



Full wwPDB X-ray Structure Validation Report

Mar 1, 2014 – 04:19 AM GMT

PDB ID : 2C42
Title : Crystal Structure Of Pyruvate-Ferredoxin Oxidoreductase From *Desulfovibrio africanus*
Authors : Cavazza, C.; Contreras-Martel, C.; Pieulle, L.; Chabriere, E.; Hatchikian, E.C.; Fontecilla-Camps, J.C.
Deposited on : 2005-10-14
Resolution : 1.78 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>

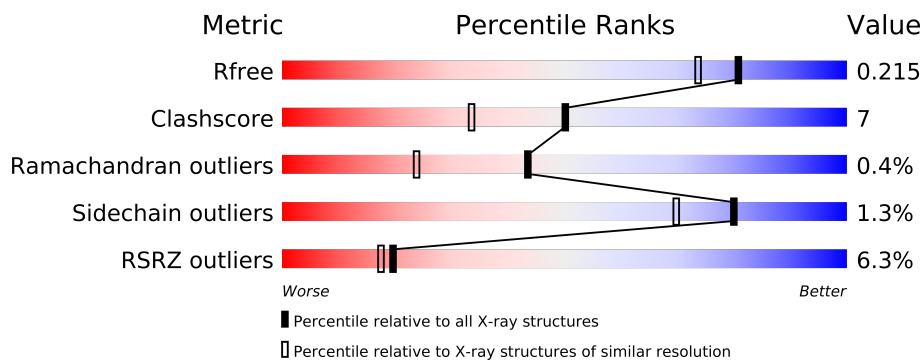
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
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) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.78 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	4987 (1.80-1.76)
Clashscore	79885	6152 (1.80-1.76)
Ramachandran outliers	78287	6074 (1.80-1.76)
Sidechain outliers	78261	6073 (1.80-1.76)
RSRZ outliers	66119	4990 (1.80-1.76)

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	1231	
1	B	1231	

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

Mol	Type	Chain	Res	Geometry	Electron density
4	PYR	B	3237	-	X
6	CA	A	3239	-	X

2 Entry composition i

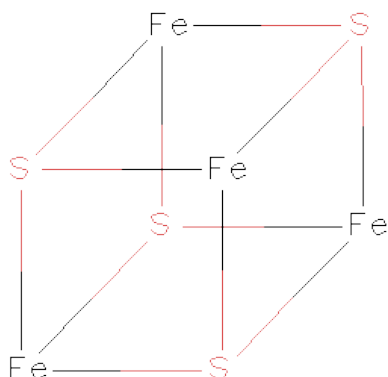
There are 7 unique types of molecules in this entry. The entry contains 20099 atoms, of which 0 are hydrogen and 0 are deuterium.

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 PYRUVATE-FERREDOXIN OXIDOREDUCTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1231	Total	C	N	O	S	0	0	0
			9383	5941	1599	1784	59			
1	B	1231	Total	C	N	O	S	0	0	0
			9383	5941	1599	1784	59			

- Molecule 2 is IRON/SULFUR CLUSTER (three-letter code: SF4) (formula: Fe₄S₄).



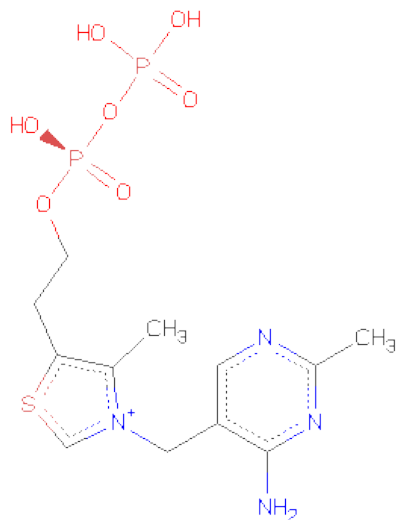
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	Fe	S	0	0
			8	4	4		
2	A	1	Total	Fe	S	0	0
			8	4	4		
2	A	1	Total	Fe	S	0	0
			8	4	4		
2	B	1	Total	Fe	S	0	0
			8	4	4		

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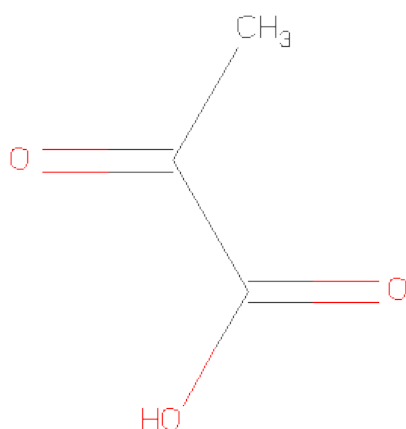
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	Fe	S	0	0
			8	4	4		
2	B	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 3 is THIAMINE DIPHOSPHATE (three-letter code: TPP) (formula: $C_{12}H_{19}N_4O_7P_2S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	S	0	0
			26	12	4	7	2	1		
3	B	1	Total	C	N	O	P	S	0	0
			26	12	4	7	2	1		

- Molecule 4 is PYRUVIC ACID (three-letter code: PYR) (formula: $C_3H_4O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	Mg	0	0
			1	1		
5	A	1	Total	Mg	0	0
			1	1		

- Molecule 6 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	Ca	0	0
			1	1		
6	A	1	Total	Ca	0	0
			1	1		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	550	Total	O	0	0
			550	550		

