



Full wwPDB X-ray Structure Validation Report ⓘ

May 21, 2014 – 11:57 PM EDT

PDB ID : 4NWP
Title : Computationally Designed Two-Component Self-Assembling Tetrahedral Cage, T33-21, Crystallized in Space Group R32
Authors : McNamara, D.E.; King, N.P.; Bale, J.B.; Sheffler, W.; Baker, D.; Yeates, T.O.
Deposited on : 2013-12-06
Resolution : 2.10 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

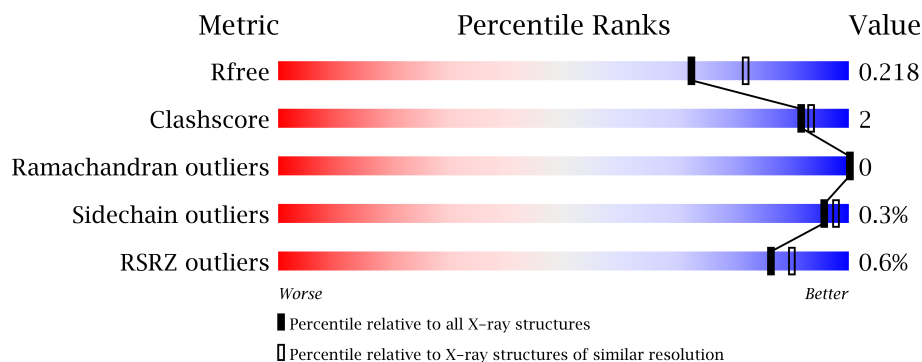
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.16 November 2013
Xtriage (Phenix) : dev-1439
EDS : stable22978
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22978

1 Overall quality at a glance

The reported resolution of this entry is 2.10 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}	66092	3012 (2.10-2.10)
Clashscore	79885	3649 (2.10-2.10)
Ramachandran outliers	78287	3610 (2.10-2.10)
Sidechain outliers	78261	3611 (2.10-2.10)
RSRZ outliers	66119	3013 (2.10-2.10)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	172	
1	B	172	
1	C	172	
1	D	172	
2	E	131	
2	F	131	
2	G	131	
2	H	131	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	SO4	D	203	-	X
3	SO4	E	203	-	X
3	SO4	F	204	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8248 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Putative uncharacterized protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	145	Total	C	N	O	S	0	1	0
			1087	700	183	199	5			
1	B	141	Total	C	N	O	S	0	0	0
			1054	677	176	196	5			
1	C	140	Total	C	N	O	S	0	0	0
			1060	681	175	199	5			
1	D	149	Total	C	N	O	S	0	0	0
			1127	726	189	207	5			

There are 52 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	85	ALA	LYS	ENGINEEREED MUTATION	UNP O58404
A	88	LEU	GLU	ENGINEEREED MUTATION	UNP O58404
A	89	LYS	GLY	ENGINEEREED MUTATION	UNP O58404
A	92	LEU	SER	ENGINEEREED MUTATION	UNP O58404
A	95	MET	GLU	ENGINEEREED MUTATION	UNP O58404
A	126	LEU	GLU	ENGINEEREED MUTATION	UNP O58404
A	130	LEU	ALA	ENGINEEREED MUTATION	UNP O58404
A	133	THR	LEU	ENGINEEREED MUTATION	UNP O58404
A	140	ALA	LYS	ENGINEEREED MUTATION	UNP O58404
A	143	ALA	LEU	ENGINEEREED MUTATION	UNP O58404
A	144	ALA	VAL	ENGINEEREED MUTATION	UNP O58404
A	147	LEU	ASN	ENGINEEREED MUTATION	UNP O58404
A	148	ALA	ARG	ENGINEEREED MUTATION	UNP O58404
B	85	ALA	LYS	ENGINEEREED MUTATION	UNP O58404
B	88	LEU	GLU	ENGINEEREED MUTATION	UNP O58404
B	89	LYS	GLY	ENGINEEREED MUTATION	UNP O58404
B	92	LEU	SER	ENGINEEREED MUTATION	UNP O58404
B	95	MET	GLU	ENGINEEREED MUTATION	UNP O58404
B	126	LEU	GLU	ENGINEEREED MUTATION	UNP O58404
B	130	LEU	ALA	ENGINEEREED MUTATION	UNP O58404
B	133	THR	LEU	ENGINEEREED MUTATION	UNP O58404

