



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:28 PM GMT

PDB ID : 4IG9
Title : Structure of NAD-dependent protein deacetylase sirtuin-1 (open state, 2.64 Å)
Authors : Davenport, A.M.; Huber, F.M.; Hoelz, A.
Deposited on : 2012-12-16
Resolution : 2.64 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

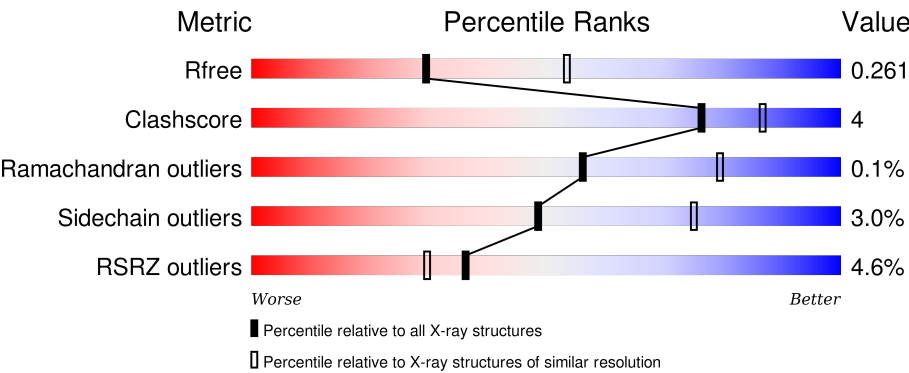
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.64 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	3377 (2.68-2.60)
Clashscore	102246	3781 (2.68-2.60)
Ramachandran outliers	100387	3722 (2.68-2.60)
Sidechain outliers	100360	3722 (2.68-2.60)
RSRZ outliers	91569	3388 (2.68-2.60)

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 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 <=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	281	<div><div></div><div><div></div><div></div><div></div><div></div></div><div>85%10% .</div></div>
1	C	281	<div><div>7%</div><div></div><div></div><div></div><div></div></div> <div>86%9% . .</div>

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Mol	Chain	Length	Quality of chain
2	D	31	<div><div><div></div><div></div><div></div></div><div><div>3%</div><div>39%</div><div>13%</div><div>48%</div></div></div>
2	F	31	<div><div><div></div><div></div><div></div></div><div><div>6%</div><div>45%</div><div>13%</div><div>42%</div></div></div>
2	H	31	<div><div><div></div><div></div><div></div></div><div><div>6%</div><div>68%</div><div>•</div><div>29%</div></div></div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9419 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 NAD-dependent protein deacetylase sirtuin-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	269	Total	C	N	O	S	0	3	0
			2162	1385	372	389	16			
1	C	269	Total	C	N	O	S	0	4	0
			2173	1391	376	390	16			
1	E	270	Total	C	N	O	S	0	4	0
			2183	1397	379	391	16			
1	G	270	Total	C	N	O	S	0	4	0
			2179	1394	377	391	17			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	EXPRESSION TAG	UNP Q96EB6
A	-3	PRO	-	EXPRESSION TAG	UNP Q96EB6
A	-2	HIS	-	EXPRESSION TAG	UNP Q96EB6
A	-1	MET	-	EXPRESSION TAG	UNP Q96EB6
C	-4	GLY	-	EXPRESSION TAG	UNP Q96EB6
C	-3	PRO	-	EXPRESSION TAG	UNP Q96EB6
C	-2	HIS	-	EXPRESSION TAG	UNP Q96EB6
C	-1	MET	-	EXPRESSION TAG	UNP Q96EB6
E	-4	GLY	-	EXPRESSION TAG	UNP Q96EB6
E	-3	PRO	-	EXPRESSION TAG	UNP Q96EB6
E	-2	HIS	-	EXPRESSION TAG	UNP Q96EB6
E	-1	MET	-	EXPRESSION TAG	UNP Q96EB6
G	-4	GLY	-	EXPRESSION TAG	UNP Q96EB6
G	-3	PRO	-	EXPRESSION TAG	UNP Q96EB6
G	-2	HIS	-	EXPRESSION TAG	UNP Q96EB6
G	-1	MET	-	EXPRESSION TAG	UNP Q96EB6

