



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 19, 2017 – 07:55 AM EST

PDB ID : 5H25  
Title : EED in complex with PRC2 allosteric inhibitor compound 11  
Authors : Zhao, K.; Zhao, M.; Luo, X.; Zhang, H.  
Deposited on : 2016-10-14  
Resolution : 2.88 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](mailto: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

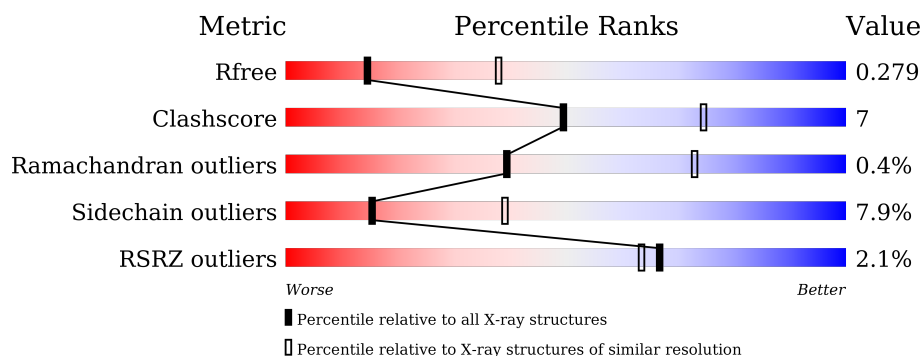
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.88 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	1945 (2.90-2.86)
Clashscore	102246	2202 (2.90-2.86)
Ramachandran outliers	100387	2149 (2.90-2.86)
Sidechain outliers	100360	2152 (2.90-2.86)
RSRZ outliers	91569	1950 (2.90-2.86)

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  $\geq 3$ , 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	367	<div> <div>2%</div> <div> <div></div> <div>74%</div> <div>23%</div> <div>..</div> </div> </div>
1	B	367	<div> <div>%</div> <div> <div></div> <div>77%</div> <div>18%</div> <div>..</div> </div> </div>
2	C	29	<div> <div>3%</div> <div> <div></div> <div>69%</div> <div>28%</div> <div>.</div> </div> </div>
2	D	29	<div> <div>3%</div> <div> <div></div> <div>83%</div> <div>17%</div> </div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6390 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 Polycomb protein EED.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	364	Total	C	N	O	S	0	0	0
			2948	1867	517	542	22			
1	B	357	Total	C	N	O	S	0	0	0
			2890	1834	503	531	22			

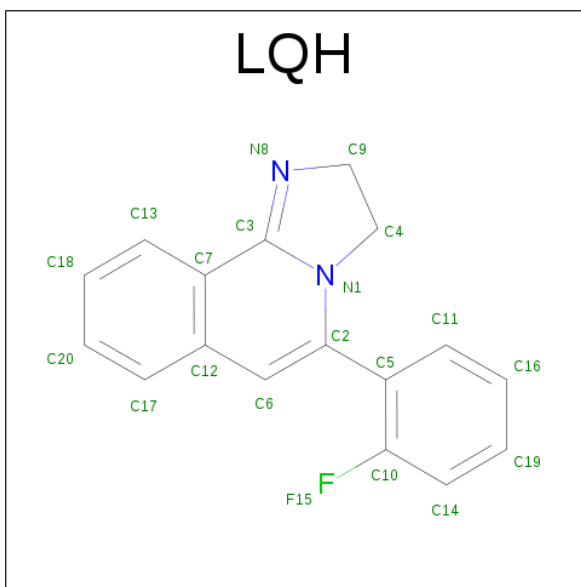
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	75	GLY	-	expression tag	UNP O75530
B	75	GLY	-	expression tag	UNP O75530

- Molecule 2 is a protein called Histone-lysine N-methyltransferase EZH2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	29	Total	C	N	O	S	0	0	0
			254	157	50	46	1			
2	D	29	Total	C	N	O	S	0	0	0
			254	157	50	46	1			