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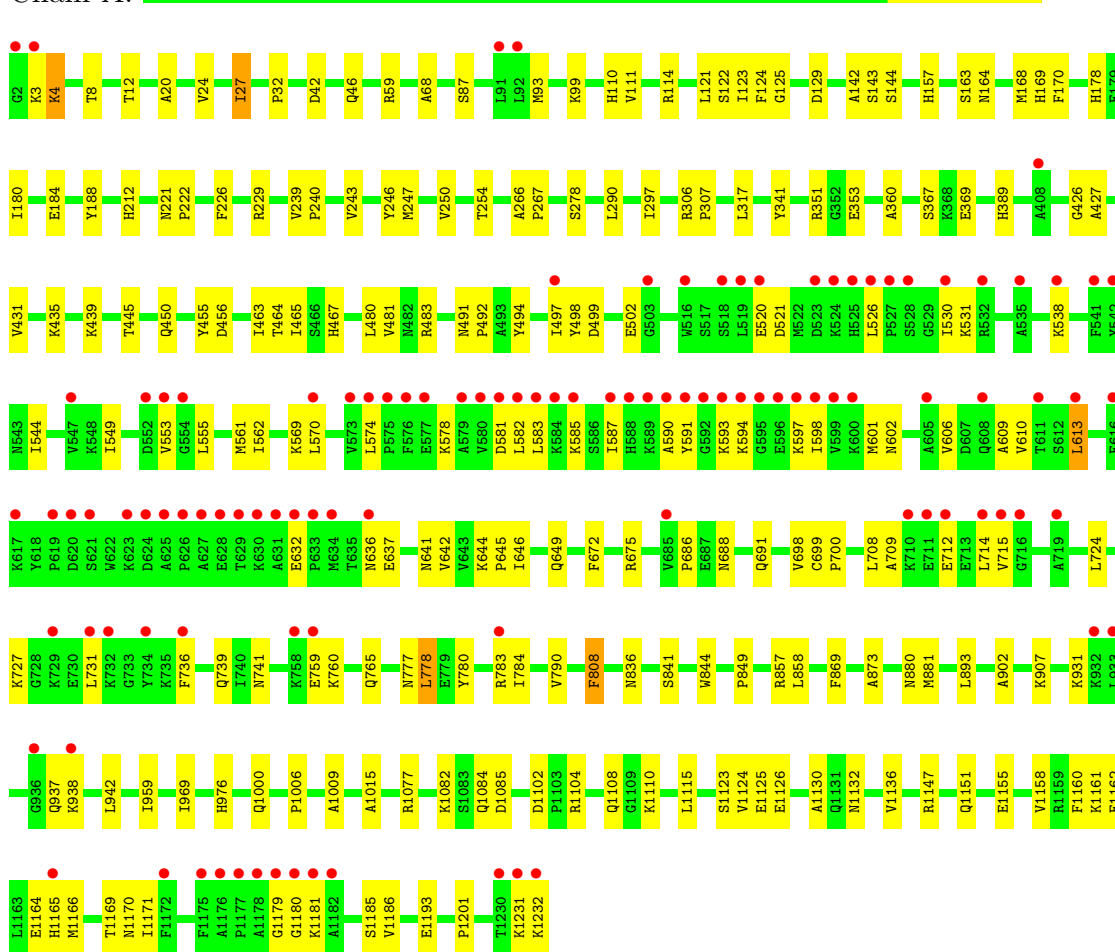
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	667	Total 667	O 667	0	0

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.

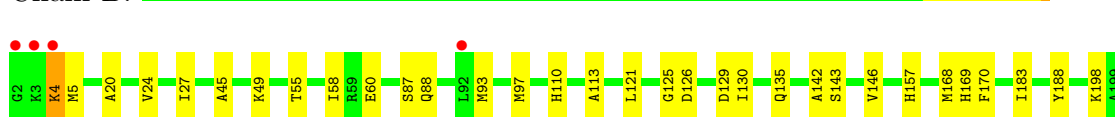
• Molecule 1: PYRUVATE-FERREDOXIN OXIDOREDUCTASE

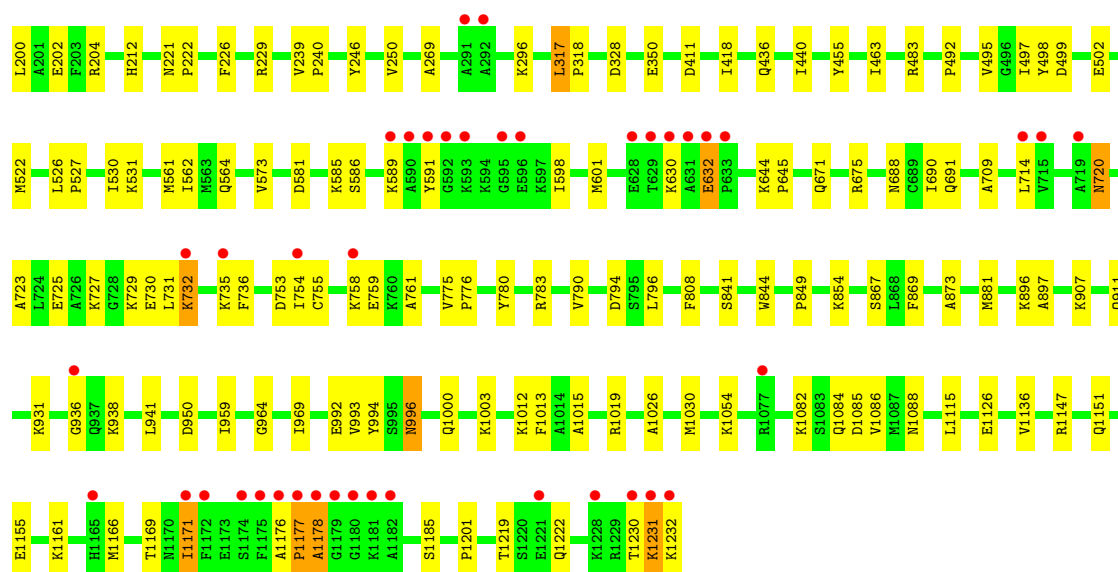
Chain A:



• Molecule 1: PYRUVATE-FERREDOXIN OXIDOREDUCTASE

Chain B:





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	86.30Å 145.98Å 211.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	51.15 – 1.78 51.15 – 1.78	Depositor EDS
% Data completeness (in resolution range)	94.4 (51.15-1.78) 94.5 (51.15-1.78)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.51 (at 1.78Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.195 , 0.219 0.192 , 0.215	Depositor DCC
R_{free} test set	12031 reflections (5.25%)	DCC
Wilson B-factor (Å ²)	18.0	Xtriage
Anisotropy	0.397	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 34.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 241140 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	20099	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 3.86% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, SF4, PYR, CA, TPP

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	0.30	0/9585	0.58	1/12954 (0.0%)
1	B	0.31	0/9585	0.59	1/12954 (0.0%)
All	All	0.31	0/19170	0.59	2/25908 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	125	GLY	N-CA-C	5.62	127.16	113.10
1	A	125	GLY	N-CA-C	5.49	126.83	113.10

There are no chirality outliers.

There are no planarity outliers.

5.2 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 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	9383	0	9262	169	0
1	B	9383	0	9262	134	0
2	A	24	0	0	0	0
2	B	24	0	0	1	0
3	A	26	0	16	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	26	0	16	1	0
4	A	6	0	3	0	0
4	B	6	0	3	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
7	A	550	0	0	8	0
7	B	667	0	0	7	0
All	All	20099	0	18562	280	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 7.