- Molecule 2 is a protein called NAD-dependent protein deacetylase sirtuin-1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	21	Total	C	N	O	0	0	0
			175	117	28	30			
2	D	16	Total	C	N	O	0	0	0
			140	95	23	21			
2	F	18	Total	C	N	O	0	2	0
			174	116	32	25			
2	H	22	Total	C	N	O	0	0	0
			183	122	29	31			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-6	GLY	-	EXPRESSION TAG	UNP Q96EB6
B	-5	PRO	-	EXPRESSION TAG	UNP Q96EB6
B	-4	HIS	-	EXPRESSION TAG	UNP Q96EB6
B	-3	MET	-	EXPRESSION TAG	UNP Q96EB6
B	-2	GLY	-	EXPRESSION TAG	UNP Q96EB6
B	-1	SER	-	EXPRESSION TAG	UNP Q96EB6
D	-6	GLY	-	EXPRESSION TAG	UNP Q96EB6
D	-5	PRO	-	EXPRESSION TAG	UNP Q96EB6
D	-4	HIS	-	EXPRESSION TAG	UNP Q96EB6
D	-3	MET	-	EXPRESSION TAG	UNP Q96EB6
D	-2	GLY	-	EXPRESSION TAG	UNP Q96EB6
D	-1	SER	-	EXPRESSION TAG	UNP Q96EB6
F	-6	GLY	-	EXPRESSION TAG	UNP Q96EB6
F	-5	PRO	-	EXPRESSION TAG	UNP Q96EB6
F	-4	HIS	-	EXPRESSION TAG	UNP Q96EB6
F	-3	MET	-	EXPRESSION TAG	UNP Q96EB6
F	-2	GLY	-	EXPRESSION TAG	UNP Q96EB6
F	-1	SER	-	EXPRESSION TAG	UNP Q96EB6
H	-6	GLY	-	EXPRESSION TAG	UNP Q96EB6
H	-5	PRO	-	EXPRESSION TAG	UNP Q96EB6
H	-4	HIS	-	EXPRESSION TAG	UNP Q96EB6
H	-3	MET	-	EXPRESSION TAG	UNP Q96EB6
H	-2	GLY	-	EXPRESSION TAG	UNP Q96EB6
H	-1	SER	-	EXPRESSION TAG	UNP Q96EB6

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	G	1	Total	Zn	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total 1	Zn 1	0	0
3	C	1	Total 1	Zn 1	0	0
3	E	1	Total 1	Zn 1	0	0

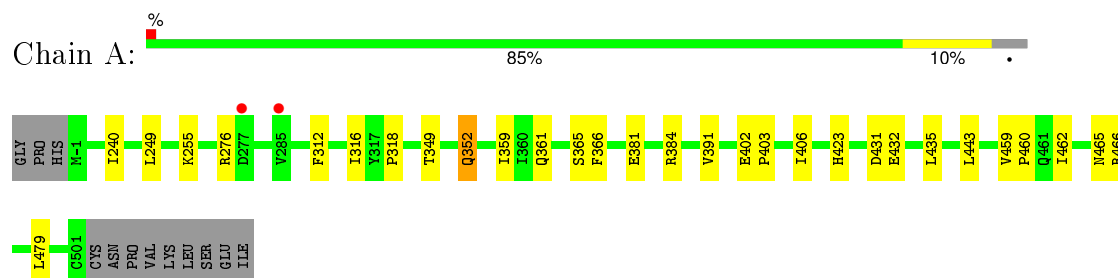
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	22	Total 22	O 22	0	0
4	B	2	Total 2	O 2	0	0
4	C	5	Total 5	O 5	0	0
4	D	2	Total 2	O 2	0	0
4	E	5	Total 5	O 5	0	0
4	F	1	Total 1	O 1	0	0
4	G	8	Total 8	O 8	0	0
4	H	1	Total 1	O 1	0	0

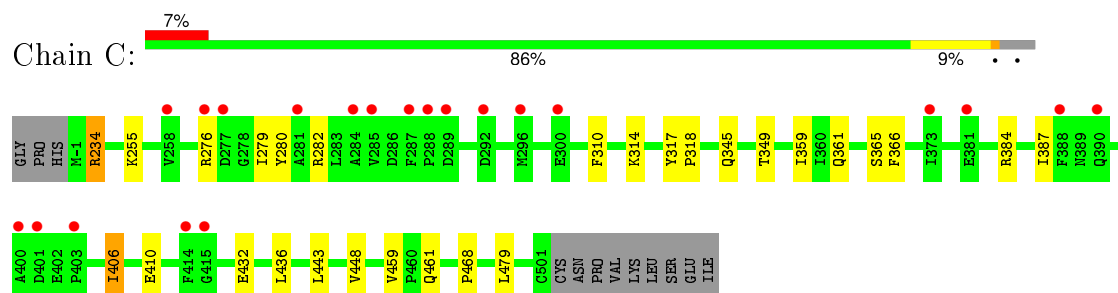
3 Residue-property plots

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.