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	140	ALA	LYS	ENGINEEREED MUTATION	UNP O58404
B	143	ALA	LEU	ENGINEEREED MUTATION	UNP O58404
B	144	ALA	VAL	ENGINEEREED MUTATION	UNP O58404
B	147	LEU	ASN	ENGINEEREED MUTATION	UNP O58404
B	148	ALA	ARG	ENGINEEREED MUTATION	UNP O58404
C	85	ALA	LYS	ENGINEEREED MUTATION	UNP O58404
C	88	LEU	GLU	ENGINEEREED MUTATION	UNP O58404
C	89	LYS	GLY	ENGINEEREED MUTATION	UNP O58404
C	92	LEU	SER	ENGINEEREED MUTATION	UNP O58404
C	95	MET	GLU	ENGINEEREED MUTATION	UNP O58404
C	126	LEU	GLU	ENGINEEREED MUTATION	UNP O58404
C	130	LEU	ALA	ENGINEEREED MUTATION	UNP O58404
C	133	THR	LEU	ENGINEEREED MUTATION	UNP O58404
C	140	ALA	LYS	ENGINEEREED MUTATION	UNP O58404
C	143	ALA	LEU	ENGINEEREED MUTATION	UNP O58404
C	144	ALA	VAL	ENGINEEREED MUTATION	UNP O58404
C	147	LEU	ASN	ENGINEEREED MUTATION	UNP O58404
C	148	ALA	ARG	ENGINEEREED MUTATION	UNP O58404
D	85	ALA	LYS	ENGINEEREED MUTATION	UNP O58404
D	88	LEU	GLU	ENGINEEREED MUTATION	UNP O58404
D	89	LYS	GLY	ENGINEEREED MUTATION	UNP O58404
D	92	LEU	SER	ENGINEEREED MUTATION	UNP O58404
D	95	MET	GLU	ENGINEEREED MUTATION	UNP O58404
D	126	LEU	GLU	ENGINEEREED MUTATION	UNP O58404
D	130	LEU	ALA	ENGINEEREED MUTATION	UNP O58404
D	133	THR	LEU	ENGINEEREED MUTATION	UNP O58404
D	140	ALA	LYS	ENGINEEREED MUTATION	UNP O58404
D	143	ALA	LEU	ENGINEEREED MUTATION	UNP O58404
D	144	ALA	VAL	ENGINEEREED MUTATION	UNP O58404
D	147	LEU	ASN	ENGINEEREED MUTATION	UNP O58404
D	148	ALA	ARG	ENGINEEREED MUTATION	UNP O58404

- Molecule 2 is a protein called Uncharacterized protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	118	Total	C	N	O	S	0	5	0
			896	561	162	170	3			
2	F	118	Total	C	N	O	S	0	4	0
			889	556	163	167	3			
2	G	118	Total	C	N	O	S	0	4	0
			890	560	160	167	3			
2	H	117	Total	C	N	O	S	0	4	0
			879	552	159	165	3			

There are 64 discrepancies between the modelled and reference sequences:

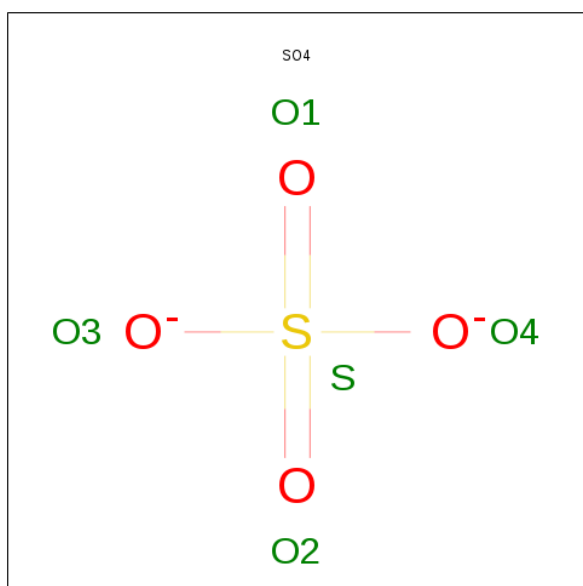
Chain	Residue	Modelled	Actual	Comment	Reference
E	27	LYS	ALA	ENGINEEREED MUTATION	UNP Q9I2D8
E	74	ILE	ALA	ENGINEEREED MUTATION	UNP Q9I2D8
E	78	THR	GLN	ENGINEEREED MUTATION	UNP Q9I2D8
E	79	LEU	ALA	ENGINEEREED MUTATION	UNP Q9I2D8
E	82	ALA	GLU	ENGINEEREED MUTATION	UNP Q9I2D8
E	86	ALA	GLU	ENGINEEREED MUTATION	UNP Q9I2D8
E	90	GLU	GLY	ENGINEEREED MUTATION	UNP Q9I2D8
E	112	LEU	ALA	ENGINEEREED MUTATION	UNP Q9I2D8
E	124	LEU	-	EXPRESSION TAG	UNP Q9I2D8
E	125	GLU	-	EXPRESSION TAG	UNP Q9I2D8
E	126	HIS	-	EXPRESSION TAG	UNP Q9I2D8
E	127	HIS	-	EXPRESSION TAG	UNP Q9I2D8
E	128	HIS	-	EXPRESSION TAG	UNP Q9I2D8
E	129	HIS	-	EXPRESSION TAG	UNP Q9I2D8
E	130	HIS	-	EXPRESSION TAG	UNP Q9I2D8
E	131	HIS	-	EXPRESSION TAG	UNP Q9I2D8
F	27	LYS	ALA	ENGINEEREED MUTATION	UNP Q9I2D8
F	74	ILE	ALA	ENGINEEREED MUTATION	UNP Q9I2D8
F	78	THR	GLN	ENGINEEREED MUTATION	UNP Q9I2D8
F	79	LEU	ALA	ENGINEEREED MUTATION	UNP Q9I2D8
F	82	ALA	GLU	ENGINEEREED MUTATION	UNP Q9I2D8
F	86	ALA	GLU	ENGINEEREED MUTATION	UNP Q9I2D8
F	90	GLU	GLY	ENGINEEREED MUTATION	UNP Q9I2D8
F	112	LEU	ALA	ENGINEEREED MUTATION	UNP Q9I2D8
F	124	LEU	-	EXPRESSION TAG	UNP Q9I2D8
F	125	GLU	-	EXPRESSION TAG	UNP Q9I2D8
F	126	HIS	-	EXPRESSION TAG	UNP Q9I2D8
F	127	HIS	-	EXPRESSION TAG	UNP Q9I2D8
F	128	HIS	-	EXPRESSION TAG	UNP Q9I2D8
F	129	HIS	-	EXPRESSION TAG	UNP Q9I2D8
F	130	HIS	-	EXPRESSION TAG	UNP Q9I2D8
F	131	HIS	-	EXPRESSION TAG	UNP Q9I2D8
G	27	LYS	ALA	ENGINEEREED MUTATION	UNP Q9I2D8
G	74	ILE	ALA	ENGINEEREED MUTATION	UNP Q9I2D8
G	78	THR	GLN	ENGINEEREED MUTATION	UNP Q9I2D8
G	79	LEU	ALA	ENGINEEREED MUTATION	UNP Q9I2D8
G	82	ALA	GLU	ENGINEEREED MUTATION	UNP Q9I2D8
G	86	ALA	GLU	ENGINEEREED MUTATION	UNP Q9I2D8
G	90	GLU	GLY	ENGINEEREED MUTATION	UNP Q9I2D8
G	112	LEU	ALA	ENGINEEREED MUTATION	UNP Q9I2D8
G	124	LEU	-	EXPRESSION TAG	UNP Q9I2D8
G	125	GLU	-	EXPRESSION TAG	UNP Q9I2D8