- Molecule 3 is 5-(2-fluorophenyl)-2,3-dihydroimidazo[2,1-a]isoquinoline (three-letter code: LQH) (formula: C<sub>17</sub>H<sub>13</sub>FN<sub>2</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	F	N	0	0
			20	17	1	2		
3	B	1	Total	C	F	N	0	0
			20	17	1	2		

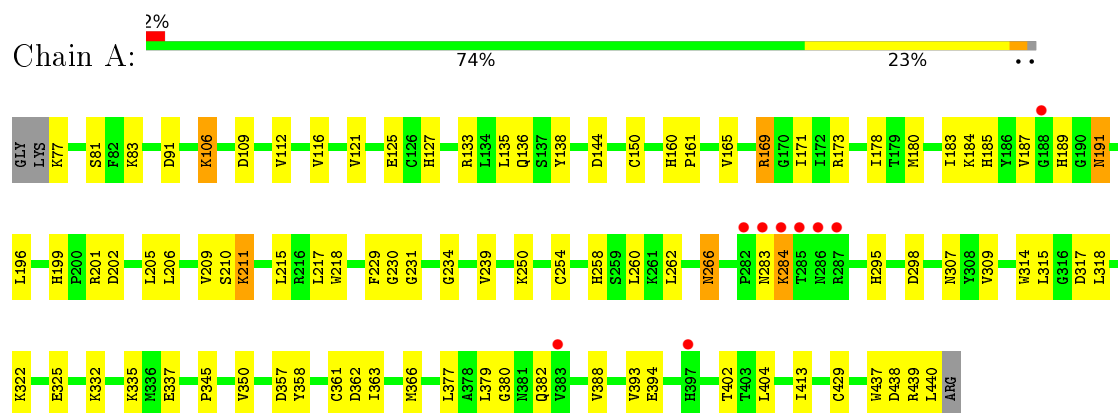
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	O	0	0
			2	2		
4	B	2	Total	O	0	0
			2	2		

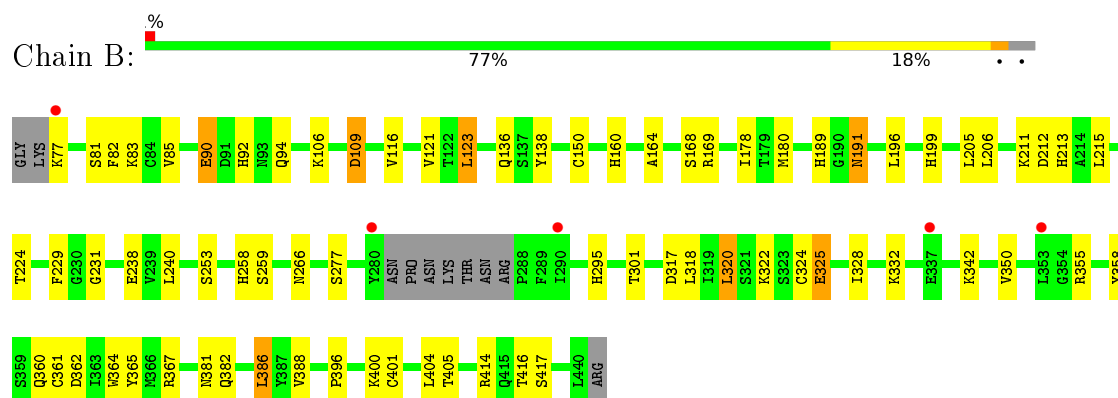
### 3 Residue-property plots [i](#)

