



wwPDB X-ray Structure Validation Summary Report ⓘ

Dec 14, 2016 – 12:10 AM EST

PDB ID : 5HZN
Title : Structure of NVP-AEW541 in complex with IGF-1R kinase
Authors : Cowan-Jacob, S.W.
Deposited on : 2016-02-02
Resolution : 2.20 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20028442
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) : rb-20028442

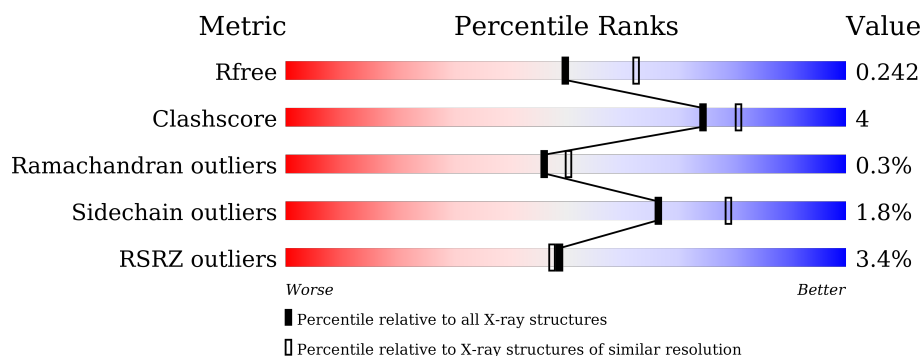
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

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 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	304	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 1%, orange 1%, orange 88%, yellow 88%, yellow 99%, grey 99%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 88% 9% . </div> </div>
1	B	304	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 7%, orange 7%, orange 83%, yellow 83%, yellow 96%, grey 96%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 7% 83% 13% .. </div> </div>
1	C	304	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 6%, orange 6%, orange 87%, yellow 87%, yellow 97%, grey 97%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 6% 87% 10% .. </div> </div>
1	D	304	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 2%, orange 2%, orange 88%, yellow 88%, yellow 99%, grey 99%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 2% 88% 11% . </div> </div>
1	E	304	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 1%, orange 1%, orange 87%, yellow 87%, yellow 97%, grey 97%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 87% 10% . </div> </div>
1	F	304	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 5%, orange 5%, orange 84%, yellow 84%, yellow 96%, grey 96%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 5% 84% 12% .. </div> </div>

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Mol	Chain	Length	Quality of chain
1	G	304	<div><div></div><div>5%</div><div>87%</div><div>10%</div><div>• •</div></div>
1	H	304	<div><div></div><div>89%</div><div>11%</div></div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 19906 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 Insulin-like growth factor 1 receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	296	Total	C	N	O	S	0	0	0
			2359	1499	394	444	22			
1	B	296	Total	C	N	O	S	0	0	0
			2359	1499	394	444	22			
1	C	296	Total	C	N	O	S	0	0	0
			2359	1499	394	444	22			
1	D	304	Total	C	N	O	S	0	0	0
			2415	1534	404	454	23			
1	E	296	Total	C	N	O	S	0	0	0
			2359	1499	394	444	22			
1	F	296	Total	C	N	O	S	0	0	0
			2359	1499	394	444	22			
1	G	296	Total	C	N	O	S	0	0	0
			2359	1499	394	444	22			
1	H	303	Total	C	N	O	S	0	0	0
			2406	1528	402	453	23			

There are 16 discrepancies between the modelled and reference sequences:

