4	1		
3	C	1	Total	O	P	0	0
			5	4	1		
3	D	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is 3-CYCLOHEXYL-1-PROPYLSULFONIC ACID (CCD ID: CXS) (formula: $C_9H_{19}NO_3S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			14	9	1	3	1		
4	A	1	Total	C	N	O	S	0	0
			14	9	1	3	1		
4	A	1	Total	C	N	O	S	0	0
			14	9	1	3	1		
4	C	1	Total	C	N	O	S	0	0
			14	9	1	3	1		
4	C	1	Total	C	N	O	S	0	0
			14	9	1	3	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	98	Total	O	0	0
			98	98		

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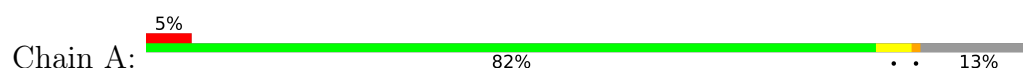
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	31	Total 31	O 31	0	0
5	C	99	Total 99	O 99	0	0
5	D	30	Total 30	O 30	0	0

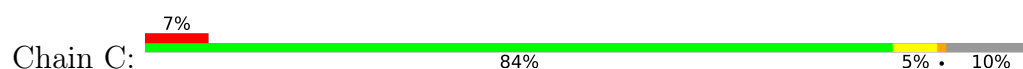
3 Residue-property plots [i](#)

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.

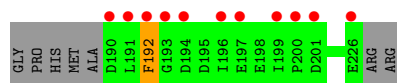
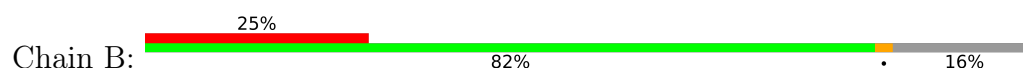
- Molecule 1: Probable ATP-dependent RNA helicase DDX6



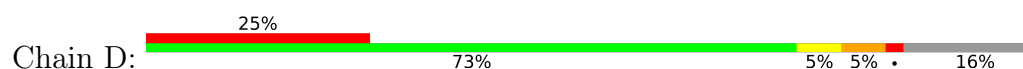
- Molecule 1: Probable ATP-dependent RNA helicase DDX6



- Molecule 2: Enhancer of mRNA-decapping protein 3



- Molecule 2: Enhancer of mRNA-decapping protein 3



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	172.43Å 47.88Å 65.76Å 90.00° 96.34° 90.00°	Depositor
Resolution (Å)	25.31 – 2.21 25.31 – 2.21	Depositor EDS
% Data completeness (in resolution range)	95.6 (25.31-2.21) 95.6 (25.31-2.21)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.62 (at 2.20Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.179 , 0.226 0.187 , 0.231	Depositor DCC
R_{free} test set	1315 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å ²)	25.8	Xtriage
Anisotropy	0.736	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 61.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3775	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.41% 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: PO4, CXS

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	0.89	1/1409 (0.1%)	1.25	6/1899 (0.3%)
1	C	0.87	1/1448 (0.1%)	1.22	2/1953 (0.1%)
2	B	0.86	0/307	1.29	1/416 (0.2%)
2	D	0.82	0/307	1.33	1/416 (0.2%)
All	All	0.87	2/3471 (0.1%)	1.25	10/4684 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	357	MET	SD-CE	-6.45	1.63	1.79
1	C	357	MET	SD-CE	-5.71	1.65	1.79

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	192	PHE	CA-CB-CG	7.55	121.35	113.80
2	B	192	PHE	CA-CB-CG	6.62	120.42	113.80
1	A	324	ASN	N-CA-C	-5.30	99.50	110.80
1	C	324	ASN	N-CA-C	-5.28	99.56	110.80
1	A	389	ILE	CA-C-N	5.27	127.60	120.38
1	A	389	ILE	C-N-CA	5.27	127.60	120.38
1	A	390	GLN	N-CA-C	5.22	117.65	111.33
1	A	462	TYR	CA-C-N	5.16	130.99	121.70
1	A	462	TYR	C-N-CA	5.16	130.99	121.70
1	C	306	THR	N-CA-C	-5.15	102.16	110.14

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	1381	0	1384	4	0
1	C	1419	0	1417	4	0
2	B	301	0	257	0	0
2	D	301	0	257	3	0
3	A	25	0	0	0	0
3	C	15	0	0	0	0
3	D	5	0	0	0	0
4	A	42	0	54	1	0
4	C	28	0	37	0	0
5	A	98	0	0	0	0
5	B	31	0	0	0	0
5	C	99	0	0	0	0
5	D	30	0	0	0	0
All	All	3775	0	3406	10	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 1.

