



# Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 12:17 am BST

PDB ID : 5FGL  
Title : Co-crystal Structure of NicR2\_Hsp  
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Deposited on : 2015-12-21  
Resolution : 2.40 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

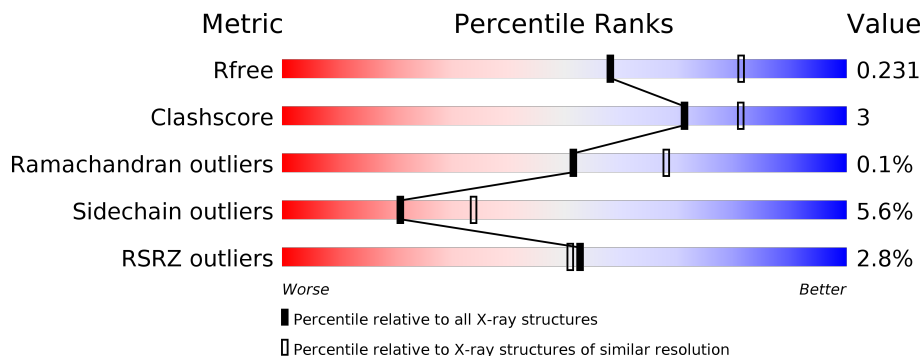
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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))
$R_{free}$	130704	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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 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	208	 4% 86% 8% ..
1	B	208	 % 88% 9% ..
1	C	208	 3% 87% 10% ..
1	D	208	 3% 85% 11% ..
1	E	208	 87% 10% ..
1	F	208	 4% 87% 8% ..

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 10186 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.

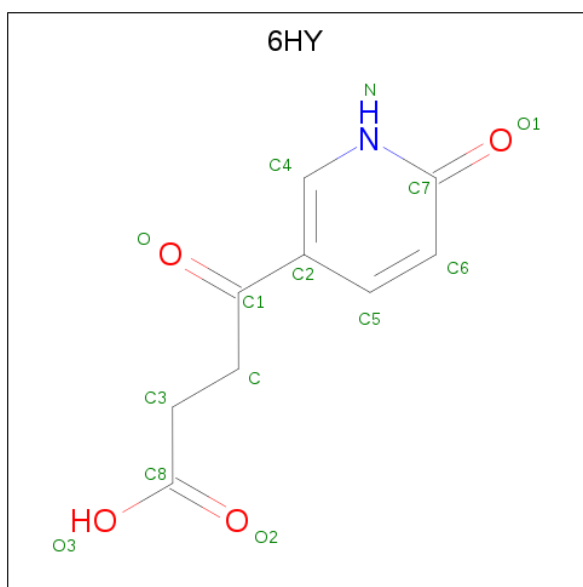
- Molecule 1 is a protein called NicR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	200	Total 1629	C 1036	N 277	O 306	S 10	0	1	0
1	B	206	Total 1661	C 1050	N 284	O 317	S 10	0	0	0
1	C	205	Total 1670	C 1058	N 287	O 315	S 10	0	0	0
1	D	203	Total 1649	C 1043	N 284	O 312	S 10	0	0	0
1	E	206	Total 1674	C 1060	N 287	O 317	S 10	0	1	0
1	F	201	Total 1631	C 1037	N 278	O 307	S 9	0	1	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	30	MET	-	initiating methionine	UNP A0A0B4KIF6
A	124	ALA	GLY	engineered mutation	UNP A0A0B4KIF6
B	30	MET	-	initiating methionine	UNP A0A0B4KIF6
B	124	ALA	GLY	engineered mutation	UNP A0A0B4KIF6
C	30	MET	-	initiating methionine	UNP A0A0B4KIF6
C	124	ALA	GLY	engineered mutation	UNP A0A0B4KIF6
D	30	MET	-	initiating methionine	UNP A0A0B4KIF6
D	124	ALA	GLY	engineered mutation	UNP A0A0B4KIF6
E	30	MET	-	initiating methionine	UNP A0A0B4KIF6
E	124	ALA	GLY	engineered mutation	UNP A0A0B4KIF6
F	30	MET	-	initiating methionine	UNP A0A0B4KIF6
F	124	ALA	GLY	engineered mutation	UNP A0A0B4KIF6

