



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 09:57 am GMT

PDB ID : 2ICR
Title : Red fluorescent protein zRFP574 from Zoanthus sp.
Authors : Pletnev, S.; Pletneva, N.; Tikhonova, T.; Pletnev, V.
Deposited on : 2006-09-13
Resolution : 1.51 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28949

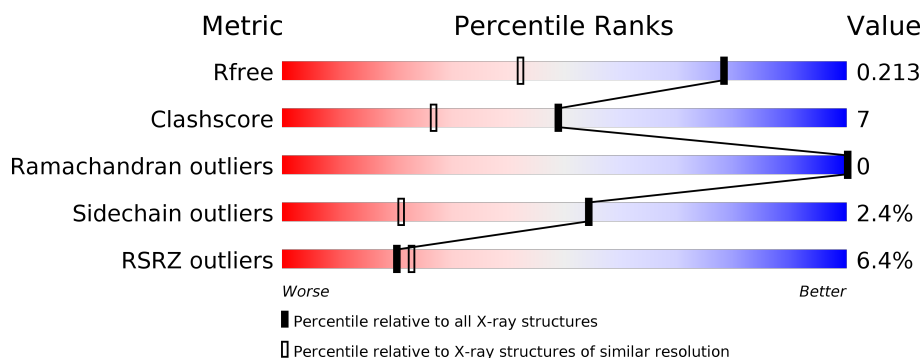
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.51 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}	100719	2964 (1.54-1.50)
Clashscore	112137	3216 (1.54-1.50)
Ramachandran outliers	110173	3145 (1.54-1.50)
Sidechain outliers	110143	3143 (1.54-1.50)
RSRZ outliers	101464	2990 (1.54-1.50)

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	237	<div> <div>3%</div> <div>86%</div> <div>8%</div> <div>5%</div> </div>
1	B	237	<div> <div>2%</div> <div>83%</div> <div>12%</div> <div>5%</div> </div>
1	C	237	<div> <div>3%</div> <div>84%</div> <div>11%</div> <div>•</div> </div>
1	D	237	<div> <div>17%</div> <div>74%</div> <div>19%</div> <div>•</div> <div>5%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7857 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 Red fluorescent protein zoanRFP.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	226	Total	C	N	O	S	0	4	0
			1843	1173	314	339	17			
1	B	226	Total	C	N	O	S	0	6	0
			1853	1181	317	339	16			
1	C	227	Total	C	N	O	S	0	7	0
			1869	1194	318	340	17			
1	D	226	Total	C	N	O	S	0	4	0
			1845	1173	315	339	18			

There are 60 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	MET	-	EXPRESSION TAG	UNP Q8T4U4
A	-6	ARG	-	EXPRESSION TAG	UNP Q8T4U4
A	-5	GLY	-	EXPRESSION TAG	UNP Q8T4U4
A	-4	SER	-	EXPRESSION TAG	UNP Q8T4U4
A	-3	HIS	-	EXPRESSION TAG	UNP Q8T4U4
A	-2	HIS	-	EXPRESSION TAG	UNP Q8T4U4
A	-1	HIS	-	EXPRESSION TAG	UNP Q8T4U4
A	0	HIS	-	EXPRESSION TAG	UNP Q8T4U4
A	1	HIS	-	EXPRESSION TAG	UNP Q8T4U4
A	2	HIS	-	EXPRESSION TAG	UNP Q8T4U4
A	3	GLY	-	EXPRESSION TAG	UNP Q8T4U4
A	5	ALA	LYS	ENGINEERED	UNP Q8T4U4
A	66	XYG	ASP	CHROMOPHORE	UNP Q8T4U4
A	66	XYG	TYR	CHROMOPHORE	UNP Q8T4U4
A	66	XYG	GLY	CHROMOPHORE	UNP Q8T4U4
B	-7	MET	-	EXPRESSION TAG	UNP Q8T4U4
B	-6	ARG	-	EXPRESSION TAG	UNP Q8T4U4
B	-5	GLY	-	EXPRESSION TAG	UNP Q8T4U4
B	-4	SER	-	EXPRESSION TAG	UNP Q8T4U4
B	-3	HIS	-	EXPRESSION TAG	UNP Q8T4U4
B	-2	HIS	-	EXPRESSION TAG	UNP Q8T4U4