All (280) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:593:LYS:HG3	1:A:594:LYS:H	1.27	0.99
1:A:110:HIS:HD2	1:A:169:HIS:HD2	1.15	0.95
1:B:27:ILE:HD13	1:B:58:ILE:HD11	1.48	0.94
1:B:723:ALA:HB3	1:B:735:LYS:HE2	1.57	0.87
1:B:110:HIS:HD2	1:B:169:HIS:HD2	1.18	0.86
1:B:1176:ALA:HB1	1:B:1177:PRO:HD2	1.59	0.85
1:A:110:HIS:HE1	1:A:157:HIS:HE1	1.23	0.84
1:A:1180:GLY:HA3	1:B:1019:ARG:HH12	1.45	0.81
1:A:1231:LYS:HG3	1:A:1232:LYS:H	1.44	0.81
1:A:351:ARG:HD3	1:A:353:GLU:HB2	1.61	0.80
1:A:110:HIS:HD2	1:A:169:HIS:CD2	2.02	0.78
1:A:594:LYS:HB3	1:A:598:ILE:HD12	1.66	0.76
1:A:110:HIS:CE1	1:A:157:HIS:HE1	2.03	0.76
1:A:1102:ASP:OD1	1:A:1104:ARG:HG2	1.85	0.76
1:B:110:HIS:HD2	1:B:169:HIS:CD2	2.03	0.74
1:B:1231:LYS:HG3	1:B:1232:LYS:H	1.54	0.73
1:A:110:HIS:HE1	1:A:157:HIS:CE1	2.06	0.72
1:A:780:TYR:O	1:A:783:ARG:HG2	1.90	0.71
1:A:644:LYS:HB3	1:A:645:PRO:HD3	1.72	0.71
1:A:1161:LYS:HB3	1:B:1171:ILE:HD12	1.73	0.69
1:A:709:ALA:HB3	1:A:714:LEU:HD11	1.74	0.69
1:B:198:LYS:O	1:B:202:GLU:HG3	1.94	0.68
1:A:1108:GLN:HE21	1:A:1110:LYS:HD2	1.59	0.67
1:B:1230:THR:O	1:B:1232:LYS:HG2	1.95	0.67
1:A:597:LYS:O	1:A:601:MET:HG3	1.95	0.67
1:A:110:HIS:CD2	1:A:169:HIS:HD2	2.06	0.67

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:463:ILE:HD13	1:A:494:TYR:CE1	2.30	0.67
1:A:708:LEU:HD21	1:A:731:LEU:HD22	1.77	0.67
1:A:691:GLN:HG2	1:A:736:PHE:CD2	2.30	0.66
1:A:739:GLN:NE2	1:A:777:ASN:HB3	2.10	0.66
1:A:739:GLN:HE22	1:A:777:ASN:HB3	1.59	0.66
1:B:143:SER:OG	1:B:169:HIS:HE1	1.78	0.65
1:B:463:ILE:HD11	1:B:498:TYR:OH	1.96	0.65
1:A:593:LYS:HG3	1:A:594:LYS:N	2.06	0.65
1:A:1166:MET:O	1:A:1169:THR:HG22	1.97	0.65
1:B:691:GLN:HG2	1:B:736:PHE:CD2	2.32	0.65
1:B:110:HIS:HE1	1:B:157:HIS:NE2	1.95	0.65
1:A:1165:HIS:CD2	1:B:1171:ILE:HG21	2.32	0.64
1:B:581:ASP:O	1:B:585:LYS:HG2	1.96	0.64
1:A:1151:GLN:O	1:A:1155:GLU:HG3	1.97	0.64
1:A:976:HIS:HD2	1:B:1003:LYS:NZ	1.96	0.63
1:A:583:LEU:O	1:A:587:ILE:HG13	1.98	0.63
1:A:1231:LYS:HG3	1:A:1232:LYS:N	2.14	0.62
1:B:110:HIS:CD2	1:B:169:HIS:HD2	2.09	0.62
1:A:1147:ARG:HB3	1:B:1178:ALA:O	1.98	0.62
1:A:445:THR:HG21	1:A:574:LEU:HD21	1.81	0.62
1:A:1185:SER:HB2	1:B:1015:ALA:HB1	1.82	0.61
1:B:992:GLU:O	1:B:993:VAL:HG13	2.01	0.61
1:B:907:LYS:O	1:B:911:GLN:HG3	2.00	0.61
1:B:561:MET:HE1	1:B:564:GLN:HG2	1.83	0.61
1:B:527:PRO:HD2	1:B:530:ILE:HD12	1.81	0.61
1:B:1219:THR:OG1	1:B:1222:GLN:HG3	2.01	0.61
1:A:1181:LYS:HB2	1:A:1181:LYS:NZ	2.16	0.60
1:A:27:ILE:HB	7:A:2040:HOH:O	2.01	0.60
1:A:937:GLN:HG2	1:A:942:LEU:HB3	1.83	0.60
1:A:902:ALA:O	1:A:907:LYS:HE3	2.02	0.60
1:A:636:ASN:ND2	1:A:672:PHE:HE1	2.00	0.59
1:A:526:LEU:HD11	1:A:530:ILE:HG21	1.83	0.59
1:B:780:TYR:HD1	1:B:783:ARG:HH21	1.50	0.59
1:A:731:LEU:HD23	1:A:790:VAL:HG11	1.84	0.59
1:A:526:LEU:O	1:A:531:LYS:HE3	2.01	0.59
1:A:465:ILE:HD11	1:A:649:GLN:NE2	2.17	0.59
1:A:27:ILE:HD13	1:A:59:ARG:O	2.02	0.59
1:A:435:LYS:O	1:A:439:LYS:HD3	2.03	0.59
1:A:110:HIS:CE1	1:A:157:HIS:CE1	2.87	0.59
1:B:714:LEU:HG	1:B:735:LYS:HD3	1.85	0.58
1:A:570:LEU:HD22	1:A:570:LEU:N	2.19	0.58
1:A:544:ILE:HD12	1:A:613:LEU:HD22	1.85	0.58

