



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 02:50 PM UTC

PDB ID : 4H9S / pdb_00004h9s
Title : Complex structure 6 of DAXX/H3.3(sub7)/H4
Authors : Elsasser, S.J.; Huang, H.; Lewis, P.W.; Chin, J.W.; Allis, D.C.; Patel, D.J.
Deposited on : 2012-09-24
Resolution : 2.60 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

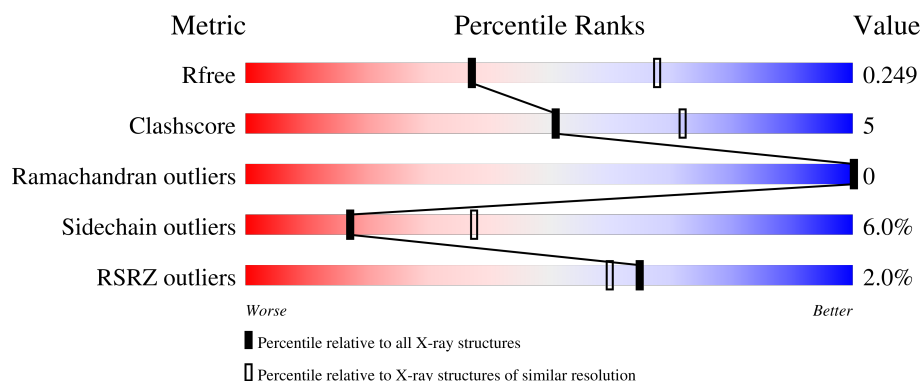
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	135	53% 13% 34%
1	B	135	59% 10% 30%
2	C	83	82% 7% 11%
2	D	83	77% 12% 11%
3	E	216	79% 12% 8%

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Mol	Chain	Length	Quality of chain
3	F	216	% 71% 19% • 6%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	PO4	C	201	-	-	X	-
4	PO4	D	201	-	-	X	-
4	PO4	F	402	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6076 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 Histone H3.3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	89	Total	C	N	O	S	0	0	0
			720	458	135	125	2			
1	B	94	Total	C	N	O	S	0	0	0
			744	474	138	130	2			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	75	CYS	ALA	conflict	UNP P84243
A	84	TRP	PHE	conflict	UNP P84243
A	96	ALA	SER	conflict	UNP P84243
A	99	PHE	TYR	conflict	UNP P84243
A	102	ALA	GLY	conflict	UNP P84243
A	111	THR	ALA	conflict	UNP P84243
A	120	PHE	MET	conflict	UNP P84243
B	75	CYS	ALA	conflict	UNP P84243
B	84	TRP	PHE	conflict	UNP P84243
B	96	ALA	SER	conflict	UNP P84243
B	99	PHE	TYR	conflict	UNP P84243
B	102	ALA	GLY	conflict	UNP P84243
B	111	THR	ALA	conflict	UNP P84243
B	120	PHE	MET	conflict	UNP P84243

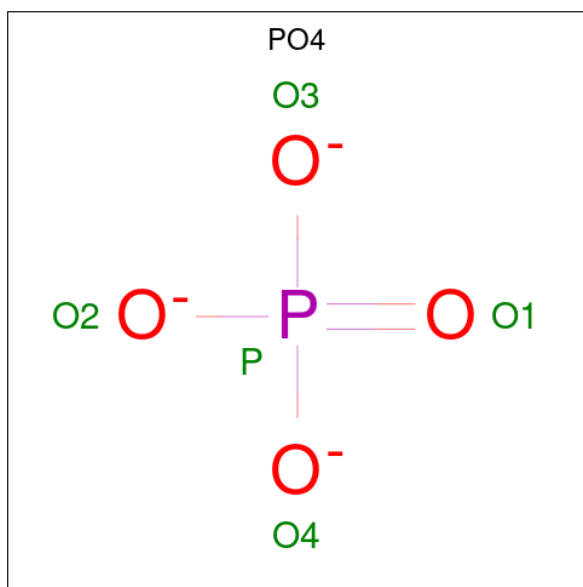
- Molecule 2 is a protein called Histone H4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	74	Total	C	N	O	S	0	0	0
			577	366	111	99	1			
2	D	74	Total	C	N	O	S	0	0	0
			577	366	111	99	1			

- Molecule 3 is a protein called Death domain-associated protein 6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	199	Total	C	N	O	S	0	0	0
			1611	1004	307	294	6			
3	F	203	Total	C	N	O	S	0	0	0
			1650	1024	313	307	6			