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-1	HIS	-	EXPRESSION TAG	UNP Q8T4U4
B	0	HIS	-	EXPRESSION TAG	UNP Q8T4U4
B	1	HIS	-	EXPRESSION TAG	UNP Q8T4U4
B	2	HIS	-	EXPRESSION TAG	UNP Q8T4U4
B	3	GLY	-	EXPRESSION TAG	UNP Q8T4U4
B	5	ALA	LYS	ENGINEERED	UNP Q8T4U4
B	66	XYG	ASP	CHROMOPHORE	UNP Q8T4U4
B	66	XYG	TYR	CHROMOPHORE	UNP Q8T4U4
B	66	XYG	GLY	CHROMOPHORE	UNP Q8T4U4
C	-7	MET	-	EXPRESSION TAG	UNP Q8T4U4
C	-6	ARG	-	EXPRESSION TAG	UNP Q8T4U4
C	-5	GLY	-	EXPRESSION TAG	UNP Q8T4U4
C	-4	SER	-	EXPRESSION TAG	UNP Q8T4U4
C	-3	HIS	-	EXPRESSION TAG	UNP Q8T4U4
C	-2	HIS	-	EXPRESSION TAG	UNP Q8T4U4
C	-1	HIS	-	EXPRESSION TAG	UNP Q8T4U4
C	0	HIS	-	EXPRESSION TAG	UNP Q8T4U4
C	1	HIS	-	EXPRESSION TAG	UNP Q8T4U4
C	2	HIS	-	EXPRESSION TAG	UNP Q8T4U4
C	3	GLY	-	EXPRESSION TAG	UNP Q8T4U4
C	5	ALA	LYS	ENGINEERED	UNP Q8T4U4
C	66	XYG	ASP	CHROMOPHORE	UNP Q8T4U4
C	66	XYG	TYR	CHROMOPHORE	UNP Q8T4U4
C	66	XYG	GLY	CHROMOPHORE	UNP Q8T4U4
D	-7	MET	-	EXPRESSION TAG	UNP Q8T4U4
D	-6	ARG	-	EXPRESSION TAG	UNP Q8T4U4
D	-5	GLY	-	EXPRESSION TAG	UNP Q8T4U4
D	-4	SER	-	EXPRESSION TAG	UNP Q8T4U4
D	-3	HIS	-	EXPRESSION TAG	UNP Q8T4U4
D	-2	HIS	-	EXPRESSION TAG	UNP Q8T4U4
D	-1	HIS	-	EXPRESSION TAG	UNP Q8T4U4
D	0	HIS	-	EXPRESSION TAG	UNP Q8T4U4
D	1	HIS	-	EXPRESSION TAG	UNP Q8T4U4
D	2	HIS	-	EXPRESSION TAG	UNP Q8T4U4
D	3	GLY	-	EXPRESSION TAG	UNP Q8T4U4
D	5	ALA	LYS	ENGINEERED	UNP Q8T4U4
D	66	XYG	ASP	CHROMOPHORE	UNP Q8T4U4
D	66	XYG	TYR	CHROMOPHORE	UNP Q8T4U4
D	66	XYG	GLY	CHROMOPHORE	UNP Q8T4U4

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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


- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	133	Total	O	0	0
			133	133		
3	B	113	Total	O	0	0
			113	113		
3	C	115	Total	O	0	0
			115	115		
3	D	66	Total	O	0	0
			66	66		

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.