- Molecule 2 is 4-oxidanylidene-4-(6-oxidanylidene-1 {H}-pyridin-3-yl)butanoic acid (three-letter code: 6HY) (formula: C<sub>9</sub>H<sub>9</sub>NO<sub>4</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	Total 14	C 9	N 1	O 4	0	0
2	B	1	Total 14	C 9	N 1	O 4	0	0
2	C	1	Total 14	C 9	N 1	O 4	0	0
2	D	1	Total 14	C 9	N 1	O 4	0	0
2	E	1	Total 14	C 9	N 1	O 4	0	0
2	E	1	Total 14	C 9	N 1	O 4	0	0

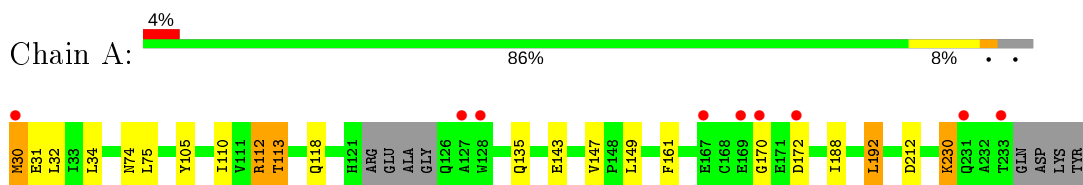
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	23	Total 23	O 23	0	0
3	B	36	Total 36	O 36	0	0
3	C	33	Total 33	O 33	0	0
3	D	35	Total 35	O 35	0	0
3	E	38	Total 38	O 38	0	0
3	F	23	Total 23	O 23	0	0

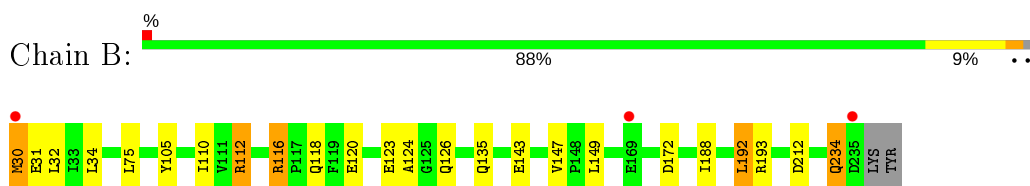
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