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

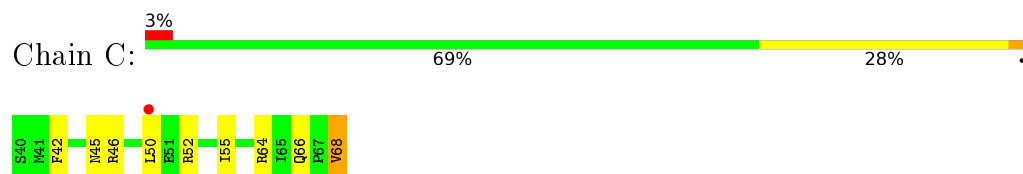
- Molecule 1: Polycomb protein EED



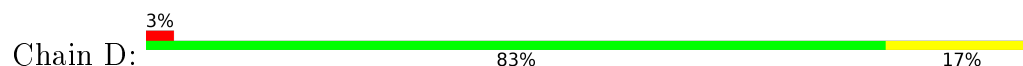
- Molecule 1: Polycomb protein EED

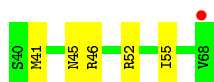


- Molecule 2: Histone-lysine N-methyltransferase EZH2



- Molecule 2: Histone-lysine N-methyltransferase EZH2





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 2 <sub>1</sub> 2 <sub>1</sub> 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	94.75 Å   179.38 Å   50.40 Å 90.00°   90.00°   90.00°	Depositor
Resolution (Å)	43.94 – 2.88 43.94 – 2.88	Depositor EDS
% Data completeness (in resolution range)	98.6 (43.94-2.88) 98.8 (43.94-2.88)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.46 (at 2.90 Å)	Xtriage
Refinement program	BUSTER	Depositor
R, $R_{free}$	0.184   ,   0.274 0.186   ,   0.279	Depositor DCC
$R_{free}$ test set	1017 reflections (5.38%)	DCC
Wilson B-factor (Å <sup>2</sup> )	46.9	Xtriage
Anisotropy	0.753	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 49.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	6390	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 32.95 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 8.6488e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup> Intensities estimated from amplitudes.

<sup>2</sup> Theoretical 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: LQH

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	0.51	0/3023	0.75	0/4093
1	B	0.50	0/2963	0.73	0/4009
2	C	0.58	0/257	0.71	0/341
2	D	0.54	0/257	0.64	0/341
All	All	0.51	0/6500	0.73	0/8784

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	2948	0	2866	50	0
1	B	2890	0	2808	37	0
2	C	254	0	260	5	0
2	D	254	0	260	3	0
3	A	20	0	0	0	0
3	B	20	0	0	0	0
4	A	2	0	0	0	0
4	B	2	0	0	0	0
All	All	6390	0	6194	89	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 7.

All (89) 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:199:HIS:HD2	1:A:201:ARG:H	1.11	0.95
1:A:215:LEU:HB2	1:A:229:PHE:HB2	1.65	0.78
1:A:125:GLU:HG3	1:A:135:LEU:HD11	1.70	0.74
1:B:317:ASP:H	2:D:45:ASN:HD21	1.34	0.73
1:A:199:HIS:CD2	1:A:201:ARG:H	2.01	0.73
1:A:358:TYR:HE1	1:A:361:CYS:HB3	1.54	0.72
1:B:191:ASN:HB3	1:B:211:LYS:HB2	1.72	0.69
1:B:92:HIS:HD2	1:B:94:GLN:H	1.40	0.69
1:A:317:ASP:H	2:C:45:ASN:HD21	1.46	0.63
1:A:127:HIS:CD2	1:A:133:ARG:HE	2.19	0.61
1:A:250:LYS:HE2	1:A:314:TRP:CH2	2.40	0.57
1:B:215:LEU:HB2	1:B:229:PHE:HB2	1.86	0.57
1:A:266:ASN:H	1:A:266:ASN:HD22	1.53	0.56
1:B:358:TYR:CE1	1:B:361:CYS:HB3	2.40	0.56
1:A:136:GLN:CB	1:A:180:MET:HG3	2.36	0.56
1:A:189:HIS:CE1	1:A:210:SER:HB2	2.40	0.55
1:B:360:GLN:O	1:B:381:ASN:HB2	2.06	0.55
1:B:92:HIS:CD2	1:B:94:GLN:H	2.24	0.55
1:A:106:LYS:O	1:A:109:ASP:HB2	2.06	0.54
1:A:191:ASN:HB3	1:A:211:LYS:HB3	1.89	0.54
1:A:77:LYS:HE3	1:A:393:VAL:HG22	1.88	0.54
1:A:136:GLN:HB3	1:A:180:MET:HG3	1.90	0.54
1:B:386:LEU:HD21	1:B:416:THR:HG21	1.90	0.54
1:A:83:LYS:HB2	1:A:440:LEU:HD21	1.89	0.54
1:A:160:HIS:CE1	1:A:178:ILE:HD12	2.43	0.53
1:A:298:ASP:O	1:A:345:PRO:HA	2.08	0.53
2:C:52:ARG:HA	2:C:55:ILE:HD12	1.91	0.53
1:A:358:TYR:CE1	1:A:361:CYS:HB3	2.39	0.53
1:B:231:GLY:HA3	1:B:295:HIS:ND1	2.23	0.53
1:B:358:TYR:HE1	1:B:361:CYS:HB3	1.73	0.53
1:A:262:LEU:HD23	1:A:298:ASP:HB2	1.91	0.52
1:B:138:TYR:HB2	1:B:180:MET:HG3	1.92	0.51
1:A:199:HIS:CD2	1:A:202:ASP:H	2.29	0.50
1:A:209:VAL:HG12	1:A:215:LEU:HG	1.94	0.50
1:A:199:HIS:HD2	1:A:201:ARG:N	1.94	0.50
1:B:388:VAL:HG13	1:B:401:CYS:HB2	1.94	0.50
1:B:82:PHE:HB2	1:B:404:LEU:HD22	1.94	0.50