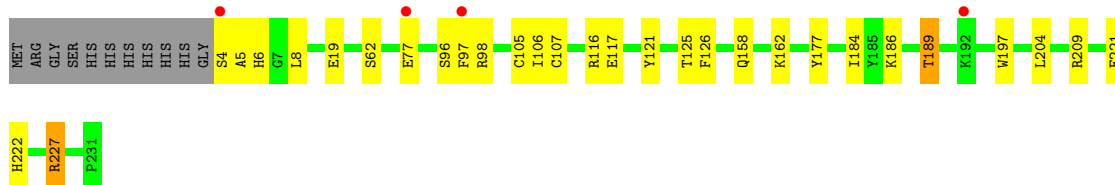
- Molecule 1: Red fluorescent protein zoanRFP

Chain A: 




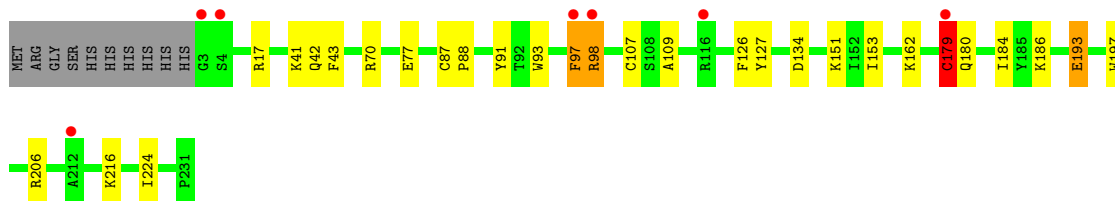
- Molecule 1: Red fluorescent protein zoanRFP

Chain B: 




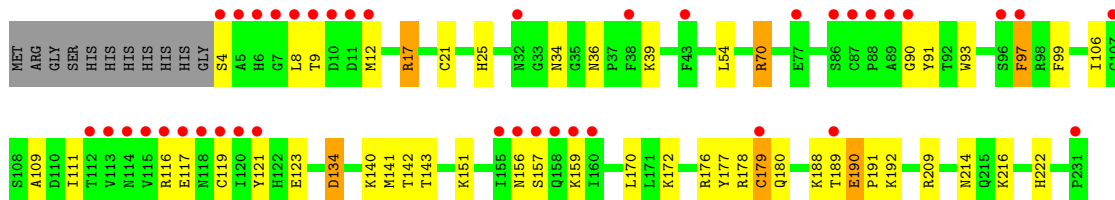
- Molecule 1: Red fluorescent protein zoanRFP

Chain C: 



- Molecule 1: Red fluorescent protein zoanRFP

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	115.19Å 146.74Å 122.93Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 1.51 45.02 – 1.46	Depositor EDS
% Data completeness (in resolution range)	99.4 (30.00-1.51) 96.4 (45.02-1.46)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.39 (at 1.46Å)	Xtriage
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.178 , 0.225 0.177 , 0.213	Depositor DCC
R_{free} test set	1613 reflections (1.01%)	DCC
Wilson B-factor (Å ²)	19.1	Xtriage
Anisotropy	0.134	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 50.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	7857	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.37% 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, XYG

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	0.60	0/1880	1.32	9/2535 (0.4%)
1	B	0.55	0/1899	1.33	14/2561 (0.5%)
1	C	0.57	0/1919	1.29	8/2590 (0.3%)
1	D	0.52	0/1882	1.37	16/2537 (0.6%)
All	All	0.56	0/7580	1.33	47/10223 (0.5%)

There are no bond length outliers.