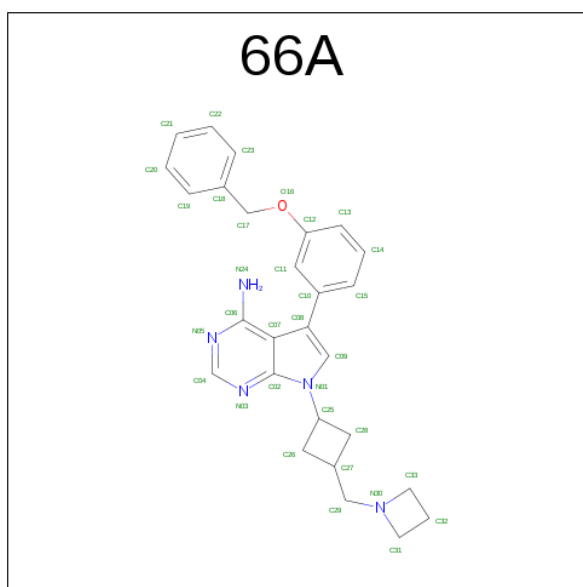
Chain	Residue	Modelled	Actual	Comment	Reference
A	1094	ALA	GLU	engineered mutation	UNP P08069
A	1096	ALA	GLU	engineered mutation	UNP P08069
B	1094	ALA	GLU	engineered mutation	UNP P08069
B	1096	ALA	GLU	engineered mutation	UNP P08069
C	1094	ALA	GLU	engineered mutation	UNP P08069
C	1096	ALA	GLU	engineered mutation	UNP P08069
D	1094	ALA	GLU	engineered mutation	UNP P08069
D	1096	ALA	GLU	engineered mutation	UNP P08069
E	1094	ALA	GLU	engineered mutation	UNP P08069
E	1096	ALA	GLU	engineered mutation	UNP P08069
F	1094	ALA	GLU	engineered mutation	UNP P08069
F	1096	ALA	GLU	engineered mutation	UNP P08069
G	1094	ALA	GLU	engineered mutation	UNP P08069

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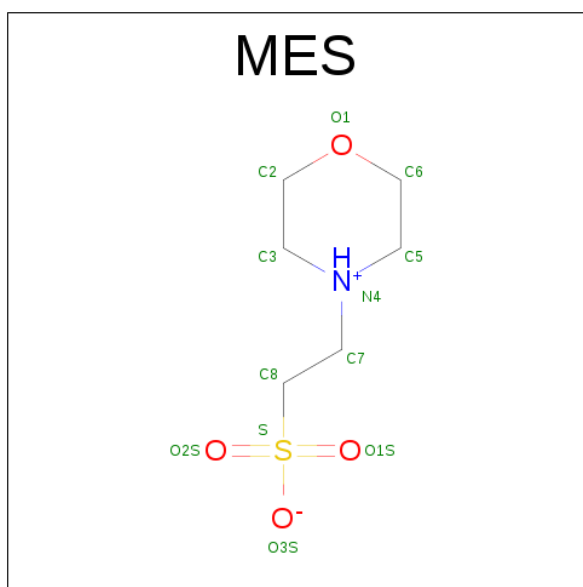
Chain	Residue	Modelled	Actual	Comment	Reference
G	1096	ALA	GLU	engineered mutation	UNP P08069
H	1094	ALA	GLU	engineered mutation	UNP P08069
H	1096	ALA	GLU	engineered mutation	UNP P08069

- Molecule 2 is 7-[cis-3-(azetidin-1-ylmethyl)cyclobutyl]-5-[3-(benzyloxy)phenyl]-7H-pyrrolo[2,3-d]pyrimidin-4-amine (three-letter code: 66A) (formula: C₂₇H₂₉N₅O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			33	27	5	1		
2	B	1	Total	C	N	O	0	0
			33	27	5	1		
2	C	1	Total	C	N	O	0	0
			33	27	5	1		
2	D	1	Total	C	N	O	0	0
			33	27	5	1		
2	E	1	Total	C	N	O	0	0
			33	27	5	1		
2	F	1	Total	C	N	O	0	0
			33	27	5	1		
2	G	1	Total	C	N	O	0	0
			33	27	5	1		
2	H	1	Total	C	N	O	0	0
			33	27	5	1		

- Molecule 3 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: C₆H₁₃NO₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
3	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
3	C	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
3	D	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
3	E	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
3	F	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
3	G	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
3	H	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	G	1	Total	O	S	0	0
			5	4	1		