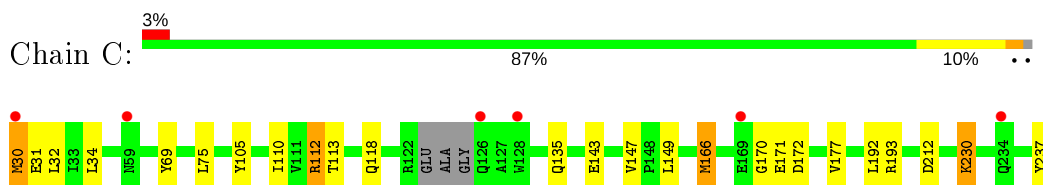
- Molecule 1: NicR



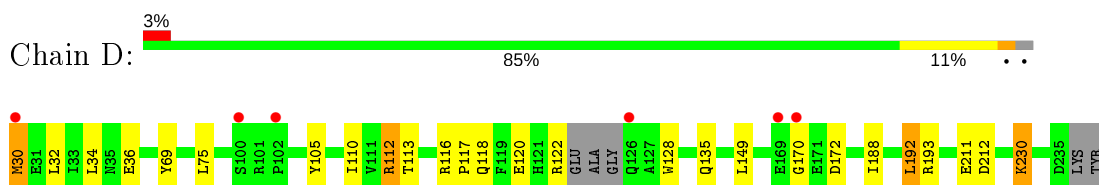
- Molecule 1: NicR



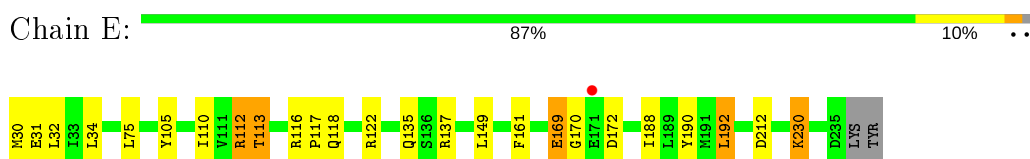
- Molecule 1: NicR



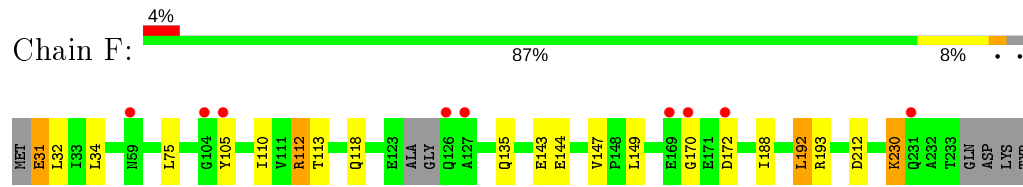
- Molecule 1: NicR



- Molecule 1: NicR



- Molecule 1: NicR



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	149.52Å 149.52Å 156.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.40 47.48 – 2.40	Depositor EDS
% Data completeness (in resolution range)	97.3 (50.00-2.40) 97.3 (47.48-2.40)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.69 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, $R_{free}$	0.200 , 0.229 0.206 , 0.231	Depositor DCC
$R_{free}$ test set	3426 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	46.7	Xtriage
Anisotropy	0.012	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 38.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.013 for -h,l,k 0.007 for -l,-k,-h	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	10186	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.06% of the height of the origin peak. No significant pseudotranslation is detected.*

<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: 6HY

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	0.46	0/1664	0.65	1/2243 (0.0%)
1	B	0.55	0/1693	0.76	3/2284 (0.1%)
1	C	0.49	0/1702	0.68	1/2293 (0.0%)
1	D	0.47	0/1680	0.70	1/2264 (0.0%)
1	E	0.50	0/1710	0.70	0/2305
1	F	0.47	0/1666	0.67	1/2247 (0.0%)
All	All	0.49	0/10115	0.69	7/13636 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	116	ARG	NE-CZ-NH1	8.12	124.36	120.30
1	D	193	ARG	CG-CD-NE	7.30	127.12	111.80
1	F	193	ARG	CG-CD-NE	7.26	127.05	111.80
1	B	116	ARG	NE-CZ-NH2	-6.46	117.07	120.30
1	B	193	ARG	NE-CZ-NH2	6.03	123.32	120.30
1	C	193	ARG	NE-CZ-NH2	5.42	123.01	120.30
1	A	30	MET	CG-SD-CE	5.14	108.43	100.20

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	1629	0	1597	10	0
1	B	1661	0	1617	9	0
1	C	1670	0	1635	13	0
1	D	1649	0	1613	14	0
1	E	1674	0	1633	14	0
1	F	1631	0	1592	10	0
2	A	14	0	0	0	0
2	B	14	0	0	0	0
2	C	14	0	0	0	0
2	D	14	0	0	0	0
2	E	28	0	0	1	0
3	A	23	0	0	1	0
3	B	36	0	0	0	0
3	C	33	0	0	0	0
3	D	35	0	0	1	0
3	E	38	0	0	0	0
3	F	23	0	0	1	0
All	All	10186	0	9687	68	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 3.