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
G	126	HIS	-	EXPRESSION TAG	UNP Q9I2D8
G	127	HIS	-	EXPRESSION TAG	UNP Q9I2D8
G	128	HIS	-	EXPRESSION TAG	UNP Q9I2D8
G	129	HIS	-	EXPRESSION TAG	UNP Q9I2D8
G	130	HIS	-	EXPRESSION TAG	UNP Q9I2D8
G	131	HIS	-	EXPRESSION TAG	UNP Q9I2D8
H	27	LYS	ALA	ENGINEEREED MUTATION	UNP Q9I2D8
H	74	ILE	ALA	ENGINEEREED MUTATION	UNP Q9I2D8
H	78	THR	GLN	ENGINEEREED MUTATION	UNP Q9I2D8
H	79	LEU	ALA	ENGINEEREED MUTATION	UNP Q9I2D8
H	82	ALA	GLU	ENGINEEREED MUTATION	UNP Q9I2D8
H	86	ALA	GLU	ENGINEEREED MUTATION	UNP Q9I2D8
H	90	GLU	GLY	ENGINEEREED MUTATION	UNP Q9I2D8
H	112	LEU	ALA	ENGINEEREED MUTATION	UNP Q9I2D8
H	124	LEU	-	EXPRESSION TAG	UNP Q9I2D8
H	125	GLU	-	EXPRESSION TAG	UNP Q9I2D8
H	126	HIS	-	EXPRESSION TAG	UNP Q9I2D8
H	127	HIS	-	EXPRESSION TAG	UNP Q9I2D8
H	128	HIS	-	EXPRESSION TAG	UNP Q9I2D8
H	129	HIS	-	EXPRESSION TAG	UNP Q9I2D8
H	130	HIS	-	EXPRESSION TAG	UNP Q9I2D8
H	131	HIS	-	EXPRESSION TAG	UNP Q9I2D8

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



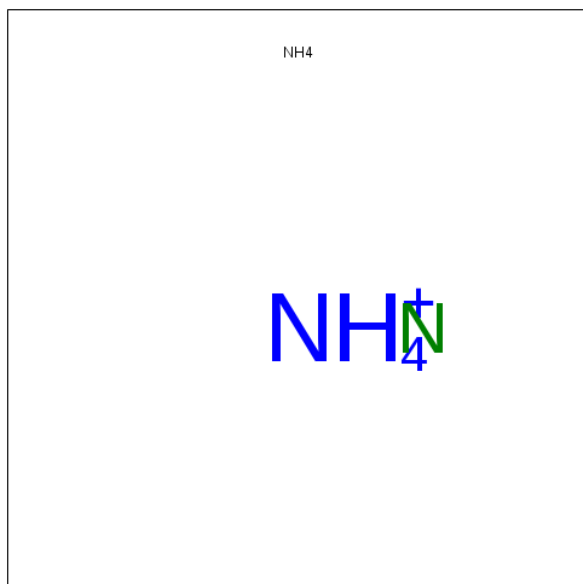
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0
3	D	1	Total O S 5 4 1	0	0
3	D	1	Total O S 5 4 1	0	0
3	D	1	Total O S 5 4 1	0	0
3	E	1	Total O S 5 4 1	0	0
3	E	1	Total O S 5 4 1	0	0
3	E	1	Total O S 5 4 1	0	0
3	E	1	Total O S 5 4 1	0	0
3	F	1	Total O S 5 4 1	0	0
3	F	1	Total O S 5 4 1	0	0
3	F	1	Total O S 5 4 1	0	0
3	F	1	Total O S 5 4 1	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	G	1	Total O S 5 4 1	0	0
3	G	1	Total O S 5 4 1	0	0
3	G	1	Total O S 5 4 1	0	0
3	H	1	Total O S 5 4 1	0	0
3	H	1	Total O S 5 4 1	0	0
3	H	1	Total O S 5 4 1	0	0