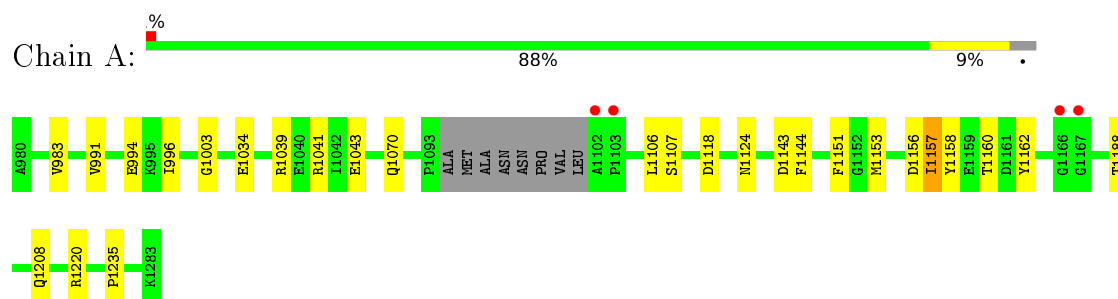
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	92	Total	O	0	0
			92	92		
5	B	42	Total	O	0	0
			42	42		
5	C	49	Total	O	0	0
			49	49		
5	D	92	Total	O	0	0
			92	92		
5	E	97	Total	O	0	0
			97	97		
5	F	43	Total	O	0	0
			43	43		
5	G	48	Total	O	0	0
			48	48		
5	H	88	Total	O	0	0
			88	88		

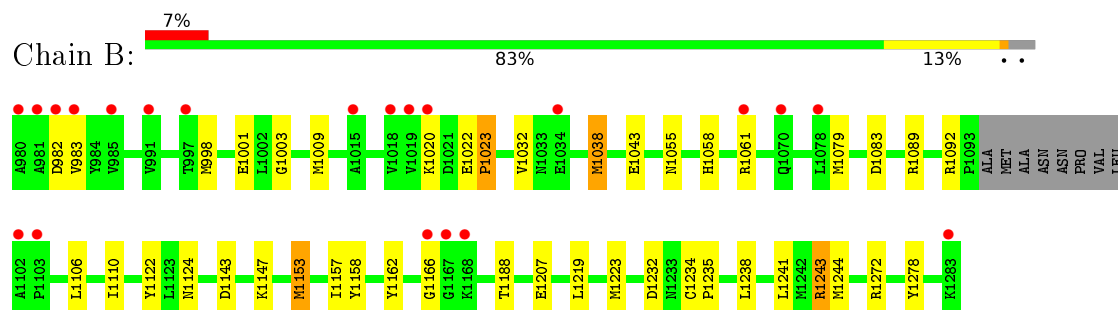
3 Residue-property plots

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.

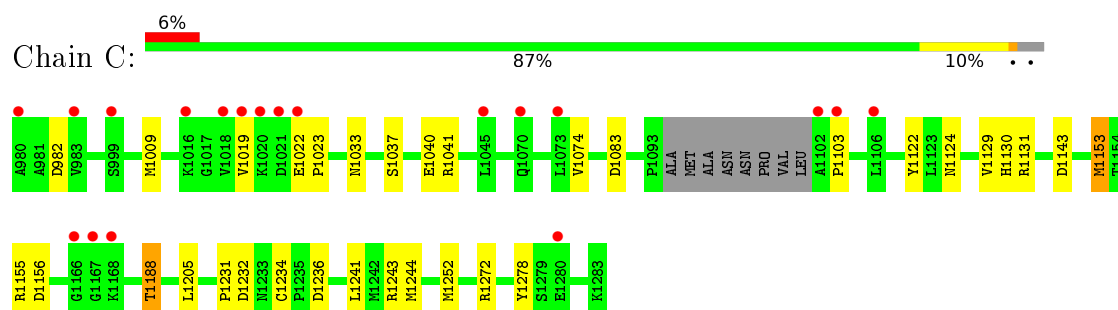
- Molecule 1: Insulin-like growth factor 1 receptor