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:731:LEU:HD23	1:B:790:VAL:HG11	1.85	0.58
1:B:723:ALA:HB1	1:B:735:LYS:HG2	1.86	0.57
1:A:553:VAL:HG23	1:A:555:LEU:HG	1.86	0.57
1:A:143:SER:OG	1:A:169:HIS:HE1	1.87	0.57
1:A:351:ARG:HD2	7:A:2205:HOH:O	2.04	0.57
1:B:931:LYS:HD2	7:B:2525:HOH:O	2.04	0.57
1:A:499:ASP:OD2	1:A:502:GLU:HB2	2.04	0.56
1:B:727:LYS:HA	1:B:727:LYS:HE2	1.88	0.56
1:A:1123:SER:O	1:A:1126:GLU:HG2	2.05	0.56
1:A:931:LYS:HD2	7:A:2421:HOH:O	2.05	0.56
1:B:1166:MET:O	1:B:1169:THR:HG22	2.05	0.56
1:A:731:LEU:CD2	1:A:790:VAL:HG11	2.36	0.56
1:A:3:LYS:NZ	1:A:254:THR:HA	2.20	0.56
1:B:794:ASP:OD1	1:B:1054:LYS:HD2	2.06	0.55
1:A:759:GLU:H	1:A:759:GLU:CD	2.08	0.55
1:A:609:ALA:O	1:A:613:LEU:HD23	2.07	0.55
1:A:497:ILE:HG13	1:A:498:TYR:CD1	2.42	0.55
1:A:142:ALA:HB2	1:A:170:PHE:CZ	2.41	0.55
1:A:27:ILE:H	1:A:27:ILE:HD13	1.72	0.54
1:B:142:ALA:HB2	1:B:170:PHE:CZ	2.43	0.54
1:A:1115:LEU:HD21	1:A:1160:PHE:CZ	2.42	0.54
1:A:1015:ALA:HB1	1:B:1185:SER:HB2	1.88	0.54
1:A:520:GLU:HG3	1:A:521:ASP:N	2.23	0.54
1:A:544:ILE:HD11	1:A:549:ILE:HD12	1.90	0.54
1:B:317:LEU:HD23	1:B:318:PRO:HD2	1.89	0.54
1:A:1158:VAL:O	1:A:1162:GLU:HG3	2.08	0.54
1:A:691:GLN:NE2	1:A:727:LYS:HG3	2.23	0.54
1:B:1232:LYS:NZ	1:B:1232:LYS:HB3	2.23	0.54
1:A:99:LYS:HE3	1:B:867:SER:O	2.08	0.54
1:A:976:HIS:HD2	1:B:1003:LYS:HZ2	1.55	0.53
1:B:411:ASP:HB2	1:B:483:ARG:HD2	1.91	0.53
1:B:492:PRO:O	1:B:495:VAL:HG22	2.07	0.53
1:A:1231:LYS:CG	1:A:1232:LYS:H	2.18	0.53
1:A:857:ARG:HG3	1:A:858:LEU:CD1	2.39	0.53
1:B:1230:THR:O	1:B:1232:LYS:N	2.42	0.52
1:B:135:GLN:H	1:B:135:GLN:NE2	2.08	0.52
1:A:465:ILE:HD11	1:A:649:GLN:HE22	1.73	0.52
1:A:562:ILE:HD12	1:A:562:ILE:N	2.25	0.52
1:A:1132:ASN:O	1:A:1136:VAL:HG12	2.10	0.52
1:A:221:ASN:HB3	1:A:222:PRO:CD	2.39	0.52
1:A:467:HIS:HD2	1:A:481:VAL:H	1.58	0.52
1:A:341:TYR:CD1	1:A:360:ALA:HB2	2.45	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:1161:LYS:CB	1:B:1171:ILE:HD12	2.39	0.51
1:A:1015:ALA:CB	1:B:1185:SER:HB2	2.41	0.51
1:A:606:VAL:O	1:A:610:VAL:HG23	2.10	0.51
1:A:1180:GLY:HA3	1:B:1019:ARG:NH1	2.22	0.51
1:B:841:SER:HA	1:B:844:TRP:CE2	2.45	0.51
1:B:675:ARG:HD3	7:B:2398:HOH:O	2.11	0.51
1:B:1231:LYS:HG3	1:B:1232:LYS:N	2.23	0.51
1:A:637:GLU:HG3	1:A:641:ASN:ND2	2.26	0.51
1:B:731:LEU:CD2	1:B:790:VAL:HG11	2.41	0.51
1:B:1177:PRO:O	1:B:1178:ALA:HB2	2.11	0.50
1:B:499:ASP:OD2	1:B:502:GLU:HB2	2.11	0.50
1:A:455:TYR:HB2	1:B:1201:PRO:HG3	1.94	0.50
1:B:586:SER:O	1:B:589:LYS:HB3	2.11	0.50
1:A:1006:PRO:HG2	1:A:1009:ALA:HB2	1.94	0.50
1:B:709:ALA:HB3	1:B:714:LEU:HD21	1.94	0.50
1:B:725:GLU:O	1:B:727:LYS:HE3	2.10	0.50
1:A:157:HIS:HD2	7:A:2004:HOH:O	1.94	0.50
1:B:729:LYS:O	1:B:732:LYS:HG3	2.11	0.50
1:B:20:ALA:HB2	1:B:188:TYR:CZ	2.47	0.50
1:A:561:MET:O	1:A:561:MET:HE2	2.12	0.50
1:A:581:ASP:OD2	1:A:585:LYS:HE3	2.12	0.49
1:A:741:ASN:CG	1:A:778:LEU:HD11	2.33	0.49
1:A:8:THR:OG1	1:A:12:THR:HB	2.11	0.49
1:A:1186:VAL:HG21	1:B:1136:VAL:HG22	1.94	0.49
1:A:494:TYR:HD2	1:A:497:ILE:HD11	1.76	0.49
1:B:873:ALA:HA	1:B:959:ILE:HD13	1.94	0.49
1:A:686:PRO:HB2	1:A:724:LEU:HD21	1.95	0.49
1:B:269:ALA:HA	1:B:296:LYS:HB3	1.95	0.49
1:A:1181:LYS:HZ2	1:A:1181:LYS:HB2	1.76	0.49
1:A:229:ARG:HD2	7:B:2059:HOH:O	2.13	0.49
1:A:389:HIS:HE1	1:B:350:GLU:OE1	1.96	0.48
1:A:578:LYS:O	1:A:582:LEU:HD13	2.14	0.48
1:A:467:HIS:CD2	1:A:481:VAL:H	2.32	0.48
1:B:239:VAL:HB	1:B:240:PRO:HD3	1.95	0.48
1:A:1232:LYS:HZ2	1:A:1232:LYS:HB3	1.79	0.48
1:A:1180:GLY:O	1:A:1181:LYS:HB2	2.14	0.48
1:B:1232:LYS:HB3	1:B:1232:LYS:HZ2	1.79	0.48
1:A:389:HIS:HD2	7:A:2059:HOH:O	1.96	0.48
1:A:538:LYS:HD2	1:A:538:LYS:N	2.29	0.48
1:B:229:ARG:HD2	7:B:2125:HOH:O	2.13	0.48
1:B:526:LEU:O	1:B:531:LYS:HE3	2.14	0.47
1:A:593:LYS:CG	1:A:594:LYS:H	2.05	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:636:ASN:ND2	1:A:672:PHE:CE1	2.80	0.47
1:B:869:PHE:CE2	1:B:969:ILE:HG21	2.48	0.47
1:B:775:VAL:N	1:B:776:PRO:HD2	2.29	0.47
1:B:630:LYS:O	1:B:630:LYS:HG3	2.13	0.47
1:A:1232:LYS:NZ	1:A:1232:LYS:HB3	2.28	0.47
1:B:1126:GLU:HG3	7:B:2618:HOH:O	2.13	0.47
1:A:1082:LYS:HE2	1:A:1085:ASP:OD1	2.14	0.47
1:A:741:ASN:ND2	1:A:778:LEU:HD11	2.29	0.47
1:B:93:MET:O	1:B:97:MET:HG3	2.15	0.47
1:B:1231:LYS:O	1:B:1232:LYS:HB2	2.14	0.47
1:A:3:LYS:HZ3	1:A:254:THR:HA	1.80	0.47
1:B:730:GLU:N	1:B:730:GLU:OE1	2.47	0.47
1:B:436:GLN:O	1:B:440:ILE:HG13	2.15	0.47
1:B:4:LYS:HE2	1:B:4:LYS:HA	1.97	0.46
1:B:221:ASN:HB3	1:B:222:PRO:CD	2.45	0.46
1:A:212:HIS:HE1	1:B:950:ASP:OD2	1.97	0.46
1:B:714:LEU:N	1:B:714:LEU:HD22	2.31	0.46
1:A:1201:PRO:HG3	1:B:455:TYR:HB2	1.97	0.46
1:B:49:LYS:NZ	1:B:55:THR:HG23	2.30	0.46
1:A:121:LEU:C	1:A:121:LEU:HD23	2.34	0.46
1:A:937:GLN:C	1:A:938:LYS:HD2	2.36	0.46
1:A:246:TYR:O	1:A:250:VAL:HG23	2.15	0.46
1:A:675:ARG:HD3	7:A:2314:HOH:O	2.13	0.46
1:B:755:CYS:SG	1:B:761:ALA:HB3	2.56	0.46