- Molecule 4 is AMMONIUM ION (three-letter code: NH4) (formula: H₄N).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	H	1	Total N 1 1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	25	Total O 25 25	0	0
5	B	29	Total O 29 29	0	0

Continued on next page...

Continued from previous page...

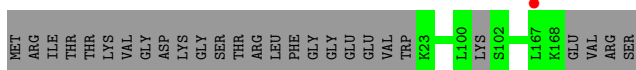
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	22	Total 22	O 22	0	0
5	D	26	Total 26	O 26	0	0
5	E	33	Total 33	O 33	0	0
5	F	29	Total 29	O 29	0	0
5	G	37	Total 37	O 37	0	0
5	H	24	Total 24	O 24	0	0

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.

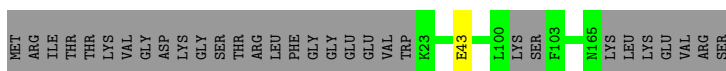
- Molecule 1: Putative uncharacterized protein

Chain A: 



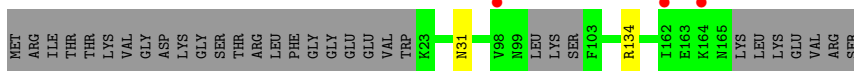
- Molecule 1: Putative uncharacterized protein

Chain B: 



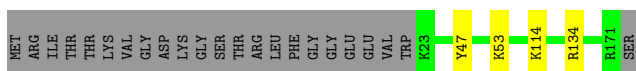
- Molecule 1: Putative uncharacterized protein

Chain C: 



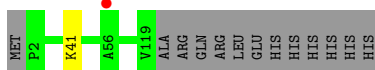
- Molecule 1: Putative uncharacterized protein

Chain D: 



- Molecule 2: Uncharacterized protein

Chain E: 



- Molecule 2: Uncharacterized protein

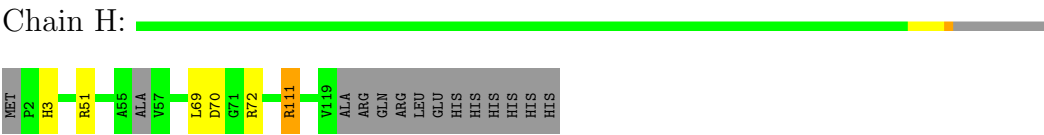
Chain F: 



● Molecule 2: Uncharacterized protein



● Molecule 2: Uncharacterized protein



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	113.35Å 113.35Å 634.88Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	93.78 – 2.10 93.78 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.9 (93.78-2.10) 99.9 (93.78-2.10)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.21 (at 2.10Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE: 1.8.4_1496)	Depositor
R, R_{free}	0.188 , 0.218 0.188 , 0.218	Depositor DCC
R_{free} test set	9242 reflections (10.00%)	DCC
Wilson B-factor (Å ²)	37.7	Xtriage
Anisotropy	0.177	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 40.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 92425 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8248	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.57% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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: NH4, SO4

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	0.25	0/1100	0.37	0/1485
1	B	0.26	0/1064	0.37	0/1436
1	C	0.26	0/1070	0.38	0/1443
1	D	0.25	0/1138	0.38	0/1534
2	E	0.25	0/920	0.40	0/1243
2	F	0.24	0/907	0.39	0/1226
2	G	0.25	0/911	0.40	0/1231
2	H	0.25	0/899	0.40	0/1214
All	All	0.25	0/8009	0.38	0/10812