All (68) 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:C:166:MET:HA	1:C:166:MET:CE	1.90	1.00
1:C:166:MET:HE2	1:C:166:MET:HA	1.54	0.89
1:B:116:ARG:HD2	1:B:120:GLU:OE1	1.86	0.76
1:D:128:TRP:HB3	3:D:428:HOH:O	1.86	0.75
1:A:118:GLN:HE22	1:A:135:GLN:HE22	1.37	0.73
1:D:118:GLN:HE22	1:D:135:GLN:HE22	1.37	0.72
1:C:118:GLN:HE22	1:C:135:GLN:HE22	1.38	0.71
1:B:234:GLN:HA	1:B:234:GLN:OE1	1.92	0.70
1:B:118:GLN:HE22	1:B:135:GLN:HE22	1.36	0.70
1:C:166:MET:HE3	1:C:166:MET:HA	1.71	0.69
1:F:118:GLN:HE22	1:F:135:GLN:HE22	1.38	0.69
1:E:118:GLN:HE22	1:E:135:GLN:HE22	1.42	0.68
1:C:166:MET:CE	1:C:171:GLU:HB3	2.25	0.67
1:D:36:GLU:OE2	1:F:144:GLU:HG3	1.96	0.66
1:C:112:ARG:NH2	1:C:212:ASP:OD1	2.29	0.66
1:C:30:MET:HB3	1:C:69:TYR:OH	1.96	0.66
1:B:112:ARG:NH2	1:B:212:ASP:OD1	2.29	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:112:ARG:NH2	1:A:212:ASP:OD1	2.30	0.65
1:F:112:ARG:NH2	1:F:212:ASP:OD1	2.30	0.65
1:D:30:MET:HB3	1:D:69:TYR:OH	1.96	0.64
1:E:112:ARG:NH2	1:E:212:ASP:OD1	2.31	0.64
1:F:31:GLU:N	3:F:301:HOH:O	2.32	0.62
1:D:112:ARG:NH2	1:D:212:ASP:OD1	2.32	0.62
1:C:166:MET:CE	1:C:166:MET:CA	2.77	0.55
1:A:113:THR:HG22	1:A:161[B]:PHE:HE2	1.74	0.51
1:E:113:THR:HG22	1:E:161[B]:PHE:HE2	1.76	0.51
1:E:137:ARG:HB3	1:E:137:ARG:NH2	2.26	0.50
1:C:149:LEU:C	1:C:149:LEU:HD23	2.32	0.50
1:E:149:LEU:HD23	1:E:149:LEU:C	2.32	0.49
1:E:105:TYR:CE2	1:E:110:ILE:HD11	2.48	0.48
1:D:149:LEU:HD23	1:D:149:LEU:C	2.34	0.48
1:D:105:TYR:CE2	1:D:110:ILE:HD11	2.49	0.47
1:F:149:LEU:C	1:F:149:LEU:HD23	2.33	0.47
1:B:30:MET:O	1:B:31:GLU:OE2	2.33	0.47
1:D:120:GLU:HG2	1:D:211:GLU:OE2	2.15	0.47
1:A:105:TYR:CE2	1:A:110:ILE:HD11	2.51	0.46
1:A:149:LEU:C	1:A:149:LEU:HD23	2.35	0.46
1:A:172:ASP:OD1	1:A:172:ASP:N	2.49	0.46
1:B:149:LEU:C	1:B:149:LEU:HD23	2.36	0.45
1:C:166:MET:CE	1:C:177:VAL:HG21	2.47	0.45
1:E:30:MET:O	1:E:31:GLU:OE2	2.35	0.45
1:A:188:ILE:HG22	1:A:192:LEU:HD22	1.99	0.44
1:B:105:TYR:CE2	1:B:110:ILE:HD11	2.52	0.44
1:D:172:ASP:OD1	1:D:172:ASP:N	2.50	0.44
1:D:116:ARG:HB2	1:D:117:PRO:HD3	1.99	0.44
1:F:188:ILE:HG22	1:F:192:LEU:HD22	2.00	0.43
1:A:143:GLU:HG3	1:A:147:VAL:HG23	2.01	0.43
1:A:170:GLY:O	1:A:230:LYS:HE2	2.18	0.43
1:E:116:ARG:HB2	1:E:117:PRO:HD3	1.99	0.43
1:E:190:TYR:OH	2:E:302:6HY:O2	2.30	0.43
1:D:170:GLY:O	1:D:230:LYS:HE2	2.19	0.43
1:C:143:GLU:HG3	1:C:147:VAL:HG23	2.00	0.43
1:E:105:TYR:HE2	1:E:110:ILE:HD11	1.84	0.43
1:E:170:GLY:O	1:E:230:LYS:HE2	2.18	0.43
1:F:170:GLY:O	1:F:230:LYS:HE2	2.19	0.42
1:C:170:GLY:O	1:C:230:LYS:HE2	2.18	0.42
1:D:105:TYR:HE2	1:D:110:ILE:HD11	1.85	0.42
1:B:143:GLU:HG3	1:B:147:VAL:HG23	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:36:GLU:CD	1:F:144:GLU:HG3	2.41	0.41
1:D:188:ILE:HG22	1:D:192:LEU:HD22	2.02	0.41
1:A:74:ASN:ND2	3:A:402:HOH:O	2.52	0.41
1:E:169:GLU:CD	1:E:169:GLU:H	2.24	0.41
1:F:143:GLU:HG3	1:F:147:VAL:HG23	2.02	0.41
1:B:188:ILE:HG22	1:B:192:LEU:HD22	2.03	0.41
1:E:137:ARG:HB3	1:E:137:ARG:HH21	1.86	0.40
1:E:188:ILE:HG22	1:E:192:LEU:HD22	2.02	0.40
1:F:105:TYR:CE1	1:F:110:ILE:HD11	2.57	0.40
1:C:105:TYR:CE1	1:C:110:ILE:HD11	2.57	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	197/208 (95%)	195 (99%)	2 (1%)	0	100	100
1	B	204/208 (98%)	199 (98%)	4 (2%)	1 (0%)	29	41
1	C	201/208 (97%)	200 (100%)	1 (0%)	0	100	100
1	D	199/208 (96%)	197 (99%)	2 (1%)	0	100	100
1	E	205/208 (99%)	200 (98%)	5 (2%)	0	100	100
1	F	198/208 (95%)	196 (99%)	2 (1%)	0	100	100
All	All	1204/1248 (96%)	1187 (99%)	16 (1%)	1 (0%)	51	68