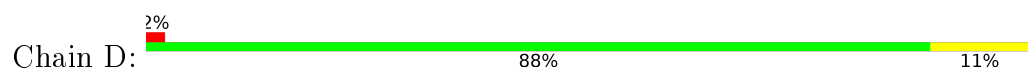
- Molecule 1: Insulin-like growth factor 1 receptor

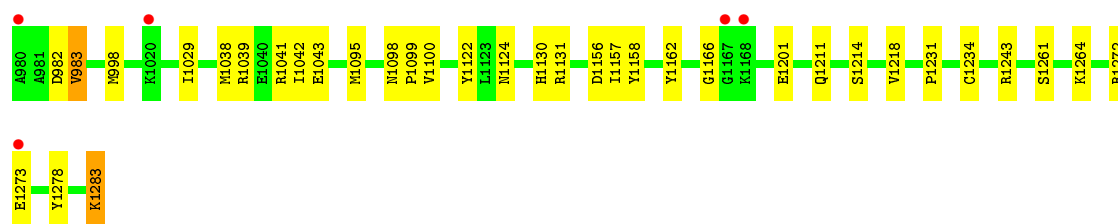


- Molecule 1: Insulin-like growth factor 1 receptor

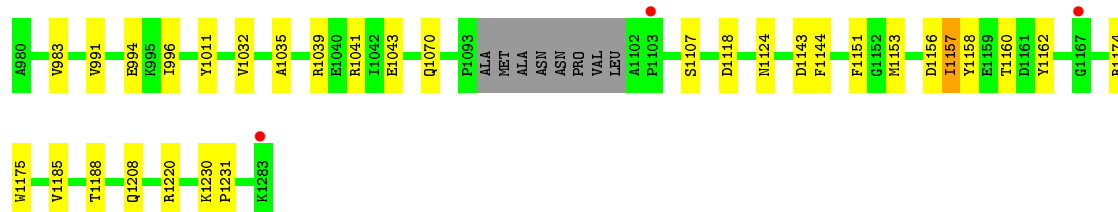
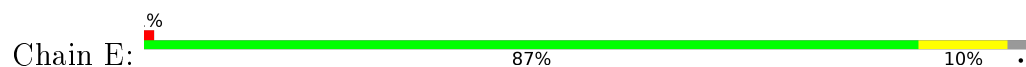


- Molecule 1: Insulin-like growth factor 1 receptor

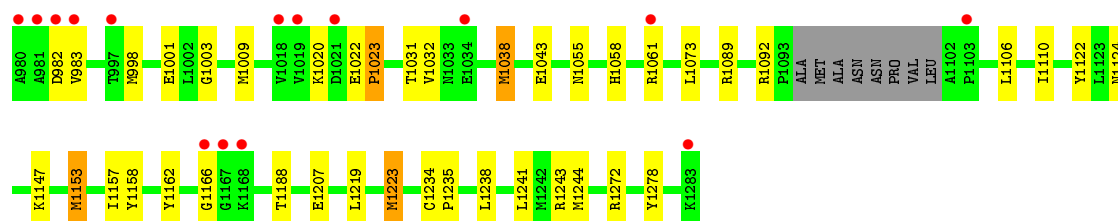
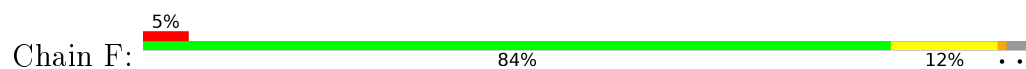