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:324:CYS:HB2	1:B:362:ASP:O	2.12	0.49
1:A:363:ILE:HG12	1:A:382:GLN:HG2	1.94	0.49
1:B:317:ASP:H	2:D:45:ASN:ND2	2.07	0.48
1:B:259:SER:HA	1:B:301:THR:O	2.13	0.48
2:D:52:ARG:HA	2:D:55:ILE:HD12	1.94	0.48
1:B:324:CYS:HB3	1:B:364:TRP:CH2	2.48	0.47
1:B:160:HIS:CE1	1:B:178:ILE:HD12	2.50	0.47
1:A:380:GLY:HA3	1:A:413:ILE:HB	1.97	0.47
1:B:215:LEU:HD21	1:B:253:SER:HB3	1.97	0.46
1:B:199:HIS:HB2	1:B:205:LEU:HB2	1.97	0.46
1:B:106:LYS:O	1:B:109:ASP:HB2	2.16	0.45
1:A:121:VAL:HB	1:A:138:TYR:HB3	1.97	0.45
1:A:231:GLY:HA3	1:A:295:HIS:ND1	2.31	0.45
1:A:322:LYS:NZ	1:A:366:MET:HB2	2.32	0.45
1:A:191:ASN:HB3	1:A:211:LYS:CB	2.46	0.44
1:A:205:LEU:HA	1:A:218:TRP:O	2.17	0.44
1:B:355:ARG:O	1:B:396:PRO:HB2	2.18	0.44
1:A:116:VAL:HG12	1:A:121:VAL:HG22	2.00	0.44
1:A:230:GLY:HA2	1:A:234:GLY:HA3	1.98	0.44
1:B:322:LYS:HB3	1:B:328:ILE:HG12	1.98	0.44
1:B:121:VAL:HB	1:B:138:TYR:HB3	2.00	0.43
1:B:90:GLU:HB3	1:B:92:HIS:CE1	2.53	0.43
1:A:254:CYS:HB2	1:A:309:VAL:HB	2.00	0.43
1:A:136:GLN:HB2	1:A:180:MET:HG3	2.01	0.43
1:A:209:VAL:HB	1:A:239:VAL:HB	2.01	0.42
1:A:169:ARG:HB3	1:A:171:ILE:HD12	2.01	0.42
1:A:106:LYS:HG3	1:A:106:LYS:H	1.68	0.42
1:A:161:PRO:HD3	2:C:66:GLN:HB2	2.01	0.42
1:B:365:TYR:HB2	1:B:414:ARG:HH21	1.84	0.42
1:B:150:CYS:HA	1:B:164:ALA:O	2.20	0.42
1:B:196:LEU:HD22	1:B:206:LEU:HD11	2.01	0.42
1:B:224:THR:OG1	1:B:277:SER:HB2	2.20	0.42
1:A:413:ILE:HG12	1:A:429:CYS:SG	2.60	0.42
1:A:377:LEU:O	1:A:388:VAL:HA	2.20	0.41
1:B:320:LEU:HA	1:B:320:LEU:HD12	1.91	0.41
1:A:112:VAL:HG21	2:C:68:VAL:HG21	2.01	0.41
1:A:196:LEU:HD22	1:A:206:LEU:HD11	2.02	0.41
1:B:92:HIS:CD2	1:B:94:GLN:HB2	2.56	0.41
1:A:315:LEU:O	1:A:318:LEU:HB2	2.20	0.41
1:B:123:LEU:HD22	1:B:180:MET:HE1	2.02	0.41
1:B:211:LYS:O	1:B:213:HIS:HD2	2.04	0.41