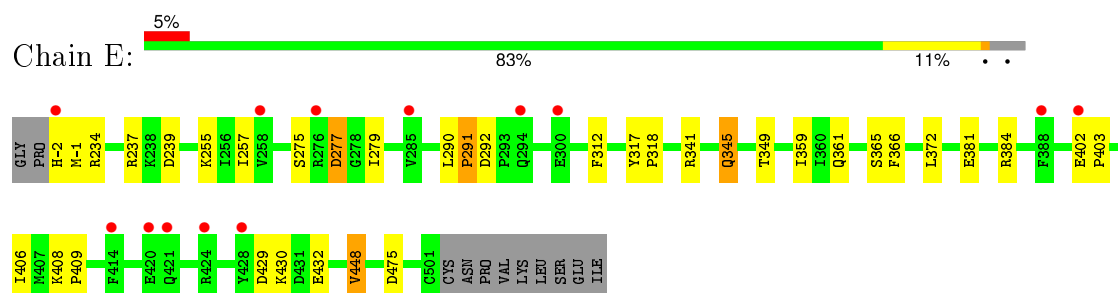
- Molecule 1: NAD-dependent protein deacetylase sirtuin-1



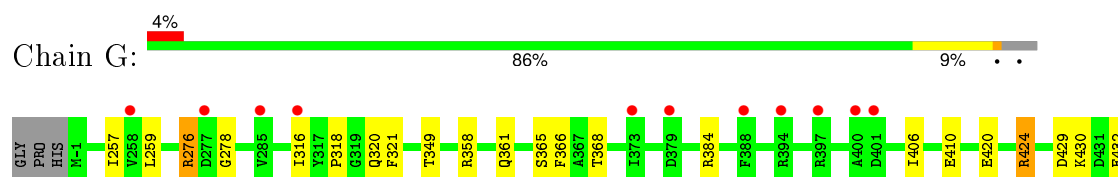
- Molecule 1: NAD-dependent protein deacetylase sirtuin-1



- Molecule 1: NAD-dependent protein deacetylase sirtuin-1



- Molecule 1: NAD-dependent protein deacetylase sirtuin-1

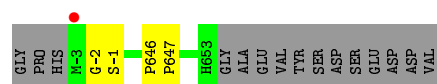
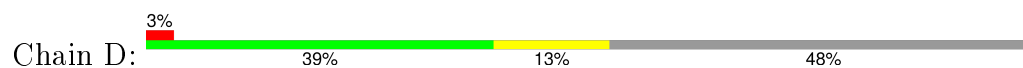




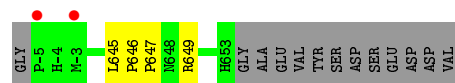
- Molecule 2: NAD-dependent protein deacetylase sirtuin-1



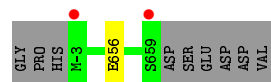
- Molecule 2: NAD-dependent protein deacetylase sirtuin-1



- Molecule 2: NAD-dependent protein deacetylase sirtuin-1



- Molecule 2: NAD-dependent protein deacetylase sirtuin-1



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	115.78Å 115.78Å 350.51Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.94 – 2.64 19.94 – 2.64	Depositor EDS
% Data completeness (in resolution range)	96.9 (19.94-2.64) 88.7 (19.94-2.64)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.26 (at 2.63Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.227 , 0.263 0.221 , 0.261	Depositor DCC
R_{free} test set	1833 reflections (2.93%)	DCC
Wilson B-factor (Å ²)	65.5	Xtriage
Anisotropy	0.469	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 37.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 68500 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9419	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.05% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