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	124	ALA

### 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	173/178 (97%)	164 (95%)	9 (5%)	23	38
1	B	175/178 (98%)	165 (94%)	10 (6%)	20	33
1	C	177/178 (99%)	165 (93%)	12 (7%)	16	25
1	D	175/178 (98%)	166 (95%)	9 (5%)	24	39
1	E	177/178 (99%)	167 (94%)	10 (6%)	21	34
1	F	172/178 (97%)	163 (95%)	9 (5%)	23	38
All	All	1049/1068 (98%)	990 (94%)	59 (6%)	21	34

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	MET
1	A	31	GLU
1	A	32	LEU
1	A	34	LEU
1	A	75	LEU
1	A	112	ARG
1	A	113	THR
1	A	192	LEU
1	A	230	LYS
1	B	30	MET
1	B	32	LEU
1	B	34	LEU
1	B	75	LEU
1	B	112	ARG
1	B	123	GLU
1	B	126	GLN
1	B	172	ASP
1	B	192	LEU
1	B	234	GLN
1	C	30	MET
1	C	31	GLU
1	C	32	LEU

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Mol	Chain	Res	Type
1	C	34	LEU
1	C	75	LEU
1	C	112	ARG
1	C	113	THR
1	C	166	MET
1	C	172	ASP
1	C	192	LEU
1	C	230	LYS
1	C	237	TYR
1	D	30	MET
1	D	32	LEU
1	D	34	LEU
1	D	75	LEU
1	D	112	ARG
1	D	113	THR
1	D	122	ARG
1	D	192	LEU
1	D	230	LYS
1	E	32	LEU
1	E	34	LEU
1	E	75	LEU
1	E	112	ARG
1	E	113	THR
1	E	122	ARG
1	E	169	GLU
1	E	172	ASP
1	E	192	LEU
1	E	230	LYS
1	F	31	GLU
1	F	32	LEU
1	F	34	LEU
1	F	75	LEU
1	F	112	ARG
1	F	113	THR
1	F	172	ASP
1	F	192	LEU
1	F	230	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	ASN

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Mol	Chain	Res	Type
1	A	62	GLN
1	A	93	ASN
1	A	135	GLN
1	B	57	ASN
1	B	62	GLN
1	B	93	ASN
1	B	135	GLN
1	C	57	ASN
1	C	62	GLN
1	C	93	ASN
1	C	135	GLN
1	D	57	ASN
1	D	62	GLN
1	D	93	ASN
1	D	126	GLN
1	D	135	GLN
1	E	57	ASN
1	E	62	GLN
1	E	74	ASN
1	E	93	ASN
1	E	135	GLN
1	F	57	ASN
1	F	62	GLN
1	F	93	ASN
1	F	135	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 carbohydrates in this entry.

## 5.6 Ligand geometry

6 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	6HY	D	301	-	10,14,14	2.50	2 (20%)	13,18,18	1.54	2 (15%)
2	6HY	B	301	-	10,14,14	2.50	2 (20%)	13,18,18	1.54	2 (15%)
2	6HY	A	301	-	10,14,14	2.49	2 (20%)	13,18,18	1.54	2 (15%)
2	6HY	C	301	-	10,14,14	2.49	2 (20%)	13,18,18	1.54	2 (15%)
2	6HY	E	301	-	10,14,14	2.49	2 (20%)	13,18,18	1.53	2 (15%)
2	6HY	E	302	-	10,14,14	2.52	2 (20%)	13,18,18	1.53	2 (15%)

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	6HY	D	301	-	-	1/7/9/9	0/1/1/1
2	6HY	B	301	-	-	1/7/9/9	0/1/1/1
2	6HY	A	301	-	-	1/7/9/9	0/1/1/1
2	6HY	C	301	-	-	1/7/9/9	0/1/1/1
2	6HY	E	301	-	-	1/7/9/9	0/1/1/1
2	6HY	E	302	-	-	1/7/9/9	0/1/1/1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	302	6HY	C2-C1	-6.91	1.38	1.49
2	B	301	6HY	C2-C1	-6.84	1.38	1.49
2	D	301	6HY	C2-C1	-6.83	1.38	1.49
2	A	301	6HY	C2-C1	-6.80	1.38	1.49
2	C	301	6HY	C2-C1	-6.80	1.38	1.49
2	E	301	6HY	C2-C1	-6.80	1.38	1.49
2	E	301	6HY	C4-N	-2.61	1.28	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	6HY	C4-N	-2.61	1.28	1.34
2	D	301	6HY	C4-N	-2.61	1.28	1.34
2	B	301	6HY	C4-N	-2.60	1.28	1.34
2	C	301	6HY	C4-N	-2.60	1.28	1.34
2	E	302	6HY	C4-N	-2.58	1.29	1.34