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	126	PHE	CB-CG-CD2	10.04	127.83	120.80
1	D	178	ARG	NE-CZ-NH2	9.47	125.04	120.30
1	B	116	ARG	NE-CZ-NH2	-8.72	115.94	120.30
1	D	209	ARG	NE-CZ-NH1	-8.66	115.97	120.30
1	A	206	ARG	NE-CZ-NH1	8.64	124.62	120.30
1	D	178	ARG	CD-NE-CZ	8.63	135.68	123.60
1	B	126	PHE	CB-CG-CD1	-8.30	114.99	120.80
1	D	116	ARG	NE-CZ-NH2	-8.28	116.16	120.30
1	B	209	ARG	NE-CZ-NH1	-7.89	116.35	120.30
1	D	177	TYR	CB-CG-CD1	6.99	125.19	121.00
1	D	117	GLU	OE1-CD-OE2	-6.90	115.02	123.30
1	C	126	PHE	CB-CG-CD2	6.75	125.53	120.80
1	B	209	ARG	NE-CZ-NH2	6.66	123.63	120.30
1	C	126	PHE	CB-CG-CD1	-6.48	116.27	120.80
1	B	117	GLU	OE1-CD-OE2	-6.40	115.62	123.30
1	D	17	ARG	NE-CZ-NH1	6.18	123.39	120.30
1	D	116	ARG	CD-NE-CZ	6.16	132.23	123.60
1	B	116	ARG	NE-CZ-NH1	6.04	123.32	120.30
1	D	177	TYR	CB-CG-CD2	-6.01	117.39	121.00
1	C	206	ARG	NE-CZ-NH2	-6.01	117.30	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	196	ASP	CB-CG-OD2	5.99	123.69	118.30
1	A	197	TRP	CG-CD2-CE3	-5.91	128.58	133.90
1	B	117	GLU	CB-CG-CD	5.89	130.11	114.20
1	B	116	ARG	CD-NE-CZ	5.85	131.79	123.60
1	D	117	GLU	CB-CG-CD	5.84	129.96	114.20
1	B	197	TRP	CG-CD2-CE3	-5.79	128.69	133.90
1	A	178	ARG	NE-CZ-NH2	-5.69	117.46	120.30
1	A	178	ARG	NE-CZ-NH1	5.64	123.12	120.30
1	B	227	ARG	NE-CZ-NH2	5.58	123.09	120.30
1	A	17	ARG	NE-CZ-NH1	-5.54	117.53	120.30
1	B	96	SER	N-CA-CB	5.43	118.64	110.50
1	C	91	TYR	CB-CG-CD2	5.42	124.25	121.00
1	B	177	TYR	CB-CG-CD2	-5.40	117.76	121.00
1	C	97	PHE	O-C-N	5.34	131.25	122.70
1	D	178	ARG	NE-CZ-NH1	-5.31	117.64	120.30
1	C	197	TRP	CG-CD2-CE3	-5.26	129.17	133.90
1	C	193	GLU	OE1-CD-OE2	-5.23	117.02	123.30
1	C	179	CYS	CB-CA-C	-5.13	100.13	110.40
1	B	121	TYR	CB-CG-CD2	5.09	124.05	121.00
1	D	134	ASP	CB-CG-OD2	5.09	122.88	118.30
1	A	177	TYR	CB-CG-CD1	5.08	124.05	121.00
1	D	214	ASN	CA-CB-CG	5.08	124.57	113.40
1	D	97	PHE	N-CA-CB	5.07	119.73	110.60
1	D	70	ARG	NE-CZ-NH2	-5.05	117.77	120.30
1	A	196	ASP	CB-CG-OD1	-5.04	113.77	118.30
1	D	179	CYS	CB-CA-C	-5.03	100.34	110.40
1	A	17	ARG	NE-CZ-NH2	5.01	122.80	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	1843	0	1774	14	0
1	B	1853	0	1795	16	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	1869	0	1814	27	0
1	D	1845	0	1772	49	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	0	0
3	A	133	0	0	1	0
3	B	113	0	0	3	0
3	C	115	0	0	4	0
3	D	66	0	0	1	0