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	0.21	0/2210	0.38	0/2987
1	C	0.21	0/2221	0.37	0/3001
1	E	0.21	0/2232	0.38	0/3016
1	G	0.21	0/2227	0.38	0/3009
2	B	0.23	0/182	0.39	0/247
2	D	0.24	0/146	0.43	0/197
2	F	0.22	0/185	0.39	0/249
2	H	0.24	0/190	0.42	0/257
All	All	0.21	0/9593	0.38	0/12963

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	2162	0	2192	18	0
1	C	2173	0	2205	15	0
1	E	2183	0	2211	19	0
1	G	2179	0	2209	15	0
2	B	175	0	164	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	140	0	136	3	0
2	F	174	0	170	3	0
2	H	183	0	173	1	0
3	A	1	0	0	0	0
3	C	1	0	0	0	0
3	E	1	0	0	0	0
3	G	1	0	0	0	0
4	A	22	0	0	0	0
4	B	2	0	0	0	0
4	C	5	0	0	0	0
4	D	2	0	0	0	0
4	E	5	0	0	0	0
4	F	1	0	0	0	0
4	G	8	0	0	0	0
4	H	1	0	0	0	0
All	All	9419	0	9460	71	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 4.

All (71) 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:G:349:THR:HG22	1:G:365:SER:HB2	1.62	0.81
1:A:349:THR:HG22	1:A:365:SER:HB2	1.63	0.81
1:E:349:THR:HG22	1:E:365:SER:HB2	1.64	0.79
1:A:431:ASP:O	1:G:358:ARG:NH2	2.21	0.72
1:A:466:ARG:NH1	1:E:-2:HIS:O	2.22	0.71
1:E:255:LYS:HD3	1:E:432:GLU:HG3	1.76	0.68
1:C:349:THR:HG22	1:C:365:SER:HB2	1.74	0.68
1:G:318:PRO:O	1:G:384[A]:ARG:NH2	2.30	0.63
1:A:255:LYS:HD3	1:A:432:GLU:HG3	1.82	0.61
1:G:466:ARG:NH1	2:H:656:GLU:OE1	2.33	0.61
1:E:372:LEU:HD21	1:E:408:LYS:HD3	1.83	0.60
1:C:255:LYS:HD3	1:C:432:GLU:HG3	1.85	0.58
1:E:318:PRO:O	1:E:384[A]:ARG:NH2	2.38	0.57
1:A:318:PRO:O	1:A:384:ARG:NH2	2.39	0.55
1:E:349:THR:HG21	1:E:366:PHE:HB2	1.90	0.53
2:B:646:PRO:HB3	2:D:-2:GLY:H	1.73	0.53
1:C:310:PHE:HZ	1:C:406:ILE:HD11	1.72	0.53
1:C:349:THR:CG2	1:C:366:PHE:H	2.22	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:318:PRO:O	1:C:384[A]:ARG:NH2	2.43	0.51
1:E:275:SER:OG	1:E:277:ASP:O	2.28	0.51
1:G:276:ARG:HA	1:G:278:GLY:H	1.75	0.51
1:A:349:THR:CG2	1:A:366:PHE:H	2.23	0.51
1:G:257:ILE:HD11	1:G:429:ASP:HB3	1.92	0.51
1:C:365:SER:N	1:C:410:GLU:O	2.44	0.51
1:E:345:GLN:HG2	1:E:448:VAL:HG21	1.93	0.51
1:A:381:GLU:OE1	1:A:384:ARG:NH1	2.44	0.50
1:G:276:ARG:HA	1:G:278:GLY:N	2.27	0.49
1:A:352:GLN:HE22	1:A:361:GLN:HE22	1.60	0.49
1:G:420:GLU:HG2	1:G:424:ARG:HD2	1.95	0.48
1:C:349:THR:HG21	1:C:366:PHE:H	1.79	0.48
1:G:494:GLY:HA2	1:G:498:ALA:HB2	1.96	0.48
1:C:314:LYS:HB3	1:C:387:ILE:HG22	1.96	0.48
1:E:237:ARG:NH2	1:E:239:ASP:OD2	2.39	0.48
1:E:381:GLU:OE1	1:E:384[A]:ARG:NH1	2.47	0.47
1:A:435:LEU:HD11	1:A:462:ILE:HD13	1.96	0.47
1:G:442:SER:H	1:G:444:LYS:HE3	1.78	0.47