1:B:897:ALA:HA	1:B:941:LEU:HD23	1.96	0.46
1:B:497:ILE:HG13	1:B:498:TYR:CD1	2.50	0.46
1:A:124:PHE:HB3	1:A:367:SER:HB2	1.97	0.46
1:A:569:LYS:HB3	1:A:570:LEU:HD22	1.98	0.46
1:A:869:PHE:CE2	1:A:969:ILE:HG21	2.50	0.46
1:A:699:CYS:HA	1:A:700:PRO:HD3	1.82	0.46
1:B:723:ALA:CB	1:B:735:LYS:HG2	2.46	0.46
1:A:688:ASN:HB3	1:A:759:GLU:O	2.16	0.46
1:A:239:VAL:HB	1:A:240:PRO:HD3	1.97	0.46
1:A:1077:ARG:HG2	1:A:1130:ALA:O	2.16	0.45
1:A:976:HIS:CD2	1:B:1003:LYS:HZ2	2.34	0.45
1:A:808:PHE:CD2	1:A:808:PHE:N	2.85	0.45
1:A:1171:ILE:HD12	1:B:1161:LYS:HD3	1.98	0.45
1:B:996:ASN:ND2	3:B:3236:TPP:S1	2.90	0.45
1:B:688:ASN:HB3	1:B:759:GLU:O	2.17	0.45
1:A:121:LEU:HD23	1:A:122:SER:N	2.31	0.45
1:A:369:GLU:OE1	1:A:480:LEU:HG	2.17	0.45
1:B:27:ILE:HD12	1:B:1013:PHE:CE2	2.51	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:522:MET:SD	1:B:526:LEU:HG	2.57	0.45
1:A:163:SER:O	1:A:164:ASN:HB2	2.17	0.45
1:A:32:PRO:HB2	1:A:178:HIS:CE1	2.51	0.45
1:A:456:ASP:OD1	1:A:463:ILE:HG22	2.17	0.45
1:A:431:VAL:CG2	1:A:464:THR:HG21	2.46	0.45
1:A:426:GLY:O	1:A:427:ALA:HB3	2.16	0.45
1:B:27:ILE:CD1	1:B:1013:PHE:CE2	3.00	0.45
1:A:42:ASP:O	1:A:46:GLN:HG3	2.17	0.45
1:B:644:LYS:HB3	1:B:645:PRO:HD3	1.99	0.45
1:B:691:GLN:NE2	1:B:727:LYS:HG2	2.32	0.44
1:A:3:LYS:O	1:A:4:LYS:HB2	2.17	0.44
1:B:897:ALA:CA	1:B:941:LEU:HD23	2.47	0.44
1:B:110:HIS:CE1	1:B:157:HIS:NE2	2.79	0.44
1:A:709:ALA:HB2	1:A:784:ILE:HG21	1.98	0.44
1:A:602:ASN:O	1:A:606:VAL:HG23	2.18	0.44
1:A:1124:VAL:HG13	1:A:1125:GLU:N	2.33	0.44
1:A:881:MET:HE1	1:B:24:VAL:HG13	1.99	0.44
1:A:110:HIS:CD2	1:A:169:HIS:CD2	2.92	0.44
1:A:544:ILE:HD12	1:A:613:LEU:CD2	2.47	0.44
1:B:1082:LYS:O	1:B:1086:VAL:HG23	2.18	0.44
1:B:796:LEU:HD23	1:B:796:LEU:C	2.37	0.44
1:A:180:ILE:O	1:A:450:GLN:HA	2.18	0.44
1:A:841:SER:HA	1:A:844:TRP:CE2	2.53	0.44
1:A:594:LYS:HB3	1:A:598:ILE:CD1	2.43	0.43
1:B:246:TYR:O	1:B:250:VAL:HG23	2.18	0.43
1:B:1147:ARG:HG3	1:B:1147:ARG:HH11	1.84	0.43
1:A:68:ALA:HB2	1:A:93:MET:HG2	2.01	0.43
1:B:130:ILE:HB	1:B:168:MET:HE1	1.99	0.43
1:A:698:VAL:HG13	1:A:1084:GLN:NE2	2.32	0.43
1:A:712:GLU:O	1:A:715:VAL:HG23	2.19	0.43
1:B:690:ILE:HG12	2:B:3233:SF4:S2	2.58	0.43
1:B:671:GLN:NE2	1:B:854:LYS:HD2	2.33	0.43
1:B:754:ILE:HD13	1:B:1084:GLN:HB2	1.99	0.43
1:A:20:ALA:HB2	1:A:188:TYR:CZ	2.54	0.43
1:A:114:ARG:NE	1:A:123:ILE:HA	2.33	0.43
1:A:306:ARG:HA	1:A:307:PRO:C	2.37	0.43
1:A:494:TYR:HA	1:A:497:ILE:CG1	2.49	0.43
1:A:1160:PHE:O	1:A:1164:GLU:HG3	2.18	0.43
1:A:873:ALA:HA	1:A:959:ILE:HD13	2.01	0.43
1:B:936:GLY:O	1:B:938:LYS:HG2	2.19	0.43
1:B:964:GLY:HA2	1:B:994:TYR:HE1	1.84	0.43
1:B:753:ASP:OD2	1:B:1085:ASP:OD2	2.37	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:200:LEU:O	1:B:204:ARG:HG2	2.19	0.42
1:A:184:GLU:HG2	7:A:2079:HOH:O	2.18	0.42
1:A:87:SER:HA	1:A:129:ASP:HB3	2.00	0.42
1:B:1012:LYS:O	1:B:1013:PHE:HB2	2.20	0.42
1:A:698:VAL:O	1:A:698:VAL:HG12	2.19	0.42
1:B:87:SER:HA	1:B:129:ASP:HB3	2.01	0.42
1:A:494:TYR:HA	1:A:497:ILE:HG12	2.02	0.42
1:B:562:ILE:HD12	1:B:562:ILE:N	2.35	0.42
1:A:317:LEU:HD11	7:A:2186:HOH:O	2.19	0.42
1:B:758:LYS:HB2	1:B:758:LYS:HE3	1.89	0.42
1:A:714:LEU:N	1:A:714:LEU:HD12	2.35	0.42
1:A:221:ASN:HB3	1:A:222:PRO:HD2	2.01	0.41
1:B:20:ALA:HB2	1:B:188:TYR:CE1	2.55	0.41
1:A:483:ARG:HH11	1:A:483:ARG:HG2	1.85	0.41
1:A:144:SER:O	1:A:278:SER:HA	2.20	0.41
1:B:1026:ALA:O	1:B:1030:MET:HG3	2.20	0.41
1:B:146:VAL:HG12	1:B:183:ILE:HD13	2.02	0.41
1:B:896:LYS:HB3	1:B:941:LEU:HD21	2.02	0.41
1:B:754:ILE:HG12	7:B:2592:HOH:O	2.20	0.41
1:A:976:HIS:HE1	1:B:60:GLU:O	2.04	0.41
1:B:212:HIS:HD2	7:B:2104:HOH:O	2.04	0.41
1:A:24:VAL:HG13	1:B:881:MET:HE2	2.01	0.41
1:A:266:ALA:HA	1:A:267:PRO:HD3	1.90	0.41
1:A:1185:SER:HB3	1:B:45:ALA:HB3	2.03	0.41
1:B:87:SER:OG	1:B:88:GLN:N	2.54	0.41
1:B:630:LYS:C	1:B:632:GLU:H	2.24	0.41
1:A:111:VAL:HG21	1:A:168:MET:HE2	2.02	0.41
1:A:1193:GLU:N	1:A:1193:GLU:OE2	2.54	0.41
1:A:243:VAL:O	1:A:247:MET:HG3	2.21	0.41
1:B:581:ASP:OD2	1:B:585:LYS:HE3	2.20	0.41
1:A:642:VAL:O	1:A:646:ILE:HG13	2.21	0.41
1:B:121:LEU:C	1:B:121:LEU:HD23	2.41	0.41
1:B:113:ALA:HB1	1:B:126:ASP:O	2.20	0.41
1:B:418:ILE:HD12	1:B:573:VAL:HA	2.03	0.41
1:A:290:LEU:HB2	1:A:297:ILE:HD11	2.03	0.41
1:B:1176:ALA:HB1	1:B:1177:PRO:CD	2.43	0.40
1:A:491:ASN:HA	1:A:492:PRO:HD2	1.96	0.40
1:B:4:LYS:HE2	1:B:5:MET:N	2.36	0.40
1:B:598:ILE:HD13	1:B:601:MET:CE	2.52	0.40
1:B:992:GLU:O	1:B:993:VAL:CG1	2.69	0.40
1:A:590:ALA:O	1:A:591:TYR:C	2.60	0.40
1:B:1151:GLN:O	1:B:1155:GLU:HG3	2.21	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:130:ILE:HB	1:B:168:MET:CE	2.51	0.40
1:B:720:ASN:H	1:B:720:ASN:ND2	2.20	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	1229/1231 (100%)	1174 (96%)	51 (4%)	4 (0%)	50	30
1	B	1229/1231 (100%)	1191 (97%)	33 (3%)	5 (0%)	43	24
All	All	2458/2462 (100%)	2365 (96%)	84 (3%)	9 (0%)	43	24