All	All	7857	0	7155	99	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 (99) 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:97:PHE:HE1	1:A:107[A]:CYS:SG	1.36	1.48
1:A:97:PHE:CE1	1:A:107[A]:CYS:SG	2.07	1.47
1:D:97:PHE:CD2	1:D:179:CYS:SG	2.15	1.39
1:C:97:PHE:CD2	1:C:179:CYS:SG	2.42	1.12
1:D:8:LEU:CD2	1:D:12[B]:MET:HE1	1.80	1.12
1:D:8:LEU:HD23	1:D:12[B]:MET:HE1	1.16	1.10
1:A:97:PHE:CZ	1:A:107[A]:CYS:SG	2.45	1.09
1:C:98:ARG:HB3	3:C:354:HOH:O	1.52	1.09
1:C:97:PHE:HD2	1:C:179:CYS:SG	1.77	1.06
1:D:8:LEU:CD2	1:D:12[B]:MET:CE	2.40	0.98
1:D:97:PHE:HD2	1:D:179:CYS:SG	1.68	0.98
1:D:8:LEU:HD23	1:D:12[B]:MET:CE	1.98	0.92
1:D:8:LEU:HD22	1:D:12[B]:MET:CE	2.06	0.85
1:D:8:LEU:HD22	1:D:12[B]:MET:HE3	1.60	0.84
1:D:97:PHE:CG	1:D:179:CYS:SG	2.75	0.80
1:B:77:GLU:HG3	3:B:322:HOH:O	1.86	0.75
1:D:17:ARG:HB3	1:D:123:GLU:OE1	1.90	0.71
1:D:97:PHE:HE2	1:D:99:PHE:CZ	2.10	0.69
1:D:97:PHE:CB	1:D:180:GLN:O	2.40	0.69
1:D:119:CYS:SG	1:D:121:TYR:HE1	2.17	0.67
1:D:97:PHE:HB3	1:D:180:GLN:O	1.95	0.66
1:B:97:PHE:HB2	1:B:105:CYS:HB2	1.77	0.65
1:C:162:LYS:HG2	1:C:184[B]:ILE:HD13	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:97:PHE:HB2	1:D:179:CYS:SG	2.36	0.64
1:A:127:TYR:HB2	1:D:106:ILE:HD11	1.80	0.64
1:D:119:CYS:SG	1:D:121:TYR:CE1	2.90	0.64
1:C:17:ARG:NH2	1:D:156:ASN:HB3	2.15	0.62
1:D:97:PHE:HA	1:D:180:GLN:O	2.01	0.61
1:A:97:PHE:HB2	1:A:105:CYS:HB2	1.83	0.60
1:A:97:PHE:HZ	1:A:107[A]:CYS:SG	2.21	0.60
1:D:8:LEU:HA	1:D:12[B]:MET:HE1	1.84	0.59
1:C:151:LYS:HE3	1:C:153[A]:ILE:HD11	1.86	0.58
1:B:106:ILE:HD11	1:C:127:TYR:HB2	1.85	0.58
1:D:170:LEU:HD23	1:D:176:ARG:HG2	1.86	0.57
1:D:97:PHE:HA	3:D:281:HOH:O	2.05	0.56
1:B:162:LYS:HD2	1:B:184:ILE:CD1	2.35	0.56
1:D:97:PHE:HE2	1:D:99:PHE:CE1	2.24	0.55
1:D:143:THR:CG2	1:D:172:LYS:HG3	2.36	0.55
1:C:41:LYS:HE2	1:C:43:PHE:CZ	2.42	0.54
1:B:106:ILE:HD11	1:C:127:TYR:CB	2.38	0.54
1:D:97:PHE:CE2	1:D:99:PHE:CE1	2.96	0.53
1:C:97:PHE:HB2	1:C:179:CYS:SG	2.49	0.53
1:C:162:LYS:HG2	1:C:184[B]:ILE:CD1	2.37	0.53
1:D:97:PHE:CE2	1:D:99:PHE:CZ	2.94	0.53
1:D:142:THR:O	1:D:172:LYS:HE3	2.07	0.52
1:A:5:ALA:O	1:A:6:HIS:HB2	2.10	0.52
1:B:189:THR:HG21	3:B:303:HOH:O	2.10	0.52
1:A:127:TYR:CB	1:D:106:ILE:HD11	2.39	0.51
1:C:151:LYS:CE	1:C:153[A]:ILE:HD11	2.40	0.51
1:D:97:PHE:CB	1:D:179:CYS:SG	2.99	0.50
1:D:54:LEU:O	1:D:140:LYS:HE2	2.12	0.50
1:C:17:ARG:HH22	1:D:156:ASN:HB3	1.75	0.50
1:D:159:LYS:HD2	1:D:190:GLU:OE2	2.11	0.50
1:C:97:PHE:HA	3:C:293:HOH:O	2.12	0.49