1:A:349:THR:HG21	1:A:366:PHE:H	1.80	0.47
1:G:349:THR:CG2	1:G:366:PHE:H	2.27	0.47
2:F:645:LEU:HB2	2:F:649[A]:ARG:HB3	1.98	0.46
1:G:259:LEU:HD21	1:G:451:ILE:HB	1.96	0.46
1:E:234:ARG:NH2	1:E:475:ASP:OD2	2.48	0.45
1:A:240:ILE:HD13	1:A:249:LEU:HD11	1.97	0.45
1:A:349:THR:HG21	1:A:366:PHE:HB2	1.98	0.45
1:C:279:ILE:HG13	1:C:282:ARG:HH21	1.81	0.45
1:C:234:ARG:NH2	1:C:459:VAL:O	2.50	0.44
2:F:645:LEU:HB2	2:F:649[B]:ARG:HB2	2.00	0.44
1:E:279:ILE:HD13	1:E:312:PHE:HD1	1.82	0.44
1:A:352:GLN:NE2	1:A:361:GLN:HE22	2.15	0.44
1:A:465:ASN:O	1:A:479:LEU:HA	2.18	0.44
2:D:-2:GLY:HA2	2:D:-1:SER:HA	1.68	0.44
2:F:646:PRO:HA	2:F:647:PRO:HA	1.77	0.43
1:C:276:ARG:HH11	1:C:280:TYR:HE2	1.66	0.43
1:E:349:THR:CG2	1:E:366:PHE:H	2.30	0.43
1:C:317:TYR:HA	1:C:318:PRO:HD2	1.80	0.43
1:G:368:THR:OG1	1:G:410:GLU:HG3	2.19	0.43
1:A:312:PHE:CE1	1:A:316:ILE:HG13	2.53	0.43
2:D:646:PRO:HA	2:D:647:PRO:HA	1.75	0.43
1:E:290:LEU:HA	1:E:291:PRO:HD3	1.86	0.43
1:A:352:GLN:HE21	1:A:352:GLN:H	1.67	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:345:GLN:HG2	1:C:448:VAL:HG21	2.00	0.43
1:C:436:LEU:HB3	1:C:461:GLN:HG2	2.01	0.42
1:E:257:ILE:HD11	1:E:429:ASP:HB3	2.02	0.42
1:C:468:PRO:HD3	1:C:479:LEU:HD13	2.02	0.42
1:G:349:THR:HG21	1:G:366:PHE:H	1.86	0.41
1:A:459:VAL:HA	1:A:460:PRO:HD3	1.93	0.41
1:E:408:LYS:HA	1:E:409:PRO:HD3	1.96	0.41
1:G:316:ILE:HA	1:G:321:PHE:HE2	1.86	0.41
1:E:402:GLU:HA	1:E:403:PRO:HD3	1.87	0.41
1:E:349:THR:HG22	1:E:366:PHE:H	1.86	0.40
1:E:341:ARG:HG3	1:E:429:ASP:OD2	2.21	0.40
1:A:402:GLU:HA	1:A:403:PRO:HD3	1.87	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	270/281 (96%)	261 (97%)	9 (3%)	0	100	100
1	C	271/281 (96%)	265 (98%)	6 (2%)	0	100	100
1	E	272/281 (97%)	261 (96%)	10 (4%)	1 (0%)	39	63
1	G	272/281 (97%)	262 (96%)	10 (4%)	0	100	100
2	B	19/31 (61%)	16 (84%)	3 (16%)	0	100	100
2	D	14/31 (45%)	12 (86%)	2 (14%)	0	100	100
2	F	18/31 (58%)	17 (94%)	1 (6%)	0	100	100
2	H	20/31 (64%)	18 (90%)	2 (10%)	0	100	100
All	All	1156/1248 (93%)	1112 (96%)	43 (4%)	1 (0%)	56	79

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	291	PRO

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	242/250 (97%)	235 (97%)	7 (3%)	50	76
1	C	243/250 (97%)	238 (98%)	5 (2%)	61	84
1	E	244/250 (98%)	234 (96%)	10 (4%)	37	64
1	G	244/250 (98%)	235 (96%)	9 (4%)	41	67
2	B	18/27 (67%)	18 (100%)	0	100	100
2	D	15/27 (56%)	15 (100%)	0	100	100
2	F	19/27 (70%)	19 (100%)	0	100	100
2	H	19/27 (70%)	19 (100%)	0	100	100
All	All	1044/1108 (94%)	1013 (97%)	31 (3%)	48	75