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	1231	LYS
1	A	760	LYS
1	A	613	LEU
1	B	591	TYR
1	B	1178	ALA
1	A	4	LYS
1	B	732	LYS
1	A	1179	GLY
1	B	1177	PRO

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	978/978 (100%)	966 (99%)	12 (1%)	82	72
1	B	978/978 (100%)	965 (99%)	13 (1%)	80	70
All	All	1956/1956 (100%)	1931 (99%)	25 (1%)	80	70

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	ILE
1	A	226	PHE
1	A	632	GLU
1	A	765	GLN
1	A	778	LEU
1	A	808	PHE
1	A	836	ASN
1	A	849	PRO
1	A	880	ASN
1	A	893	LEU
1	A	1000	GLN
1	A	1170	ASN
1	B	4	LYS
1	B	226	PHE
1	B	317	LEU
1	B	328	ASP
1	B	632	GLU
1	B	720	ASN
1	B	808	PHE
1	B	849	PRO
1	B	996	ASN
1	B	1000	GLN
1	B	1088	ASN
1	B	1115	LEU
1	B	1171	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (42) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	46	GLN
1	A	110	HIS
1	A	157	HIS
1	A	169	HIS
1	A	212	HIS

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Mol	Chain	Res	Type
1	A	389	HIS
1	A	421	GLN
1	A	434	ASN
1	A	467	HIS
1	A	536	ASN
1	A	543	ASN
1	A	636	ASN
1	A	683	GLN
1	A	688	ASN
1	A	739	GLN
1	A	765	GLN
1	A	777	ASN
1	A	937	GLN
1	A	976	HIS
1	A	1000	GLN
1	A	1108	GLN
1	A	1165	HIS
1	A	1170	ASN
1	B	46	GLN
1	B	54	GLN
1	B	110	HIS
1	B	135	GLN
1	B	164	ASN
1	B	169	HIS
1	B	197	GLN
1	B	212	HIS
1	B	220	GLN
1	B	434	ASN
1	B	683	GLN
1	B	688	ASN
1	B	720	ASN
1	B	836	ASN
1	B	866	ASN
1	B	996	ASN
1	B	1000	GLN
1	B	1088	ASN
1	B	1108	GLN

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 14 ligands modelled in this entry, 4 are monoatomic - leaving 10 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	$\# Z > 2$
2	SF4	A	3233	1	12,12,12	4.43	12 (100%)	0,24,24	0.00	-
2	SF4	A	3234	1	12,12,12	4.20	11 (91%)	0,24,24	0.00	-
2	SF4	A	3235	1	12,12,12	4.36	10 (83%)	0,24,24	0.00	-
3	TPP	A	3236	5	27,27,27	2.93	10 (37%)	40,40,40	1.76	9 (22%)
4	PYR	A	3237	-	5,5,5	1.07	0	6,6,6	2.36	3 (50%)
2	SF4	B	3233	1	12,12,12	4.47	10 (83%)	0,24,24	0.00	-
2	SF4	B	3234	1	12,12,12	4.95	12 (100%)	0,24,24	0.00	-
2	SF4	B	3235	1	12,12,12	5.11	11 (91%)	0,24,24	0.00	-
3	TPP	B	3236	5	27,27,27	2.48	9 (33%)	40,40,40	1.99	8 (20%)
4	PYR	B	3237	-	5,5,5	1.04	0	6,6,6	2.34	3 (50%)

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	SF4	A	3233	1	-	0/0/48/48	0/0/5/5
2	SF4	A	3234	1	-	0/0/48/48	0/0/5/5
2	SF4	A	3235	1	-	0/0/48/48	0/0/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	TPP	A	3236	5	-	0/17/17/17	0/2/2/2
4	PYR	A	3237	-	-	0/4/4/4	0/0/0/0
2	SF4	B	3233	1	-	0/0/48/48	0/0/5/5
2	SF4	B	3234	1	-	0/0/48/48	0/0/5/5
2	SF4	B	3235	1	-	0/0/48/48	0/0/5/5
3	TPP	B	3236	5	-	0/17/17/17	0/2/2/2
4	PYR	B	3237	-	-	0/4/4/4	0/0/0/0