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1087	0	0	0	0
1	B	1054	0	0	1	0
1	C	1060	0	0	1	0
1	D	1127	0	0	4	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	896	0	0	1	0
2	F	889	0	0	3	0
2	G	890	0	0	2	0
2	H	879	0	0	5	0
3	A	25	0	0	0	0
3	B	15	0	0	0	0
3	C	15	0	0	1	0
3	D	15	0	0	3	0
3	E	20	0	0	1	0
3	F	20	0	0	2	0
3	G	15	0	0	0	0
3	H	15	0	0	3	0
4	H	1	0	0	0	0
5	A	25	0	0	0	0
5	B	29	0	0	0	0
5	C	22	0	0	0	0
5	D	26	0	0	0	0
5	E	33	0	0	0	0
5	F	29	0	0	0	0
5	G	37	0	0	0	0
5	H	24	0	0	0	0
All	All	8248	0	0	15	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 2.

All (15) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:111:ARG:NH2	3:H:202:SO4:O3	2.22	0.72
2:G:34:GLN:OE1	2:G:72[B]:ARG:NH2	2.30	0.64
2:H:3:HIS:NE2	3:H:203:SO4:O3	2.32	0.63
2:H:70:ASP:O	2:H:111:ARG:NH1	2.36	0.59
2:H:51:ARG:NH1	3:H:201:SO4:O1	2.39	0.56
1:C:134:ARG:NH2	3:C:202:SO4:O3	2.39	0.56
2:F:72[A]:ARG:NH2	3:F:202:SO4:S	2.80	0.54
2:H:69:LEU:O	2:H:72[A]:ARG:NH2	2.42	0.53
1:D:53:LYS:NZ	3:D:202:SO4:O1	2.43	0.52
1:D:134:ARG:NH2	3:D:201:SO4:O1	2.48	0.46
2:F:37:GLU:OE2	2:G:59:ARG:NH1	2.50	0.44
2:E:41:LYS:NZ	3:E:204:SO4:O4	2.51	0.43
1:D:47:TYR:OH	3:D:203:SO4:O4	2.38	0.41
1:B:43:GLU:OE2	1:D:114:LYS:NZ	2.54	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:F:72[A]:ARG:NH2	3:F:202:SO4:O4	2.53	0.41

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	142/172 (83%)	141 (99%)	1 (1%)	0	100	100
1	B	137/172 (80%)	137 (100%)	0	0	100	100
1	C	136/172 (79%)	136 (100%)	0	0	100	100
1	D	147/172 (86%)	146 (99%)	1 (1%)	0	100	100
2	E	121/131 (92%)	121 (100%)	0	0	100	100
2	F	120/131 (92%)	119 (99%)	1 (1%)	0	100	100
2	G	120/131 (92%)	119 (99%)	1 (1%)	0	100	100
2	H	117/131 (89%)	117 (100%)	0	0	100	100
All	All	1040/1212 (86%)	1036 (100%)	4 (0%)	0	100	100

There are no Ramachandran outliers to report.

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	106/145 (73%)	106 (100%)	0	100	100
1	B	103/145 (71%)	103 (100%)	0	100	100
1	C	106/145 (73%)	105 (99%)	1 (1%)	87	91

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	113/145 (78%)	113 (100%)	0	100	100
2	E	90/102 (88%)	90 (100%)	0	100	100
2	F	88/102 (86%)	88 (100%)	0	100	100
2	G	89/102 (87%)	89 (100%)	0	100	100
2	H	89/102 (87%)	88 (99%)	1 (1%)	84	88
All	All	784/988 (79%)	782 (100%)	2 (0%)	96	98

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	31	ASN
2	H	111	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA chains 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 29 ligands modelled in this entry, 1 is modelled with single atom - leaving 28 for Mogul analysis.