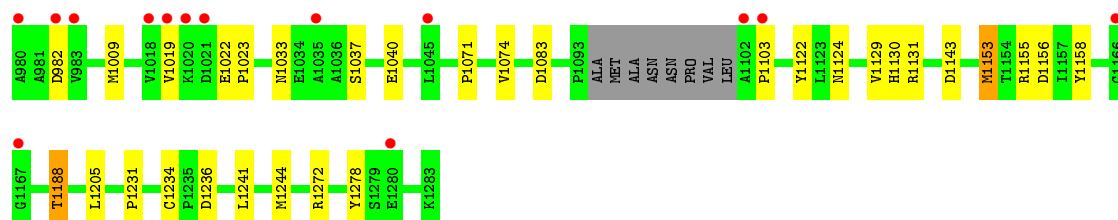
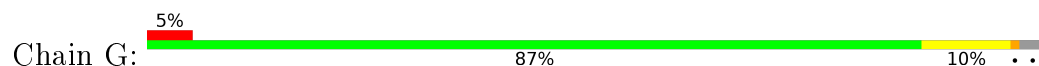
- Molecule 1: Insulin-like growth factor 1 receptor



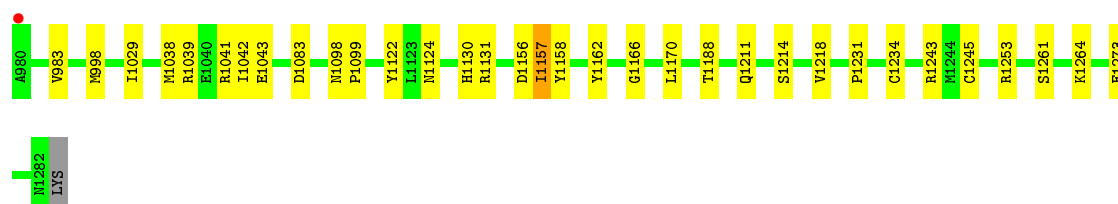
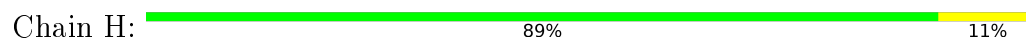
- Molecule 1: Insulin-like growth factor 1 receptor



- Molecule 1: Insulin-like growth factor 1 receptor



- Molecule 1: Insulin-like growth factor 1 receptor



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	111.44Å 190.04Å 155.49Å 90.00° 90.22° 90.00°	Depositor
Resolution (Å)	30.00 – 2.20 29.87 – 2.20	Depositor EDS
% Data completeness (in resolution range)	90.6 (30.00-2.20) 89.6 (29.87-2.20)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.98 (at 2.20Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.194 , 0.218 0.221 , 0.242	Depositor DCC
R_{free} test set	7400 reflections (5.01%)	DCC
Wilson B-factor (Å ²)	40.0	Xtriage
Anisotropy	0.088	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 25.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.011 for -1/2*h+1/2*k,3/2*h+1/2*k,-l 0.011 for -1/2*h-1/2*k,-3/2*h+1/2*k,-l 0.010 for 1/2*h+1/2*k,3/2*h-1/2*k,-l 0.009 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.467 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	19906	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.09% 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: SO4, MES, 66A

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	0.65	2/2409 (0.1%)	0.75	3/3250 (0.1%)
1	B	0.55	3/2409 (0.1%)	0.74	5/3250 (0.2%)
1	C	0.56	4/2409 (0.2%)	0.75	7/3250 (0.2%)
1	D	0.71	10/2467 (0.4%)	0.77	4/3332 (0.1%)
1	E	0.64	1/2409 (0.0%)	0.76	3/3250 (0.1%)
1	F	0.54	4/2409 (0.2%)	0.73	2/3250 (0.1%)
1	G	0.56	4/2409 (0.2%)	0.74	6/3250 (0.2%)
1	H	0.69	7/2458 (0.3%)	0.77	5/3321 (0.2%)
All	All	0.61	35/19379 (0.2%)	0.75	35/26153 (0.1%)

The worst 5 of 35 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	1243	ARG	CZ-NH1	8.16	1.43	1.33
1	C	1122	TYR	CE1-CZ	-7.29	1.29	1.38
1	G	1122	TYR	CE2-CZ	-7.22	1.29	1.38
1	B	1122	TYR	CE1-CZ	-6.54	1.30	1.38
1	D	1162	TYR	CE1-CZ	-6.39	1.30	1.38