All (85) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	3235	SF4	S3-FE1	-10.11	2.26	2.33
3	A	3236	TPP	C4-N3	9.99	1.50	1.39
2	B	3235	SF4	S4-FE2	-8.17	2.27	2.33
2	B	3234	SF4	S3-FE1	-8.05	2.27	2.33
2	B	3234	SF4	S3-FE2	-7.34	2.28	2.33
3	B	3236	TPP	C4-N3	7.15	1.47	1.39
2	A	3233	SF4	S1-FE2	-6.74	2.28	2.33
2	B	3234	SF4	S2-FE4	-6.68	2.28	2.33
2	A	3234	SF4	S3-FE1	-6.63	2.28	2.33
2	B	3233	SF4	S1-FE2	-6.57	2.28	2.33
2	B	3233	SF4	S4-FE3	-6.44	2.28	2.33
2	A	3235	SF4	S3-FE2	-6.24	2.29	2.33
2	A	3233	SF4	S4-FE3	-6.15	2.29	2.33
2	A	3235	SF4	S3-FE1	-5.98	2.29	2.33
2	B	3233	SF4	S2-FE1	-5.87	2.29	2.33
2	B	3235	SF4	S1-FE2	-5.86	2.29	2.33
2	B	3233	SF4	S3-FE4	-5.68	2.29	2.33
2	B	3233	SF4	S1-FE4	-5.57	2.29	2.33
2	A	3234	SF4	S3-FE4	-5.56	2.29	2.33
2	A	3233	SF4	S1-FE4	-5.51	2.29	2.33
2	A	3235	SF4	S4-FE2	-5.40	2.29	2.33
2	A	3234	SF4	S3-FE2	-5.35	2.29	2.33
2	A	3234	SF4	S1-FE2	-5.22	2.29	2.33
2	A	3235	SF4	S4-FE3	-5.12	2.29	2.33
2	B	3235	SF4	S1-FE3	-5.12	2.29	2.33
2	A	3235	SF4	S3-FE4	-5.11	2.29	2.33
2	B	3234	SF4	S1-FE4	-5.10	2.29	2.33
2	A	3233	SF4	S4-FE2	-5.00	2.29	2.33
3	B	3236	TPP	C4'-N3'	4.95	1.43	1.35
2	A	3235	SF4	S1-FE2	-4.92	2.29	2.33
2	B	3234	SF4	S3-FE4	-4.86	2.30	2.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	3236	TPP	C5'-C4'	4.82	1.54	1.42
2	A	3233	SF4	S1-FE3	-4.75	2.30	2.33
2	B	3234	SF4	S1-FE2	-4.58	2.30	2.33
3	B	3236	TPP	C2'-N1'	4.53	1.42	1.34
2	A	3233	SF4	S3-FE1	-4.41	2.30	2.33
2	A	3234	SF4	S2-FE1	-4.36	2.30	2.33
2	B	3235	SF4	S3-FE2	-4.18	2.30	2.33
3	A	3236	TPP	PB-O3A	4.17	1.67	1.60
2	B	3233	SF4	S4-FE1	-4.09	2.30	2.33
2	A	3234	SF4	S4-FE2	-4.09	2.30	2.33
2	B	3234	SF4	S4-FE3	-3.99	2.30	2.33
2	B	3235	SF4	S3-FE4	-3.99	2.30	2.33
2	B	3235	SF4	S2-FE1	-3.83	2.30	2.33
2	A	3234	SF4	S1-FE3	-3.81	2.30	2.33
2	A	3233	SF4	S2-FE1	-3.79	2.30	2.33
2	B	3235	SF4	S4-FE1	-3.79	2.30	2.33
3	A	3236	TPP	C2'-N1'	3.79	1.41	1.34
2	B	3234	SF4	S4-FE2	-3.76	2.30	2.33
3	A	3236	TPP	C6'-C5'	3.76	1.46	1.37
2	B	3234	SF4	S4-FE1	-3.56	2.30	2.33
3	A	3236	TPP	C6'-N1'	3.54	1.42	1.34
3	A	3236	TPP	C6-C5	3.54	1.55	1.51
2	A	3235	SF4	S4-FE1	-3.52	2.30	2.33
3	A	3236	TPP	C5'-C4'	3.51	1.51	1.42
3	B	3236	TPP	C2'-N3'	3.39	1.40	1.34
2	A	3235	SF4	S2-FE1	-3.38	2.31	2.33
2	B	3233	SF4	S4-FE2	-3.34	2.31	2.33
2	A	3233	SF4	S3-FE4	-3.32	2.31	2.33
3	A	3236	TPP	C2'-N3'	3.30	1.40	1.34
3	A	3236	TPP	C4'-N3'	3.24	1.40	1.35
2	B	3233	SF4	S1-FE3	-3.16	2.31	2.33
2	A	3234	SF4	S2-FE4	-3.14	2.31	2.33
2	B	3233	SF4	S3-FE2	-3.12	2.31	2.33
2	B	3235	SF4	S4-FE3	-3.11	2.31	2.33
2	A	3235	SF4	S1-FE3	-3.11	2.31	2.33
2	A	3234	SF4	S4-FE3	-3.01	2.31	2.33
2	A	3235	SF4	S1-FE4	-2.94	2.31	2.33
2	A	3233	SF4	S4-FE1	-2.92	2.31	2.33
2	B	3233	SF4	S2-FE3	-2.80	2.31	2.33
3	B	3236	TPP	C6-C5	2.76	1.54	1.51
2	A	3233	SF4	S2-FE3	-2.74	2.31	2.33
2	A	3233	SF4	S2-FE4	-2.65	2.31	2.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	3235	SF4	S1-FE4	-2.63	2.31	2.33
2	B	3234	SF4	S2-FE1	-2.58	2.31	2.33
2	A	3233	SF4	S3-FE2	-2.56	2.31	2.33
2	B	3234	SF4	S1-FE3	-2.53	2.31	2.33
2	A	3234	SF4	S4-FE1	-2.37	2.31	2.33
2	A	3234	SF4	S1-FE4	-2.34	2.31	2.33
3	B	3236	TPP	O7-C7	-2.19	1.35	1.44
2	B	3234	SF4	S2-FE3	-2.11	2.31	2.33
3	B	3236	TPP	PB-O3B	-2.10	1.47	1.54
3	B	3236	TPP	C6'-C5'	2.05	1.42	1.37
3	A	3236	TPP	C7'-C5'	-2.02	1.46	1.51
2	B	3235	SF4	S2-FE4	-2.02	2.31	2.33