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:240:LEU:HD22	1:B:365:TYR:HE1	1.85	0.41
1:A:205:LEU:HB3	1:A:217:LEU:HD11	2.02	0.41
1:A:382:GLN:HB2	1:A:382:GLN:HE21	1.66	0.41
1:B:211:LYS:HA	1:B:238:GLU:HG2	2.03	0.41
1:B:92:HIS:HD2	1:B:94:GLN:N	2.12	0.41
1:A:394:GLU:HA	2:C:42:PHE:CE2	2.56	0.41
1:B:83:LYS:HD3	1:B:85:VAL:HG12	2.02	0.40
1:A:183:ILE:HG13	1:A:184:LYS:HG2	2.03	0.40
1:A:258:HIS:CE1	1:A:307:ASN:HA	2.56	0.40
1:A:404:LEU:HB3	1:A:437:TRP:CZ3	2.56	0.40
1:A:150:CYS:HB3	1:A:165:VAL:HG12	2.04	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	362/367 (99%)	343 (95%)	17 (5%)	2 (1%)	30	65
1	B	353/367 (96%)	334 (95%)	18 (5%)	1 (0%)	46	78
2	C	27/29 (93%)	25 (93%)	2 (7%)	0	100	100
2	D	27/29 (93%)	25 (93%)	2 (7%)	0	100	100
All	All	769/792 (97%)	727 (94%)	39 (5%)	3 (0%)	39	73

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	284	LYS
1	A	325	GLU
1	B	325	GLU

### 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	327/329 (99%)	303 (93%)	24 (7%)	17	43
1	B	320/329 (97%)	294 (92%)	26 (8%)	15	38
2	C	29/29 (100%)	25 (86%)	4 (14%)	4	11
2	D	29/29 (100%)	27 (93%)	2 (7%)	19	46
All	All	705/716 (98%)	649 (92%)	56 (8%)	15	39

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	81	SER
1	A	91	ASP
1	A	106	LYS
1	A	144	ASP
1	A	169	ARG
1	A	173	ARG
1	A	185	HIS
1	A	187	VAL
1	A	191	ASN
1	A	211	LYS
1	A	260	LEU
1	A	266	ASN
1	A	283	ASN
1	A	284	LYS
1	A	332	LYS
1	A	335	LYS
1	A	337	GLU
1	A	350	VAL
1	A	357	ASP
1	A	362	ASP
1	A	379	LEU
1	A	402	THR
1	A	438	ASP
1	A	439	ARG

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	B	77	LYS
1	B	81	SER
1	B	90	GLU
1	B	109	ASP
1	B	116	VAL
1	B	123	LEU
1	B	136	GLN
1	B	168	SER
1	B	169	ARG
1	B	189	HIS
1	B	191	ASN
1	B	212	ASP
1	B	258	HIS
1	B	266	ASN
1	B	318	LEU
1	B	320	LEU
1	B	325	GLU
1	B	332	LYS
1	B	342	LYS
1	B	350	VAL
1	B	367	ARG
1	B	382	GLN
1	B	386	LEU
1	B	400	LYS
1	B	405	THR
1	B	417	SER
2	C	46	ARG
2	C	50	LEU
2	C	64	ARG
2	C	68	VAL
2	D	41	MET
2	D	46	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	100	GLN
1	A	199	HIS
1	A	213	HIS
1	A	258	HIS
1	A	266	ASN
1	B	92	HIS