- Molecule 4 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	20	Total	O	0	0
			20	20		
5	B	24	Total	O	0	0
			24	24		
5	C	18	Total	O	0	0
			18	18		

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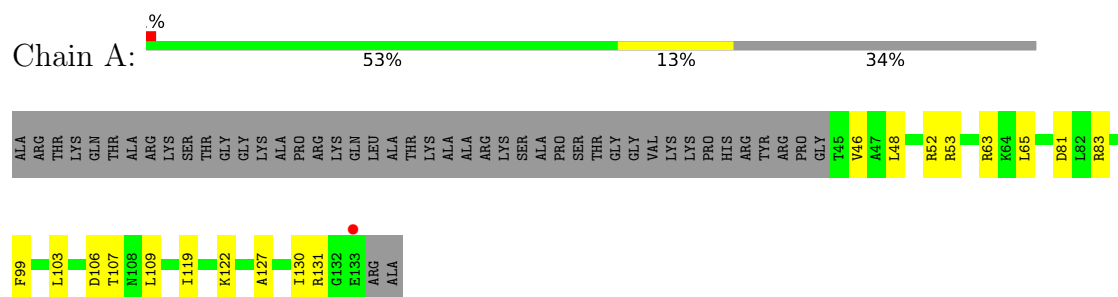
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	D	20	Total	O	0	0
			20	20		
5	E	45	Total	O	0	0
			45	45		
5	F	45	Total	O	0	0
			45	45		

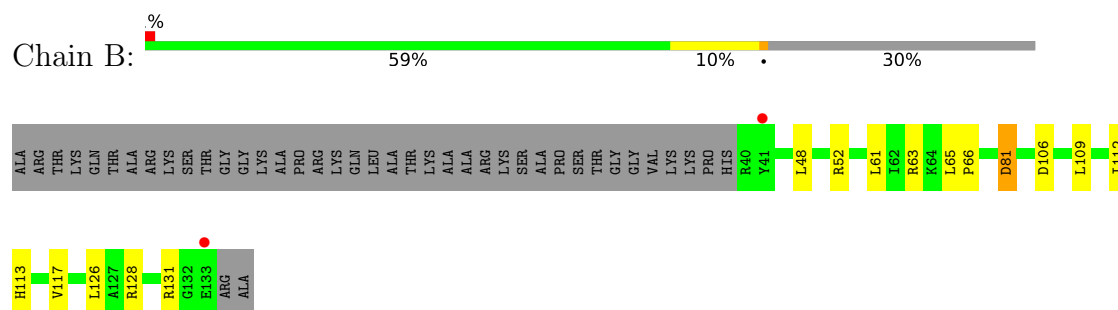
3 Residue-property plots [i](#)

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.