The worst 5 of 35 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	1083	ASP	CB-CG-OD2	10.48	127.73	118.30
1	C	1083	ASP	CB-CG-OD2	9.90	127.21	118.30
1	B	1243	ARG	NE-CZ-NH2	9.44	125.02	120.30
1	F	1243	ARG	NE-CZ-NH2	8.83	124.72	120.30
1	D	1039	ARG	NE-CZ-NH2	8.64	124.62	120.30

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	2359	0	2318	15	0
1	B	2359	0	2318	31	0
1	C	2359	0	2318	17	0
1	D	2415	0	2377	19	0
1	E	2359	0	2318	17	0
1	F	2359	0	2318	32	0
1	G	2359	0	2318	15	0
1	H	2406	0	2364	17	0
2	A	33	0	0	0	0
2	B	33	0	0	0	0
2	C	33	0	0	0	0
2	D	33	0	0	0	0
2	E	33	0	0	0	0
2	F	33	0	0	0	0
2	G	33	0	0	0	0
2	H	33	0	0	0	0
3	A	12	0	12	0	0
3	B	12	0	12	0	0
3	C	12	0	12	1	0
3	D	12	0	12	0	0
3	E	12	0	12	0	0
3	F	12	0	12	0	0
3	G	12	0	12	1	0
3	H	12	0	12	0	0
4	A	5	0	0	0	0
4	C	5	0	0	0	0
4	E	5	0	0	0	0
4	G	5	0	0	0	0
5	A	92	0	0	1	0
5	B	42	0	0	0	0
5	C	49	0	0	3	0
5	D	92	0	0	1	0
5	E	97	0	0	1	0
5	F	43	0	0	1	0
5	G	48	0	0	2	0
5	H	88	0	0	1	0
All	All	19906	0	18745	153	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.

The worst 5 of 153 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:1124:ASN:OD1	1:G:1188:THR:HB	1.60	1.01
1:C:1124:ASN:OD1	1:C:1188:THR:HB	1.71	0.91
1:B:1003:GLY:H	1:B:1009:MET:HE3	1.41	0.85
1:F:1003:GLY:H	1:F:1009:MET:HE3	1.43	0.84
1:F:1219:LEU:HD21	1:F:1223:MET:HE1	1.64	0.80

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	292/304 (96%)	281 (96%)	11 (4%)	0	100	100
1	B	292/304 (96%)	280 (96%)	10 (3%)	2 (1%)	26	25
1	C	292/304 (96%)	277 (95%)	14 (5%)	1 (0%)	46	50
1	D	302/304 (99%)	294 (97%)	7 (2%)	1 (0%)	46	50
1	E	292/304 (96%)	282 (97%)	10 (3%)	0	100	100
1	F	292/304 (96%)	279 (96%)	11 (4%)	2 (1%)	26	25
1	G	292/304 (96%)	276 (94%)	15 (5%)	1 (0%)	46	50
1	H	301/304 (99%)	293 (97%)	7 (2%)	1 (0%)	46	50
All	All	2355/2432 (97%)	2262 (96%)	85 (4%)	8 (0%)	46	50

5 of 8 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	1023	PRO

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Mol	Chain	Res	Type
1	F	1023	PRO
1	C	1023	PRO
1	G	1023	PRO
1	D	1166	GLY

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	256/263 (97%)	252 (98%)	4 (2%)	70	82
1	B	256/263 (97%)	248 (97%)	8 (3%)	47	59
1	C	256/263 (97%)	252 (98%)	4 (2%)	70	82
1	D	262/263 (100%)	259 (99%)	3 (1%)	80	89
1	E	256/263 (97%)	251 (98%)	5 (2%)	63	76
1	F	256/263 (97%)	249 (97%)	7 (3%)	52	64
1	G	256/263 (97%)	252 (98%)	4 (2%)	70	82
1	H	261/263 (99%)	258 (99%)	3 (1%)	80	89
All	All	2059/2104 (98%)	2021 (98%)	38 (2%)	66	79