All (10) 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:A:455:LYS:HB2	1:A:463:HIS:HB2	1.54	0.89
1:C:357:MET:HE2	1:C:362:ARG:HG3	1.74	0.69
2:D:195:ASP:HB2	2:D:198:GLU:HB2	1.83	0.61
2:D:196:ILE:HA	2:D:199:ILE:HD12	1.90	0.52
1:C:353:ILE:HG22	1:C:357:MET:HE1	1.94	0.49
1:C:308:ARG:HB2	1:C:308:ARG:CZ	2.45	0.46
1:A:353:ILE:HG22	1:A:357:MET:HE1	1.97	0.46
1:C:339:LEU:HB2	2:D:192:PHE:HZ	1.82	0.44
1:A:348:TYR:CD1	4:A:507:CXS:H71	2.56	0.40
1:A:459:VAL:HG12	1:A:462:TYR:HB2	2.04	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	166/193 (86%)	161 (97%)	5 (3%)	0	100	100
1	C	171/193 (89%)	167 (98%)	4 (2%)	0	100	100
2	B	35/44 (80%)	34 (97%)	1 (3%)	0	100	100
2	D	35/44 (80%)	34 (97%)	1 (3%)	0	100	100
All	All	407/474 (86%)	396 (97%)	11 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	152/176 (86%)	152 (100%)	0	100	100
1	C	157/176 (89%)	152 (97%)	5 (3%)	34	45
2	B	32/37 (86%)	31 (97%)	1 (3%)	35	46
2	D	32/37 (86%)	29 (91%)	3 (9%)	8	8
All	All	373/426 (88%)	364 (98%)	9 (2%)	43	56

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

Mol	Chain	Res	Type
2	B	192	PHE
1	C	308	ARG

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Mol	Chain	Res	Type
1	C	417	PHE
1	C	426	LEU
1	C	464	SER
1	C	467	VAL
2	D	192	PHE
2	D	196	ILE
2	D	198	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	300	GLN
1	A	315	ASN
1	A	376	ASN
1	A	409	HIS
1	A	463	HIS
1	C	300	GLN
1	C	315	ASN
1	C	359	GLN
1	C	376	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

14 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
3	PO4	A	501	-	4,4,4	2.54	3 (75%)	6,6,6	0.65	0
3	PO4	D	301	-	4,4,4	2.50	2 (50%)	6,6,6	0.89	0
3	PO4	A	505	-	4,4,4	2.57	2 (50%)	6,6,6	0.91	0
4	CXS	A	507	-	14,14,14	0.58	0	18,18,18	1.27	2 (11%)
4	CXS	C	504	-	14,14,14	0.57	0	18,18,18	0.78	0
3	PO4	C	502	-	4,4,4	2.57	1 (25%)	6,6,6	0.29	0
4	CXS	A	506	-	14,14,14	0.55	0	18,18,18	1.26	2 (11%)
3	PO4	C	503	-	4,4,4	2.60	2 (50%)	6,6,6	0.56	0
4	CXS	C	505	-	14,14,14	0.29	0	18,18,18	1.05	2 (11%)
4	CXS	A	508	-	14,14,14	0.28	0	18,18,18	0.90	1 (5%)
3	PO4	A	503	-	4,4,4	2.58	3 (75%)	6,6,6	0.55	0
3	PO4	A	502	-	4,4,4	2.57	2 (50%)	6,6,6	0.69	0
3	PO4	C	501	-	4,4,4	2.52	2 (50%)	6,6,6	0.72	0
3	PO4	A	504	-	4,4,4	2.00	1 (25%)	6,6,6	0.52	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	CXS	A	507	-	-	4/8/16/16	0/1/1/1
4	CXS	C	504	-	-	0/8/16/16	0/1/1/1
4	CXS	A	506	-	-	1/8/16/16	0/1/1/1
4	CXS	C	505	-	-	4/8/16/16	0/1/1/1
4	CXS	A	508	-	-	1/8/16/16	0/1/1/1

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	505	PO4	P-O1	4.32	1.60	1.50
3	C	503	PO4	P-O1	4.29	1.60	1.50
3	A	502	PO4	P-O1	4.28	1.60	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	503	PO4	P-O1	4.26	1.60	1.50
3	C	502	PO4	P-O1	4.25	1.60	1.50
3	A	501	PO4	P-O1	4.18	1.60	1.50
3	C	501	PO4	P-O1	4.13	1.60	1.50
3	D	301	PO4	P-O1	4.12	1.60	1.50
3	A	503	PO4	P-O2	2.05	1.60	1.54
3	A	505	PO4	P-O4	2.04	1.60	1.54
3	A	504	PO4	P-O1	2.04	1.55	1.50
3	C	503	PO4	P-O3	2.03	1.60	1.54
3	A	501	PO4	P-O2	2.03	1.60	1.54
3	D	301	PO4	P-O3	2.02	1.60	1.54
3	A	503	PO4	P-O4	2.02	1.60	1.54
3	C	501	PO4	P-O2	2.02	1.60	1.54
3	A	502	PO4	P-O2	2.02	1.60	1.54
3	A	501	PO4	P-O4	2.01	1.60	1.54