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	SO4	A	201	-	4,4,4	0.27	0	6,6,6	0.07	0
3	SO4	A	202	-	4,4,4	0.20	0	6,6,6	0.14	0
3	SO4	A	203	-	4,4,4	0.19	0	6,6,6	0.10	0
3	SO4	A	204	-	4,4,4	0.23	0	6,6,6	0.08	0
3	SO4	A	205	-	4,4,4	0.25	0	6,6,6	0.07	0
3	SO4	B	201	-	4,4,4	0.26	0	6,6,6	0.11	0
3	SO4	B	202	-	4,4,4	0.21	0	6,6,6	0.15	0
3	SO4	B	203	-	4,4,4	0.23	0	6,6,6	0.07	0
3	SO4	C	201	-	4,4,4	0.20	0	6,6,6	0.07	0
3	SO4	C	202	-	4,4,4	0.22	0	6,6,6	0.09	0
3	SO4	C	203	-	4,4,4	0.21	0	6,6,6	0.09	0
3	SO4	D	201	-	4,4,4	0.25	0	6,6,6	0.07	0
3	SO4	D	202	-	4,4,4	0.20	0	6,6,6	0.08	0
3	SO4	D	203	-	4,4,4	0.21	0	6,6,6	0.08	0
3	SO4	E	201	-	4,4,4	0.26	0	6,6,6	0.15	0
3	SO4	E	202	-	4,4,4	0.21	0	6,6,6	0.08	0
3	SO4	E	203	-	4,4,4	0.24	0	6,6,6	0.09	0
3	SO4	E	204	-	4,4,4	0.20	0	6,6,6	0.10	0
3	SO4	F	201	-	4,4,4	0.23	0	6,6,6	0.14	0
3	SO4	F	202	-	4,4,4	0.22	0	6,6,6	0.13	0
3	SO4	F	203	-	4,4,4	0.26	0	6,6,6	0.17	0
3	SO4	F	204	-	4,4,4	0.22	0	6,6,6	0.06	0
3	SO4	G	201	-	4,4,4	0.27	0	6,6,6	0.09	0
3	SO4	G	202	-	4,4,4	0.22	0	6,6,6	0.09	0
3	SO4	G	203	-	4,4,4	0.25	0	6,6,6	0.14	0
3	SO4	H	201	-	4,4,4	0.20	0	6,6,6	0.09	0
3	SO4	H	202	-	4,4,4	0.21	0	6,6,6	0.13	0
3	SO4	H	203	-	4,4,4	0.20	0	6,6,6	0.10	0

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	SO4	A	201	-	-	0/0/0/0	0/0/0/0
3	SO4	A	202	-	-	0/0/0/0	0/0/0/0
3	SO4	A	203	-	-	0/0/0/0	0/0/0/0
3	SO4	A	204	-	-	0/0/0/0	0/0/0/0
3	SO4	A	205	-	-	0/0/0/0	0/0/0/0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SO4	B	201	-	-	0/0/0/0	0/0/0/0
3	SO4	B	202	-	-	0/0/0/0	0/0/0/0
3	SO4	B	203	-	-	0/0/0/0	0/0/0/0
3	SO4	C	201	-	-	0/0/0/0	0/0/0/0
3	SO4	C	202	-	-	0/0/0/0	0/0/0/0
3	SO4	C	203	-	-	0/0/0/0	0/0/0/0
3	SO4	D	201	-	-	0/0/0/0	0/0/0/0
3	SO4	D	202	-	-	0/0/0/0	0/0/0/0
3	SO4	D	203	-	-	0/0/0/0	0/0/0/0
3	SO4	E	201	-	-	0/0/0/0	0/0/0/0
3	SO4	E	202	-	-	0/0/0/0	0/0/0/0
3	SO4	E	203	-	-	0/0/0/0	0/0/0/0
3	SO4	E	204	-	-	0/0/0/0	0/0/0/0
3	SO4	F	201	-	-	0/0/0/0	0/0/0/0
3	SO4	F	202	-	-	0/0/0/0	0/0/0/0
3	SO4	F	203	-	-	0/0/0/0	0/0/0/0
3	SO4	F	204	-	-	0/0/0/0	0/0/0/0
3	SO4	G	201	-	-	0/0/0/0	0/0/0/0
3	SO4	G	202	-	-	0/0/0/0	0/0/0/0
3	SO4	G	203	-	-	0/0/0/0	0/0/0/0
3	SO4	H	201	-	-	0/0/0/0	0/0/0/0
3	SO4	H	202	-	-	0/0/0/0	0/0/0/0
3	SO4	H	203	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 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	145/172 (84%)	-0.27	1 (0%) 84 88	26, 39, 65, 80	0
1	B	141/172 (81%)	-0.29	0 100 100	24, 38, 68, 90	0
1	C	140/172 (81%)	-0.18	3 (2%) 60 65	25, 42, 72, 87	0
1	D	149/172 (86%)	-0.21	0 100 100	25, 39, 69, 87	0
2	E	118/131 (90%)	-0.33	1 (0%) 83 87	26, 37, 55, 67	0
2	F	118/131 (90%)	-0.30	1 (0%) 83 87	25, 43, 69, 82	0
2	G	118/131 (90%)	-0.34	0 100 100	24, 39, 59, 75	0
2	H	117/131 (89%)	-0.30	0 100 100	26, 42, 67, 82	0
All	All	1046/1212 (86%)	-0.27	6 (0%) 86 90	24, 40, 68, 90	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	162	ILE	4.2
2	F	56	ALA	4.0
1	A	167	LEU	3.0
2	E	56	ALA	2.5
1	C	164	LYS	2.2
1	C	98	VAL	2.1