5 of 38 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	1211	GLN
1	E	1160	THR
1	H	983	VAL
1	E	1107	SER
1	E	1185	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 9 such sidechains are listed below:

Mol	Chain	Res	Type
1	E	1046	ASN

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Mol	Chain	Res	Type
1	H	1068	GLN
1	G	1033	ASN
1	C	1033	ASN
1	F	1046	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

20 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	66A	A	1301	-	32,38,38	1.67	3 (9%)	35,54,54	2.46	8 (22%)
3	MES	A	1302	-	12,12,12	1.76	1 (8%)	15,16,16	2.47	5 (33%)
4	SO4	A	1303	-	4,4,4	0.19	0	6,6,6	0.14	0
2	66A	B	1301	-	32,38,38	1.71	3 (9%)	35,54,54	2.13	4 (11%)
3	MES	B	1302	-	12,12,12	1.96	1 (8%)	15,16,16	2.50	6 (40%)
2	66A	C	1301	-	32,38,38	1.79	3 (9%)	35,54,54	2.33	4 (11%)
3	MES	C	1302	-	12,12,12	1.86	1 (8%)	15,16,16	2.78	5 (33%)
4	SO4	C	1303	-	4,4,4	0.13	0	6,6,6	0.27	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	66A	D	1301	-	32,38,38	1.80	3 (9%)	35,54,54	2.23	7 (20%)
3	MES	D	1302	-	12,12,12	1.65	1 (8%)	15,16,16	2.46	4 (26%)
2	66A	E	1301	-	32,38,38	1.62	3 (9%)	35,54,54	2.39	9 (25%)
3	MES	E	1302	-	12,12,12	1.74	1 (8%)	15,16,16	2.95	6 (40%)
4	SO4	E	1303	-	4,4,4	0.20	0	6,6,6	0.15	0
2	66A	F	1301	-	32,38,38	1.71	4 (12%)	35,54,54	2.04	4 (11%)
3	MES	F	1302	-	12,12,12	1.91	1 (8%)	15,16,16	2.82	5 (33%)
2	66A	G	1301	-	32,38,38	1.80	3 (9%)	35,54,54	2.36	5 (14%)
3	MES	G	1302	-	12,12,12	1.89	1 (8%)	15,16,16	2.82	5 (33%)
4	SO4	G	1303	-	4,4,4	0.13	0	6,6,6	0.31	0
2	66A	H	1301	-	32,38,38	1.82	3 (9%)	35,54,54	2.15	4 (11%)
3	MES	H	1302	-	12,12,12	1.76	1 (8%)	15,16,16	3.04	5 (33%)

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	66A	A	1301	-	-	0/11/31/31	0/6/6/6
3	MES	A	1302	-	-	0/6/14/14	0/1/1/1
4	SO4	A	1303	-	-	0/0/0/0	0/0/0/0
2	66A	B	1301	-	-	0/11/31/31	0/6/6/6
3	MES	B	1302	-	-	0/6/14/14	0/1/1/1
2	66A	C	1301	-	-	0/11/31/31	0/6/6/6
3	MES	C	1302	-	-	0/6/14/14	0/1/1/1
4	SO4	C	1303	-	-	0/0/0/0	0/0/0/0
2	66A	D	1301	-	-	0/11/31/31	0/6/6/6
3	MES	D	1302	-	-	0/6/14/14	0/1/1/1
2	66A	E	1301	-	-	0/11/31/31	0/6/6/6
3	MES	E	1302	-	-	0/6/14/14	0/1/1/1
4	SO4	E	1303	-	-	0/0/0/0	0/0/0/0
2	66A	F	1301	-	-	0/11/31/31	0/6/6/6
3	MES	F	1302	-	-	0/6/14/14	0/1/1/1
2	66A	G	1301	-	-	0/11/31/31	0/6/6/6
3	MES	G	1302	-	-	0/6/14/14	0/1/1/1
4	SO4	G	1303	-	-	0/0/0/0	0/0/0/0
2	66A	H	1301	-	-	0/11/31/31	0/6/6/6
3	MES	H	1302	-	-	0/6/14/14	0/1/1/1