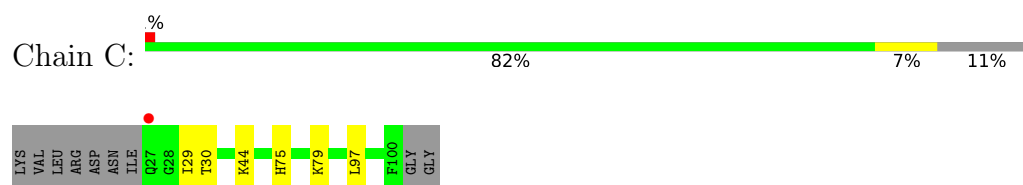
- Molecule 1: Histone H3.3



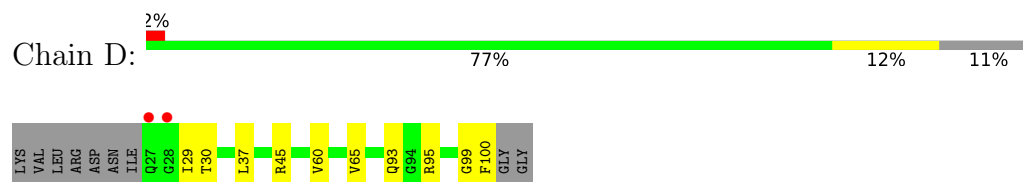
- Molecule 1: Histone H3.3



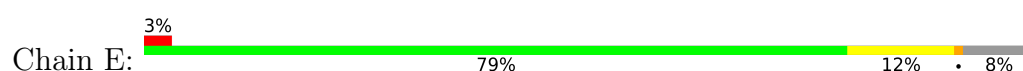
- Molecule 2: Histone H4

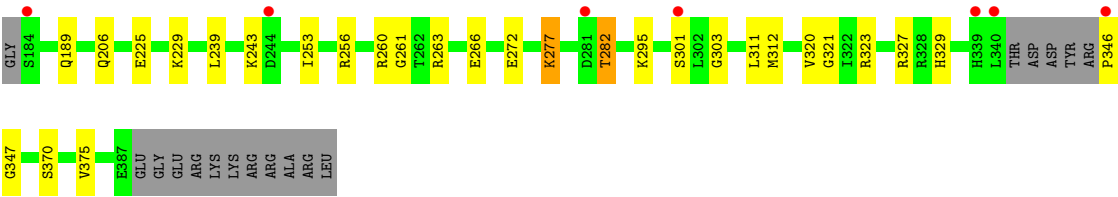


- Molecule 2: Histone H4

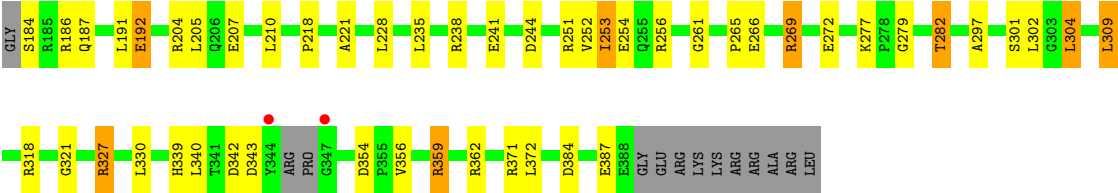


- Molecule 3: Death domain-associated protein 6





● Molecule 3: Death domain-associated protein 6



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	96.40Å 99.17Å 100.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.09 – 2.60 44.09 – 2.60	Depositor EDS
% Data completeness (in resolution range)	100.0 (44.09-2.60) 96.5 (44.09-2.60)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.39 (at 2.29Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.2_869)	Depositor
R, R_{free}	0.195 , 0.250 0.194 , 0.249	Depositor DCC
R_{free} test set	2209 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	41.3	Xtriage
Anisotropy	0.103	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 42.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.065 for -h,l,k 0.024 for -l,-k,-h 0.026 for k,h,-l 0.009 for k,l,h 0.009 for l,h,k	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6076	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.35% 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: PO4

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	0.39	0/730	0.80	1/985 (0.1%)
1	B	0.37	0/755	0.77	0/1020
2	C	0.38	0/583	0.76	0/782
2	D	0.39	0/583	0.77	0/782
3	E	0.39	0/1632	0.82	3/2194 (0.1%)
3	F	0.38	0/1671	0.84	2/2247 (0.1%)
All	All	0.38	0/5954	0.81	6/8010 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	279	GLY	CA-C-N	8.49	128.19	119.19
3	F	279	GLY	C-N-CA	8.49	128.19	119.19
1	A	130	ILE	N-CA-C	5.57	115.81	110.74
3	E	277	LYS	CA-C-N	5.23	125.15	119.76
3	E	277	LYS	C-N-CA	5.23	125.15	119.76
3	E	303	GLY	N-CA-C	-5.12	101.03	113.18

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	720	0	741	16	0
1	B	744	0	757	10	0
2	C	577	0	612	4	0
2	D	577	0	612	7	0
3	E	1611	0	1645	14	0
3	F	1650	0	1669	28	0
4	C	5	0	0	2	0
4	D	5	0	0	2	0
4	E	5	0	0	1	0
4	F	10	0	0	3	0
5	A	20	0	0	0	0
5	B	24	0	0	1	0
5	C	18	0	0	0	0
5	D	20	0	0	0	0
5	E	45	0	0	1	0
5	F	45	0	0	1	0
All	All	6076	0	6036	62	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 5.