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 ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	SO4	E	203	5/5	0.16	4.53	28,29,29,33	5
3	SO4	F	204	5/5	0.17	4.31	98,105,105,106	0
3	SO4	D	203	5/5	0.19	3.42	97,98,99,100	0
3	SO4	H	203	5/5	0.15	1.98	103,105,110,116	0
3	SO4	C	202	5/5	0.16	1.85	45,47,54,54	5
3	SO4	F	201	5/5	0.13	1.84	29,29,32,33	0
3	SO4	A	204	5/5	0.12	1.12	41,41,42,42	5
3	SO4	C	203	5/5	0.19	1.03	95,99,100,100	5
3	SO4	D	202	5/5	0.14	0.97	76,78,82,86	5
3	SO4	C	201	5/5	0.14	0.88	77,82,83,84	5
3	SO4	G	202	5/5	0.13	0.85	101,103,106,110	0
3	SO4	D	201	5/5	0.15	0.76	42,44,50,58	5
3	SO4	G	203	5/5	0.16	0.75	49,53,60,64	5
3	SO4	A	205	5/5	0.11	0.70	44,45,45,46	5
4	NH4	H	204	1/1	0.10	0.68	22,22,22,22	0
3	SO4	E	204	5/5	0.11	0.58	104,104,107,109	0
3	SO4	B	203	5/5	0.12	0.46	79,81,85,87	0
3	SO4	F	202	5/5	0.14	0.45	51,53,59,62	5
3	SO4	F	203	5/5	0.15	0.40	39,40,45,49	5
3	SO4	E	201	5/5	0.13	0.34	35,42,48,50	0
3	SO4	B	202	5/5	0.13	0.33	40,50,55,59	5
3	SO4	A	203	5/5	0.12	0.25	63,64,67,69	5
3	SO4	B	201	5/5	0.12	0.16	38,39,44,51	0
3	SO4	H	201	5/5	0.12	0.11	70,73,76,78	5
3	SO4	A	201	5/5	0.12	-0.20	41,44,48,51	0
3	SO4	G	201	5/5	0.16	-0.20	34,40,46,47	0
3	SO4	A	202	5/5	0.11	-0.74	49,49,55,58	5
3	SO4	E	202	5/5	0.08	-1.19	55,56,67,67	5
3	SO4	H	202	5/5	0.07	-1.86	43,56,63,64	0

6.5 Other polymers ⓘ

There are no such residues in this entry.