1:D:93:TRP:NE1	1:D:109:ALA:HB3	2.26	0.49
1:B:162:LYS:HD2	1:B:184:ILE:HD13	1.95	0.49
1:D:12[A]:MET:O	1:D:34:ASN:HA	2.14	0.47
1:A:59:ASP:HB3	1:A:169:LEU:HD21	1.97	0.46
1:C:98:ARG:CB	3:C:354:HOH:O	2.33	0.46
1:A:222:HIS:HE1	3:A:333:HOH:O	1.99	0.45
1:C:87:CYS:HB3	1:C:88:PRO:HA	1.98	0.45
1:C:184[B]:ILE:CG2	1:C:186:LYS:HE2	2.46	0.45
1:D:141[B]:MET:HA	1:D:172:LYS:HD2	1.99	0.44
1:D:9:THR:H	1:D:12[B]:MET:HE2	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:36:ASN:ND2	1:D:39:LYS:HG3	2.32	0.44
1:A:4:SER:HB3	1:A:7:GLY:HA2	1.99	0.44
1:B:162:LYS:HD2	1:B:184:ILE:HD11	1.99	0.44
1:C:97:PHE:HA	1:C:180:GLN:O	2.17	0.44
1:C:97:PHE:CB	1:C:180:GLN:O	2.65	0.43
1:B:5:ALA:O	1:B:6:HIS:HB2	2.18	0.43
1:B:98:ARG:HD2	1:B:98:ARG:HH21	1.65	0.43
1:D:97:PHE:CA	1:D:180:GLN:O	2.67	0.43
1:D:190:GLU:HA	1:D:191:PRO:HD3	1.72	0.43
1:D:90:GLY:HA3	1:D:111:ILE:O	2.19	0.42
1:A:143:THR:HG21	3:B:348:HOH:O	2.18	0.42
1:C:77:GLU:HG3	3:C:328:HOH:O	2.20	0.42
1:C:184[B]:ILE:HG21	1:C:186:LYS:HE2	2.00	0.42
1:D:91:TYR:HA	1:D:188:LYS:HG3	2.01	0.42
1:B:221:GLU:HG2	1:B:222:HIS:N	2.33	0.42
1:B:4:SER:HA	1:B:8:LEU:O	2.20	0.42
1:B:19:GLU:O	1:B:125:THR:HA	2.20	0.42
1:C:70:ARG:HA	1:C:70:ARG:NE	2.35	0.42
1:D:111:ILE:HD12	1:D:111:ILE:N	2.34	0.42
1:D:141[A]:MET:HA	1:D:172:LYS:HD2	2.02	0.42
1:D:106:ILE:HD13	1:D:106:ILE:HG21	1.84	0.41
1:D:8:LEU:CA	1:D:12[B]:MET:HE1	2.49	0.41
1:D:70:ARG:NE	1:D:70:ARG:HA	2.35	0.41
1:B:227:ARG:HD2	1:B:227:ARG:HA	1.85	0.41
1:C:93:TRP:NE1	1:C:109:ALA:HB3	2.36	0.41
1:D:21:CYS:HA	1:D:25:HIS:O	2.20	0.41
1:A:221:GLU:HG2	1:A:222:HIS:N	2.36	0.41
1:C:17:ARG:NH1	1:D:157:SER:HA	2.36	0.41
1:A:114:ASN:HD22	1:A:117:GLU:H	1.69	0.40
1:C:41:LYS:HG3	1:C:42:GLN:N	2.37	0.40
1:B:62:SER:HB3	1:B:204:LEU:HD21	2.04	0.40
1:D:8:LEU:HB3	1:D:12[B]:MET:CE	2.52	0.40
1:B:184:ILE:HG21	1:B:186:LYS:HE2	2.02	0.40
1:C:97:PHE:HB3	1:C:180:GLN:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	227/237 (96%)	223 (98%)	4 (2%)	0	100	100
1	B	229/237 (97%)	225 (98%)	4 (2%)	0	100	100
1	C	232/237 (98%)	227 (98%)	5 (2%)	0	100	100
1	D	227/237 (96%)	223 (98%)	4 (2%)	0	100	100
All	All	915/948 (96%)	898 (98%)	17 (2%)	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	200/205 (98%)	198 (99%)	2 (1%)	80	59
1	B	202/205 (98%)	198 (98%)	4 (2%)	60	27
1	C	204/205 (100%)	197 (97%)	7 (3%)	42	11
1	D	200/205 (98%)	192 (96%)	8 (4%)	36	7
All	All	806/820 (98%)	785 (97%)	21 (3%)	54	18