The worst 5 of 33 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1302	MES	C8-S	-6.46	1.67	1.77
3	F	1302	MES	C8-S	-6.22	1.68	1.77
3	G	1302	MES	C8-S	-6.13	1.68	1.77
3	C	1302	MES	C8-S	-6.02	1.68	1.77
3	A	1302	MES	C8-S	-5.52	1.69	1.77

The worst 5 of 86 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1301	66A	N03-C04-N05	-9.59	121.33	128.87
2	G	1301	66A	N03-C04-N05	-9.34	121.54	128.87
2	H	1301	66A	N03-C04-N05	-9.20	121.65	128.87
2	A	1301	66A	N03-C04-N05	-9.05	121.76	128.87
2	C	1301	66A	N03-C04-N05	-8.98	121.82	128.87

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	1302	MES	1	0
3	G	1302	MES	1	0

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	296/304 (97%)	0.01	4 (1%) 78 77	28, 44, 62, 72	0
1	B	296/304 (97%)	0.26	21 (7%) 19 18	30, 46, 64, 74	0
1	C	296/304 (97%)	0.14	19 (6%) 23 22	29, 46, 65, 74	0
1	D	304/304 (100%)	-0.02	5 (1%) 74 73	27, 44, 63, 73	0
1	E	296/304 (97%)	0.08	3 (1%) 84 83	28, 44, 62, 73	0
1	F	296/304 (97%)	0.20	15 (5%) 32 31	30, 46, 64, 73	0
1	G	296/304 (97%)	0.07	14 (4%) 35 34	29, 45, 65, 74	0
1	H	303/304 (99%)	0.01	1 (0%) 94 94	27, 44, 62, 74	0
All	All	2383/2432 (97%)	0.09	82 (3%) 49 47	27, 45, 64, 74	0

The worst 5 of 82 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	1102	ALA	6.5
1	G	1102	ALA	5.5
1	B	1019	VAL	5.2
1	B	980	ALA	5.1
1	B	981	ALA	5.0

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 ⓘ

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
4	SO4	E	1303	5/5	0.87	0.14	1.05	87,87,89,90	0
2	66A	F	1301	33/33	0.92	0.15	0.69	39,43,51,52	0
4	SO4	A	1303	5/5	0.88	0.13	0.52	86,87,88,89	0
2	66A	B	1301	33/33	0.93	0.15	0.43	35,43,50,51	0
2	66A	D	1301	33/33	0.93	0.14	0.37	37,40,48,49	0
2	66A	H	1301	33/33	0.94	0.15	0.25	37,40,50,50	0
3	MES	F	1302	12/12	0.81	0.17	0.21	74,78,84,84	0
3	MES	B	1302	12/12	0.89	0.17	0.17	85,87,91,91	0
2	66A	G	1301	33/33	0.92	0.13	0.01	42,45,51,51	0
3	MES	C	1302	12/12	0.96	0.14	-0.01	52,59,69,69	0
2	66A	A	1301	33/33	0.94	0.13	-0.13	31,36,48,49	0
3	MES	G	1302	12/12	0.94	0.14	-0.15	53,59,71,71	0
2	66A	C	1301	33/33	0.89	0.13	-0.19	43,45,53,53	0
3	MES	D	1302	12/12	0.91	0.12	-0.30	52,57,67,68	0
2	66A	E	1301	33/33	0.95	0.14	-0.32	30,35,48,49	0
3	MES	E	1302	12/12	0.94	0.13	-0.52	49,55,60,61	0
3	MES	H	1302	12/12	0.93	0.12	-0.55	52,56,69,69	0
4	SO4	C	1303	5/5	0.97	0.11	-0.57	64,65,67,67	0
3	MES	A	1302	12/12	0.94	0.13	-0.59	45,50,59,60	0
4	SO4	G	1303	5/5	0.98	0.08	-1.83	61,61,61,62	0

6.5 Other polymers [i](#)

There are no such residues in this entry.