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	276	ARG
1	A	352	GLN
1	A	359	ILE
1	A	391	VAL
1	A	406	ILE
1	A	423	HIS
1	A	443	LEU
1	C	234	ARG
1	C	359	ILE
1	C	361	GLN
1	C	406	ILE
1	C	443	LEU
1	E	-1	MET
1	E	277	ASP
1	E	292	ASP

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Mol	Chain	Res	Type
1	E	317	TYR
1	E	345	GLN
1	E	359	ILE
1	E	361	GLN
1	E	406	ILE
1	E	430	LYS
1	E	448	VAL
1	G	276	ARG
1	G	320	GLN
1	G	361	GLN
1	G	406	ILE
1	G	424	ARG
1	G	430	LYS
1	G	432	GLU
1	G	443	LEU
1	G	444	LYS

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

Mol	Chain	Res	Type
1	A	361	GLN
2	F	-4	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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 ⓘ

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

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 [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	269/281 (95%)	-0.06	2 (0%) 89 87	26, 44, 77, 105	0
1	C	269/281 (95%)	0.37	21 (7%) 16 11	30, 62, 122, 144	0
1	E	270/281 (96%)	0.30	13 (4%) 34 28	30, 62, 109, 130	0
1	G	270/281 (96%)	0.09	12 (4%) 38 31	27, 50, 105, 129	0
2	B	21/31 (67%)	-0.38	0 100 100	28, 38, 52, 78	0
2	D	16/31 (51%)	-0.18	1 (6%) 23 18	30, 36, 69, 85	0
2	F	18/31 (58%)	0.40	2 (11%) 7 4	29, 40, 70, 88	0
2	H	22/31 (70%)	-0.10	2 (9%) 11 8	31, 39, 65, 83	0
All	All	1155/1248 (92%)	0.16	53 (4%) 36 30	26, 54, 109, 144	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	285	VAL	6.8
2	F	-5	PRO	4.8
1	C	277	ASP	4.7
1	E	424	ARG	4.5
1	E	428	TYR	4.4
1	A	277	ASP	4.4
1	E	420	GLU	4.3
1	C	414	PHE	4.2
1	C	281	ALA	4.1
1	C	289	ASP	4.0
1	C	276	ARG	4.0
1	E	285	VAL	3.7
1	C	400	ALA	3.5
1	C	401	ASP	3.5
1	E	294	GLN	3.4
2	F	-3	MET	3.3

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Mol	Chain	Res	Type	RSRZ
1	C	373	ILE	3.3
1	G	502	CYS	3.2
1	G	285	VAL	3.1
1	G	401	ASP	3.1
1	E	402	GLU	3.0
1	E	-2	HIS	3.0
1	C	284	ALA	3.0
1	C	390	GLN	2.9
1	G	388	PHE	2.8
1	C	287	PHE	2.8
2	H	-3	MET	2.7
1	C	403	PRO	2.7
1	G	397	ARG	2.7
1	G	277	ASP	2.7
1	C	288	PRO	2.6
1	C	292	ASP	2.5
1	C	415	GLY	2.5
1	E	276	ARG	2.5
1	E	421	GLN	2.5
1	G	258	VAL	2.5
1	C	300	GLU	2.3
1	C	381	GLU	2.3
1	G	373	ILE	2.3
1	G	394	ARG	2.3
1	E	300	GLU	2.3
1	E	388	PHE	2.3
1	E	414	PHE	2.2
2	H	659	SER	2.2
1	A	285	VAL	2.2
1	C	296	MET	2.2
2	D	-3	MET	2.1
1	C	388	PHE	2.1
1	G	400	ALA	2.1
1	C	258	VAL	2.0
1	E	258	VAL	2.0
1	G	316	ILE	2.0
1	G	379	ASP	2.0

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

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(Å ²)	Q<0.9
3	ZN	E	601	1/1	0.96	0.12	-0.85	69,69,69,69	0
3	ZN	A	601	1/1	0.99	0.09	-1.43	71,71,71,71	0
3	ZN	G	601	1/1	0.90	0.07	-1.84	112,112,112,112	0
3	ZN	C	601	1/1	0.90	0.06	-2.00	107,107,107,107	0

6.5 Other polymers [i](#)

There are no such residues in this entry.