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	114	ASN
1	A	222	HIS

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Mol	Chain	Res	Type
1	B	107[A]	CYS
1	B	107[B]	CYS
1	B	158	GLN
1	B	189	THR
1	C	98	ARG
1	C	107[A]	CYS
1	C	107[B]	CYS
1	C	134	ASP
1	C	179	CYS
1	C	193	GLU
1	C	216	LYS
1	D	4	SER
1	D	134	ASP
1	D	151	LYS
1	D	189	THR
1	D	190	GLU
1	D	192	LYS
1	D	216	LYS
1	D	222	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	114	ASN
1	A	222	HIS
1	B	222	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	XYG	A	66	1	20,21,22	1.02	2 (10%)	22,29,31	1.67	6 (27%)
1	XYG	B	66	1	20,21,22	0.98	2 (10%)	22,29,31	1.79	6 (27%)
1	XYG	C	66	1	20,21,22	1.15	2 (10%)	22,29,31	2.09	8 (36%)
1	XYG	D	66	1	20,21,22	1.13	2 (10%)	22,29,31	2.33	9 (40%)

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
1	XYG	A	66	1	-	0/7/27/28	0/2/2/2
1	XYG	B	66	1	-	0/7/27/28	0/2/2/2
1	XYG	C	66	1	-	0/7/27/28	0/2/2/2
1	XYG	D	66	1	-	0/7/27/28	0/2/2/2

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	66	XYG	CB1-CA1	2.00	1.53	1.50
1	D	66	XYG	CB1-CA1	2.06	1.54	1.50
1	B	66	XYG	CB1-CA1	2.26	1.54	1.50
1	C	66	XYG	CA1-N	2.40	1.32	1.28
1	C	66	XYG	CB1-CA1	2.51	1.54	1.50
1	B	66	XYG	CA1-N	2.58	1.32	1.28
1	D	66	XYG	CA1-N	3.00	1.33	1.28
1	A	66	XYG	CA1-N	3.04	1.33	1.28

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	66	XYG	O2-C2-CA2	-4.81	128.38	130.97
1	B	66	XYG	CA2-C2-N3	-3.79	101.60	103.30
1	A	66	XYG	N3-C1-N2	-3.72	108.53	113.31
1	B	66	XYG	N3-C1-N2	-3.68	108.58	113.31
1	D	66	XYG	N3-C1-N2	-3.63	108.65	113.31
1	C	66	XYG	N3-C1-N2	-3.55	108.75	113.31
1	D	66	XYG	CB2-CA2-C2	-3.55	117.86	122.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	66	XYG	O-C-CA3	-3.30	115.49	126.38
1	D	66	XYG	C2-CA2-N2	-3.17	106.61	108.93
1	D	66	XYG	O-C-CA3	-3.13	116.04	126.38
1	C	66	XYG	CB2-CA2-C2	-3.13	118.39	122.32
1	C	66	XYG	C2-CA2-N2	-2.95	106.77	108.93
1	B	66	XYG	CB2-CA2-C2	-2.94	118.63	122.32
1	A	66	XYG	O-C-CA3	-2.74	117.32	126.38
1	B	66	XYG	O-C-CA3	-2.68	117.53	126.38
1	C	66	XYG	CE1-CD1-CG2	-2.48	118.03	121.28
1	A	66	XYG	C2-CA2-N2	-2.27	107.27	108.93
1	D	66	XYG	CA3-N3-C2	-2.06	119.38	123.94
1	C	66	XYG	CA3-N3-C2	-2.04	119.42	123.94
1	D	66	XYG	CE1-CD1-CG2	-2.04	118.61	121.28
1	A	66	XYG	O2-C2-CA2	2.32	132.22	130.97
1	A	66	XYG	CB2-CA2-N2	2.54	132.63	128.79
1	B	66	XYG	CB2-CA2-N2	2.54	132.63	128.79
1	B	66	XYG	CA2-N2-C1	2.60	109.07	104.21
1	A	66	XYG	CA2-N2-C1	2.84	109.52	104.21
1	C	66	XYG	CA2-N2-C1	3.30	110.37	104.21
1	D	66	XYG	CA2-N2-C1	3.40	110.56	104.21
1	C	66	XYG	CB2-CA2-N2	3.95	134.76	128.79
1	D	66	XYG	CB2-CA2-N2	4.38	135.41	128.79