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	506	CXS	C3-N-C4	4.21	122.29	114.18
4	A	507	CXS	C9-C4-N	3.78	129.46	110.56
4	C	505	CXS	C9-C4-N	2.91	125.10	110.56
4	A	507	CXS	O2-S-C1	-2.51	102.94	106.73
4	A	508	CXS	C3-N-C4	2.50	119.00	114.18
4	A	506	CXS	C9-C4-N	2.25	121.79	110.56
4	C	505	CXS	C2-C1-S	-2.19	109.88	113.25

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	506	CXS	C9-C4-N-C3
4	A	507	CXS	C9-C4-N-C3
4	C	505	CXS	C9-C4-N-C3
4	C	505	CXS	C1-C2-C3-N
4	A	507	CXS	C2-C1-S-O2
4	C	505	CXS	C2-C3-N-C4
4	A	507	CXS	C2-C1-S-O3
4	A	508	CXS	C2-C3-N-C4
4	C	505	CXS	C5-C4-N-C3
4	A	507	CXS	S-C1-C2-C3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	507	CXS	1	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	168/193 (87%)	-0.05	10 (5%) 27 25	18, 31, 66, 118	0
1	C	173/193 (89%)	-0.10	13 (7%) 20 17	18, 28, 79, 103	0
2	B	37/44 (84%)	1.11	11 (29%) 1 1	21, 35, 106, 108	0
2	D	37/44 (84%)	0.94	11 (29%) 1 1	24, 35, 112, 114	0
All	All	415/474 (87%)	0.12	45 (10%) 11 8	18, 30, 89, 118	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	196	ILE	7.0
2	B	196	ILE	6.5
2	B	191	LEU	6.2
2	B	192	PHE	5.5
2	D	190	ASP	4.9
2	B	193	GLY	4.7
2	D	193	GLY	4.7
1	C	467	VAL	4.6
1	A	463	HIS	4.3
2	B	190	ASP	4.2
2	D	194	ASP	3.9
2	B	194	ASP	3.7
1	A	462	TYR	3.7
1	C	296	LYS	3.7
2	D	192	PHE	3.6
2	D	191	LEU	3.6
1	C	417	PHE	3.6
1	C	414	SER	3.4
1	C	465	GLU	3.4
2	B	226	GLU	3.2
2	D	197	GLU	3.0

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Mol	Chain	Res	Type	RSRZ
1	C	466	PRO	2.9
1	C	463	HIS	2.9
1	A	417	PHE	2.9
2	B	199	ILE	2.8
1	C	415	GLY	2.7
1	A	387	ILE	2.7
1	C	420	LEU	2.7
1	A	419	HIS	2.6
2	D	195	ASP	2.5
1	A	416	ARG	2.5
1	A	414	SER	2.4
2	B	201	ASP	2.4
1	C	416	ARG	2.4
2	D	199	ILE	2.3
2	D	198	GLU	2.3
1	A	296	LYS	2.3
1	A	388	ASP	2.2
1	C	468	GLU	2.2
2	D	226	GLU	2.2
1	A	415	GLY	2.2
1	C	419	HIS	2.2
2	B	197	GLU	2.1
1	C	464	SER	2.1
2	B	200	PRO	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
4	CXS	A	508	14/14	0.69	0.26	78,88,109,109	0
3	PO4	A	503	5/5	0.80	0.16	106,107,107,108	0
3	PO4	A	502	5/5	0.80	0.16	80,81,83,85	0
3	PO4	A	501	5/5	0.82	0.14	90,90,91,91	0
4	CXS	C	505	14/14	0.84	0.13	48,60,86,87	0
4	CXS	A	507	14/14	0.86	0.13	43,53,84,85	0
3	PO4	C	501	5/5	0.87	0.13	76,77,79,79	0
3	PO4	A	505	5/5	0.87	0.19	66,67,69,70	0
3	PO4	C	503	5/5	0.88	0.13	74,75,77,77	0
4	CXS	C	504	14/14	0.88	0.12	44,57,83,83	0
3	PO4	A	504	5/5	0.88	0.13	82,82,83,83	0
3	PO4	C	502	5/5	0.90	0.14	66,67,67,71	0
3	PO4	D	301	5/5	0.95	0.11	49,55,59,62	0
4	CXS	A	506	14/14	0.97	0.07	26,30,34,36	0

6.5 Other polymers [i](#)

There are no such residues in this entry.