All (62) 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:52:ARG:NH2	3:E:261:GLY:O	2.20	0.74
1:A:106:ASP:OD1	1:A:131:ARG:NH1	2.28	0.66
1:B:109:LEU:HD21	3:E:321:GLY:HA3	1.78	0.65
1:B:52:ARG:NH2	3:F:261:GLY:O	2.29	0.65
3:F:186:ARG:NH2	3:F:187:GLN:OE1	2.31	0.63
2:C:75:HIS:NE2	3:F:354:ASP:OD2	2.29	0.63
1:B:106:ASP:OD1	1:B:131:ARG:NH1	2.32	0.62
3:F:252:VAL:N	4:F:402:PO4:O1	2.33	0.62
1:B:128:ARG:NH1	5:B:219:HOH:O	2.32	0.62
3:E:256:ARG:NH1	3:E:272:GLU:OE1	2.32	0.61
1:A:109:LEU:HD21	3:F:321:GLY:HA3	1.82	0.61
1:A:122:LYS:NZ	3:F:384:ASP:OD2	2.29	0.61
3:F:238:ARG:NH2	5:F:507:HOH:O	2.33	0.60
3:F:253:ILE:HG12	3:F:327:ARG:HD3	1.82	0.60
1:A:48:LEU:N	4:E:401:PO4:O2	2.25	0.59
3:F:191:LEU:HD22	3:F:235:LEU:HD11	1.85	0.58
3:F:304:LEU:HB3	3:F:309:LEU:HD13	1.87	0.57
3:E:263:ARG:NH1	3:F:192:GLU:OE2	2.38	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:83:ARG:HH22	2:C:79:LYS:HE3	1.70	0.56
3:E:225:GLU:HG2	3:E:229:LYS:HE3	1.87	0.56
1:A:63:ARG:N	4:C:201:PO4:O1	2.36	0.56
1:B:81:ASP:OD1	1:B:81:ASP:N	2.38	0.56
2:D:30:THR:N	4:D:201:PO4:O4	2.36	0.56
2:C:30:THR:N	4:C:201:PO4:O3	2.25	0.55
3:F:207:GLU:HG2	3:F:340:LEU:HD21	1.89	0.55
1:B:112:ILE:O	1:B:113:HIS:ND1	2.39	0.54
3:F:277:LYS:HB2	3:F:282:THR:HG23	1.88	0.54
3:F:251:ARG:HD2	4:F:402:PO4:O2	2.08	0.53
2:D:100:PHE:O	3:F:318:ARG:NH2	2.41	0.53
1:A:131:ARG:NH2	2:D:99:GLY:O	2.41	0.52
2:D:65:VAL:HG22	2:D:93:GLN:HG3	1.91	0.52
1:A:46:VAL:O	3:E:263:ARG:NH2	2.40	0.52
3:F:266:GLU:H	3:F:266:GLU:CD	2.16	0.51
3:F:297:ALA:HA	3:F:302:LEU:HD12	1.95	0.49
3:F:210:LEU:HD21	3:F:221:ALA:HB3	1.95	0.49
1:B:63:ARG:N	4:D:201:PO4:O2	2.37	0.47
3:F:256:ARG:NH1	3:F:272:GLU:OE1	2.47	0.47
1:A:53:ARG:NH2	3:F:342:ASP:OD2	2.35	0.47
1:A:81:ASP:OD2	1:A:81:ASP:N	2.39	0.47
3:F:359:ARG:HG3	3:F:362:ARG:HH12	1.79	0.46
1:B:48:LEU:N	4:F:401:PO4:O4	2.31	0.46
2:D:60:VAL:HG21	3:E:375:VAL:HG21	1.97	0.46
3:E:323:ARG:NH1	5:E:521:HOH:O	2.36	0.45
3:F:251:ARG:O	3:F:254:GLU:HG2	2.17	0.44
3:F:265:PRO:O	3:F:269:ARG:HD3	2.17	0.44
3:F:205:LEU:HD12	3:F:228:LEU:HD12	1.99	0.44
1:A:99:PHE:HZ	3:F:372:LEU:HD21	1.82	0.44
3:E:312:MET:HE3	3:E:312:MET:HB2	1.93	0.43
1:A:107:THR:HA	1:A:127:ALA:HB2	2.00	0.43
3:E:206:GLN:HG3	3:E:225:GLU:OE1	2.18	0.43
3:F:339:HIS:CE1	3:F:340:LEU:HG	2.54	0.43
3:E:346:PRO:HA	3:E:347:GLY:HA3	1.67	0.43
1:A:65:LEU:HD13	3:F:218:PRO:HA	2.01	0.42
2:D:95:ARG:O	3:E:329:HIS:HE1	2.02	0.42
3:E:239:LEU:O	3:E:243:LYS:HB2	2.20	0.42
1:B:65:LEU:HB3	1:B:66:PRO:HD3	2.02	0.42
3:F:309:LEU:HD12	3:F:309:LEU:HA	1.88	0.41
1:A:48:LEU:HD12	1:A:48:LEU:HA	1.86	0.41
1:A:119:ILE:HD13	1:A:119:ILE:HA	1.88	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:61:LEU:HD12	2:D:37:LEU:HD23	2.03	0.40
3:E:277:LYS:O	3:E:282:THR:HG21	2.21	0.40
2:C:97:LEU:HD23	2:C:97:LEU:HA	1.71	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	87/135 (64%)	85 (98%)	2 (2%)	0	100	100
1	B	92/135 (68%)	91 (99%)	1 (1%)	0	100	100
2	C	72/83 (87%)	72 (100%)	0	0	100	100
2	D	72/83 (87%)	72 (100%)	0	0	100	100
3	E	195/216 (90%)	193 (99%)	2 (1%)	0	100	100
3	F	199/216 (92%)	196 (98%)	3 (2%)	0	100	100
All	All	717/868 (83%)	709 (99%)	8 (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	74/110 (67%)	73 (99%)	1 (1%)	59	81
1	B	75/110 (68%)	72 (96%)	3 (4%)	28	55
2	C	58/68 (85%)	56 (97%)	2 (3%)	32	60
2	D	58/68 (85%)	56 (97%)	2 (3%)	32	60
3	E	172/190 (90%)	161 (94%)	11 (6%)	16	35
3	F	177/190 (93%)	159 (90%)	18 (10%)	7	15
All	All	614/736 (83%)	577 (94%)	37 (6%)	17	37