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

4 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	SO4	A	250	-	4,4,4	0.28	0	6,6,6	0.25	0
2	SO4	B	250	-	4,4,4	0.43	0	6,6,6	0.31	0
2	SO4	C	250	-	4,4,4	0.37	0	6,6,6	0.24	0
2	SO4	D	250	-	4,4,4	0.33	0	6,6,6	0.21	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
2	SO4	A	250	-	-	0/0/0/0	0/0/0/0
2	SO4	B	250	-	-	0/0/0/0	0/0/0/0
2	SO4	C	250	-	-	0/0/0/0	0/0/0/0
2	SO4	D	250	-	-	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.

No monomer is involved in short contacts.

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	225/237 (94%)	0.20	7 (3%) 49 55	12, 19, 32, 56	0
1	B	225/237 (94%)	0.08	4 (1%) 69 74	16, 23, 35, 60	0
1	C	226/237 (95%)	0.13	7 (3%) 49 55	15, 22, 36, 57	0
1	D	225/237 (94%)	1.16	40 (17%) 2 2	18, 32, 54, 81	0
All	All	901/948 (95%)	0.39	58 (6%) 20 23	12, 23, 44, 81	0

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	119	CYS	19.4
1	D	4	SER	10.7
1	D	120	ILE	10.3
1	D	115	VAL	9.1
1	D	116	ARG	9.0
1	D	113	VAL	9.0
1	D	117	GLU	8.5
1	D	156	ASN	8.4
1	A	212	ALA	7.5
1	A	4	SER	7.1
1	D	114	ASN	6.8
1	D	112	THR	6.5
1	D	160	ILE	6.1
1	D	97	PHE	6.0
1	D	118	ASN	5.9
1	D	9	THR	5.8
1	D	155	ILE	5.3
1	B	4	SER	5.0
1	D	189	THR	5.0
1	D	5	ALA	4.7
1	D	8	LEU	4.5

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Mol	Chain	Res	Type	RSRZ
1	D	157	SER	4.2
1	C	97	PHE	4.1
1	D	77	GLU	4.0
1	D	121	TYR	3.9
1	D	88	PRO	3.7
1	A	213	LYS	3.7
1	D	87	CYS	3.7
1	C	116	ARG	3.4
1	D	159	LYS	3.3
1	A	5	ALA	3.2
1	C	179	CYS	3.2
1	D	231	PRO	3.1
1	D	7	GLY	3.1
1	A	97	PHE	3.1
1	D	6	HIS	3.1
1	D	32	ASN	3.0
1	D	11	ASP	3.0
1	B	97	PHE	2.9
1	D	158	GLN	2.9
1	C	3	GLY	2.8
1	D	89	ALA	2.8
1	C	212	ALA	2.7
1	C	4	SER	2.6
1	D	107[A]	CYS	2.5
1	D	38	PHE	2.5
1	B	192	LYS	2.5
1	D	43	PHE	2.5
1	D	86	SER	2.4
1	A	211	ASP	2.4
1	C	98	ARG	2.4
1	A	165	VAL	2.3
1	D	10	ASP	2.3
1	D	90	GLY	2.3
1	D	96	SER	2.2
1	D	179	CYS	2.2
1	B	77	GLU	2.1
1	D	12[A]	MET	2.0

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

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
1	XYG	A	66	20/21	0.96	0.09	-	13,16,18,21	0
1	XYG	C	66	20/21	0.96	0.08	-	17,18,22,25	0
1	XYG	B	66	20/21	0.96	0.06	-	17,18,22,24	0
1	XYG	D	66	20/21	0.95	0.07	-	22,26,30,32	0

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	SO4	A	250	5/5	0.96	0.11	1.13	32,33,42,44	0
2	SO4	D	250	5/5	0.94	0.15	0.71	44,45,48,56	0
2	SO4	B	250	5/5	0.97	0.11	0.50	32,33,37,38	0
2	SO4	C	250	5/5	0.98	0.06	-1.48	25,27,32,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.