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	6HY	C4-N-C7	4.08	122.52	116.79
2	C	301	6HY	C4-N-C7	4.07	122.51	116.79
2	B	301	6HY	C4-N-C7	4.06	122.49	116.79
2	D	301	6HY	C4-N-C7	4.05	122.48	116.79
2	E	301	6HY	C4-N-C7	4.05	122.47	116.79
2	E	302	6HY	C4-N-C7	4.04	122.46	116.79
2	A	301	6HY	C6-C7-N	-2.84	119.91	124.79
2	D	301	6HY	C6-C7-N	-2.84	119.92	124.79
2	B	301	6HY	C6-C7-N	-2.84	119.92	124.79
2	E	302	6HY	C6-C7-N	-2.83	119.93	124.79
2	E	301	6HY	C6-C7-N	-2.83	119.94	124.79
2	C	301	6HY	C6-C7-N	-2.82	119.94	124.79

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	301	6HY	C1-C-C3-C8
2	B	301	6HY	C1-C-C3-C8
2	A	301	6HY	C1-C-C3-C8
2	C	301	6HY	C1-C-C3-C8
2	E	301	6HY	C1-C-C3-C8
2	E	302	6HY	C1-C-C3-C8

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	302	6HY	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, 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	200/208 (96%)	0.15	9 (4%) 33 31	31, 54, 99, 119	1 (0%)
1	B	206/208 (99%)	-0.19	3 (1%) 73 72	31, 49, 89, 110	1 (0%)
1	C	205/208 (98%)	-0.01	6 (2%) 51 50	31, 53, 94, 123	2 (0%)
1	D	203/208 (97%)	-0.00	6 (2%) 50 49	32, 52, 97, 122	1 (0%)
1	E	206/208 (99%)	-0.13	1 (0%) 91 89	33, 51, 92, 116	1 (0%)
1	F	201/208 (96%)	0.16	9 (4%) 33 31	34, 56, 97, 130	2 (0%)
All	All	1221/1248 (97%)	-0.01	34 (2%) 53 51	31, 52, 97, 130	8 (0%)

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	169	GLU	5.1
1	D	30	MET	4.8
1	A	169	GLU	4.7
1	C	30	MET	4.5
1	F	126	GLN	4.5
1	A	30	MET	4.3
1	D	170	GLY	4.2
1	A	127	ALA	4.1
1	F	172	ASP	4.1
1	B	30	MET	4.0
1	A	170	GLY	3.8
1	D	102	PRO	3.6
1	F	170	GLY	3.5
1	B	235	ASP	3.5
1	D	126	GLN	3.4
1	F	127	ALA	3.3
1	B	169	GLU	3.3
1	A	231	GLN	3.1
1	F	105	TYR	3.1

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Mol	Chain	Res	Type	RSRZ
1	C	169	GLU	3.0
1	F	59	ASN	2.8
1	C	234	GLN	2.7
1	C	128	TRP	2.7
1	F	169	GLU	2.6
1	A	233	THR	2.6
1	F	104	GLY	2.6
1	C	126	GLN	2.6
1	A	167	GLU	2.6
1	A	128	TRP	2.4
1	C	59	ASN	2.2
1	D	100	SER	2.2
1	E	171	GLU	2.2
1	F	231	GLN	2.1
1	A	172	ASP	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. 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	B-factors(Å <sup>2</sup> )	Q<0.9
2	6HY	C	301	14/14	0.92	0.12	20,20,20,20	0
2	6HY	A	301	14/14	0.94	0.13	20,20,20,20	0
2	6HY	E	302	14/14	0.95	0.13	20,20,20,20	0
2	6HY	E	301	14/14	0.96	0.14	20,20,20,20	0
2	6HY	D	301	14/14	0.96	0.09	20,20,20,20	0
2	6HY	B	301	14/14	0.97	0.12	20,20,20,20	0

## 6.5 Other polymers

There are no such residues in this entry.