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	103	LEU
1	B	81	ASP
1	B	117	VAL
1	B	126	LEU
2	C	29	ILE
2	C	44	LYS
2	D	29	ILE
2	D	45	ARG
3	E	189	GLN
3	E	253	ILE
3	E	260	ARG
3	E	266	GLU
3	E	282	THR
3	E	295	LYS
3	E	301	SER
3	E	311	LEU
3	E	320	VAL
3	E	327	ARG
3	E	370	SER
3	F	184	SER
3	F	192	GLU
3	F	204	ARG
3	F	241	GLU
3	F	244	ASP
3	F	253	ILE
3	F	269	ARG
3	F	282	THR
3	F	301	SER
3	F	304	LEU
3	F	309	LEU

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Mol	Chain	Res	Type
3	F	327	ARG
3	F	330	LEU
3	F	343	ASP
3	F	356	VAL
3	F	359	ARG
3	F	371	ARG
3	F	387	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	85	GLN
3	E	308	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](#)

5 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	PO4	E	401	-	4,4,4	1.06	0	6,6,6	0.45	0
4	PO4	D	201	-	4,4,4	1.03	0	6,6,6	0.47	0
4	PO4	F	402	-	4,4,4	1.02	0	6,6,6	0.47	0
4	PO4	C	201	-	4,4,4	0.90	0	6,6,6	0.67	0
4	PO4	F	401	-	4,4,4	0.92	0	6,6,6	0.55	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.

5 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	E	401	PO4	1	0
4	D	201	PO4	2	0
4	F	402	PO4	2	0
4	C	201	PO4	2	0
4	F	401	PO4	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	89/135 (65%)	-0.20	1 (1%) 78 74	34, 48, 81, 96	0
1	B	94/135 (69%)	-0.08	2 (2%) 63 58	35, 50, 89, 107	0
2	C	74/83 (89%)	-0.27	1 (1%) 73 69	35, 50, 75, 110	0
2	D	74/83 (89%)	-0.17	2 (2%) 56 50	34, 48, 69, 100	0
3	E	199/216 (92%)	0.03	7 (3%) 47 41	37, 57, 95, 122	0
3	F	203/216 (93%)	-0.01	2 (0%) 79 76	36, 57, 90, 110	0
All	All	733/868 (84%)	-0.07	15 (2%) 65 60	34, 54, 89, 122	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	133	GLU	3.8
1	B	133	GLU	3.4
3	F	347	GLY	3.4
3	E	346	PRO	3.3
3	F	344	TYR	3.3
3	E	340	LEU	3.2
3	E	244	ASP	3.1
2	D	28	GLY	2.9
2	C	27	GLN	2.7
3	E	339	HIS	2.5
2	D	27	GLN	2.4
1	B	41	TYR	2.2
3	E	184	SER	2.2
3	E	281	ASP	2.1
3	E	301	SER	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	PO4	F	401	5/5	0.88	0.29	75,77,92,96	0
4	PO4	F	402	5/5	0.88	0.33	81,87,89,90	0
4	PO4	C	201	5/5	0.89	0.39	77,78,84,95	0
4	PO4	D	201	5/5	0.89	0.37	74,80,87,89	0
4	PO4	E	401	5/5	0.90	0.29	69,78,93,98	0

6.5 Other polymers [i](#)

There are no such residues in this entry.