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	B	100	GLN
1	B	213	HIS
1	B	266	ASN
1	B	415	GLN
2	C	45	ASN
2	C	58	GLN
2	D	45	ASN

### 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 ⓘ

2 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	LQH	A	501	-	21,23,23	2.30	7 (33%)	26,33,33	1.61	5 (19%)
3	LQH	B	501	-	21,23,23	2.32	5 (23%)	26,33,33	2.01	7 (26%)

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
3	LQH	A	501	-	-	0/4/10/10	0/3/4/4
3	LQH	B	501	-	-	0/4/10/10	0/3/4/4

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	501	LQH	C2-N1	-2.66	1.33	1.37
3	A	501	LQH	C2-N1	-2.52	1.33	1.37
3	A	501	LQH	C18-C13	2.00	1.41	1.36
3	A	501	LQH	C20-C17	2.40	1.42	1.36
3	A	501	LQH	C5-C10	3.00	1.42	1.38
3	B	501	LQH	C5-C2	3.13	1.53	1.48
3	A	501	LQH	C5-C2	3.28	1.53	1.48
3	B	501	LQH	C5-C10	3.44	1.43	1.38
3	A	501	LQH	C6-C12	3.93	1.51	1.42
3	B	501	LQH	C6-C12	4.15	1.51	1.42
3	A	501	LQH	C3-C7	6.78	1.52	1.41
3	B	501	LQH	C3-C7	6.80	1.52	1.41

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	501	LQH	C9-C4-N1	-5.52	98.18	102.37
3	A	501	LQH	C9-C4-N1	-3.63	99.62	102.37
3	B	501	LQH	C14-C10-C5	-3.18	120.13	123.44
3	B	501	LQH	C6-C2-C5	-2.80	115.01	120.24
3	A	501	LQH	C14-C10-C5	-2.63	120.70	123.44
3	B	501	LQH	C6-C12-C7	-2.42	116.87	119.83
3	A	501	LQH	C6-C12-C7	-2.16	117.19	119.83
3	A	501	LQH	C11-C5-C10	3.19	119.46	115.99
3	B	501	LQH	C11-C5-C10	3.36	119.65	115.99
3	B	501	LQH	C5-C2-N1	3.45	124.74	120.65
3	B	501	LQH	C3-N1-C2	4.10	123.69	118.05
3	A	501	LQH	C3-N1-C2	4.30	123.96	118.05

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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	364/367 (99%)	-0.14	9 (2%) 61 57	24, 44, 70, 103	0
1	B	357/367 (97%)	-0.15	5 (1%) 78 76	29, 47, 72, 83	0
2	C	29/29 (100%)	0.28	1 (3%) 49 42	38, 56, 107, 114	0
2	D	29/29 (100%)	0.29	1 (3%) 49 42	40, 58, 83, 92	0
All	All	779/792 (98%)	-0.11	16 (2%) 67 63	24, 46, 73, 114	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	283	ASN	4.0
1	A	286	ASN	3.4
1	A	287	ARG	3.3
2	D	68	VAL	3.3
1	B	280	TYR	3.2
1	A	383	VAL	3.2
1	A	188	GLY	3.0
1	A	285	THR	3.0
1	A	284	LYS	2.8
1	B	290	ILE	2.8
1	A	397	HIS	2.8
2	C	50	LEU	2.5
1	A	282	PRO	2.4
1	B	77	LYS	2.2
1	B	353	LEU	2.1
1	B	337	GLU	2.0

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

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

### 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

### 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
3	LQH	A	501	20/20	0.93	0.19	0.75	44,52,60,62	0
3	LQH	B	501	20/20	0.94	0.17	-0.32	49,57,64,65	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.