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	3236	TPP	C4-C5-S1	6.22	115.01	109.93
3	A	3236	TPP	C4-C5-S1	5.60	114.50	109.93
3	B	3236	TPP	C6-C5-S1	-4.52	115.96	122.28
3	B	3236	TPP	C6'-N1'-C2'	4.49	123.56	115.68
3	A	3236	TPP	C6-C5-S1	-4.26	116.33	122.28
4	A	3237	PYR	O2-C1-C2	4.02	123.22	114.37
4	B	3237	PYR	O2-C1-C2	3.94	123.06	114.37
3	B	3236	TPP	CM4-C4-N3	3.50	127.10	122.53
4	A	3237	PYR	O1-C1-C2	-3.15	111.65	120.73
4	B	3237	PYR	O1-C1-C2	-3.04	111.98	120.73
3	B	3236	TPP	N1'-C2'-N3'	-2.81	120.58	125.65
3	B	3236	TPP	CM2-C2'-N3'	2.79	122.32	117.31
3	A	3236	TPP	N4'-C4'-N3'	-2.77	112.84	116.88
3	A	3236	TPP	C6'-N1'-C2'	2.56	120.18	115.68
3	A	3236	TPP	N1'-C2'-N3'	-2.50	121.14	125.65
3	A	3236	TPP	CM4-C4-N3	2.42	125.69	122.53
3	A	3236	TPP	C6-C5-C4	2.37	129.17	127.44
3	B	3236	TPP	C7-C6-C5	2.37	118.84	112.48
3	A	3236	TPP	C7-C6-C5	2.35	118.81	112.48
4	B	3237	PYR	O3-C2-C1	2.32	122.95	118.75
3	A	3236	TPP	CM2-C2'-N1'	2.29	119.75	117.02
3	B	3236	TPP	C6-C5-C4	2.19	129.03	127.44
4	A	3237	PYR	O3-C2-C1	2.14	122.62	118.75

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	1231/1231 (100%)	0.43	109 (8%) 10 9	9, 23, 64, 106	0
1	B	1231/1231 (100%)	0.15	45 (3%) 39 37	10, 19, 48, 96	0
All	All	2462/2462 (100%)	0.29	154 (6%) 19 17	9, 21, 58, 106	0

All (154) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	594	LYS	9.1
1	A	576	PHE	8.9
1	A	630	LYS	8.2
1	A	1182	ALA	8.2
1	A	1179	GLY	8.1
1	A	631	ALA	7.8
1	A	1178	ALA	7.4
1	B	631	ALA	7.1
1	A	593	LYS	7.1
1	B	1177	PRO	7.0
1	A	1232	LYS	7.0
1	A	591	TYR	6.9
1	B	1232	LYS	6.8
1	A	1231	LYS	6.6
1	A	629	THR	6.5
1	A	626	PRO	6.1
1	B	732	LYS	6.0
1	B	2	GLY	5.8
1	A	582	LEU	5.7
1	A	628	GLU	5.7
1	B	1176	ALA	5.6
1	B	1178	ALA	5.6
1	A	715	VAL	5.5
1	B	629	THR	5.5

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Mol	Chain	Res	Type	RSRZ
1	A	592	GLY	5.4
1	A	588	HIS	5.2
1	A	574	LEU	5.2
1	A	596	GLU	5.0
1	A	575	PRO	5.0
1	A	595	GLY	4.8
1	B	3	LYS	4.7
1	B	1175	PHE	4.6
1	B	590	ALA	4.6
1	A	627	ALA	4.5
1	B	630	LYS	4.5
1	A	597	LYS	4.4
1	A	1175	PHE	4.4
1	A	633	PRO	4.3
1	A	577	GLU	4.3
1	B	628	GLU	4.3
1	A	589	LYS	4.3
1	A	1181	LYS	4.2
1	A	621	SER	4.1
1	A	590	ALA	4.1
1	A	1180	GLY	4.0
1	B	592	GLY	4.0
1	A	632	GLU	3.9
1	B	715	VAL	3.9
1	A	554	GLY	3.8
1	A	598	ILE	3.7
1	B	1231	LYS	3.7
1	A	1230	THR	3.7
1	A	2	GLY	3.7
1	A	619	PRO	3.7
1	A	719	ALA	3.7
1	A	1165	HIS	3.6
1	A	3	LYS	3.6
1	A	580	VAL	3.4
1	A	613	LEU	3.4
1	B	632	GLU	3.4
1	B	1171	ILE	3.4
1	A	714	LEU	3.4
1	A	936	GLY	3.3
1	A	620	ASP	3.3
1	A	732	LYS	3.3
1	B	591	TYR	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	1228	LYS	3.3
1	A	716	GLY	3.2
1	A	585	LYS	3.2
1	B	4	LYS	3.2
1	B	593	LYS	3.2
1	A	520	GLU	3.2
1	A	758	LYS	3.2
1	B	1165	HIS	3.1
1	A	710	LYS	3.1
1	A	623	LYS	3.1
1	A	1177	PRO	3.1
1	B	596	GLU	3.1
1	A	599	VAL	3.0
1	A	547	VAL	3.0
1	B	1174	SER	3.0
1	B	589	LYS	3.0
1	A	519	LEU	3.0
1	A	525	HIS	3.0
1	B	719	ALA	2.9
1	A	524	LYS	2.9
1	A	624	ASP	2.9
1	A	1176	ALA	2.8
1	A	938	LYS	2.8
1	B	595	GLY	2.8
1	B	1230	THR	2.8
1	A	518	SER	2.8
1	A	625	ALA	2.8
1	A	729	LYS	2.8
1	A	573	VAL	2.7
1	A	552	ASP	2.7
1	A	553	VAL	2.7
1	A	783	ARG	2.7
1	A	541	PHE	2.7
1	A	759	GLU	2.6
1	A	523	ASP	2.6
1	B	1182	ALA	2.6
1	B	1179	GLY	2.6
1	B	1180	GLY	2.6
1	B	633	PRO	2.6
1	A	526	LEU	2.5
1	A	584	LYS	2.5
1	B	1181	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	497	ILE	2.5
1	A	605	ALA	2.5
1	B	754	ILE	2.5
1	A	634	MET	2.5
1	A	528	SER	2.5
1	A	608	GLN	2.5
1	A	538	LYS	2.5
1	A	581	ASP	2.4
1	A	516	TRP	2.4
1	B	714	LEU	2.4
1	B	1077	ARG	2.4
1	A	616	PHE	2.4
1	A	1172	PHE	2.4
1	A	535	ALA	2.3
1	A	408	ALA	2.3
1	A	503	GLY	2.3
1	A	736	PHE	2.3
1	A	731	LEU	2.3
1	A	92	LEU	2.2
1	B	758	LYS	2.2
1	A	685	VAL	2.2
1	B	1172	PHE	2.2
1	A	91	LEU	2.2
1	B	92	LEU	2.2
1	A	636	ASN	2.2
1	A	542	TYR	2.2
1	A	611	THR	2.2
1	B	292	ALA	2.2
1	B	1221	GLU	2.2
1	A	734	TYR	2.2
1	A	587	ILE	2.2
1	A	532	ARG	2.1
1	A	527	PRO	2.1
1	A	712	GLU	2.1
1	A	530	ILE	2.1
1	A	600	LYS	2.1
1	B	291	ALA	2.1
1	A	617	LYS	2.1
1	B	936	GLY	2.0
1	A	711	GLU	2.0
1	A	932	LYS	2.0
1	B	735	LYS	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	579	ALA	2.0
1	A	570	LEU	2.0
1	A	933	LEU	2.0
1	A	583	LEU	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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
6	CA	A	3239	1/1	0.13	5.35	40,40,40,40	0
4	PYR	B	3237	6/6	0.18	3.06	16,19,24,26	0
5	MG	B	3238	1/1	0.11	0.10	9,9,9,9	0
3	TPP	B	3236	26/26	0.12	-0.51	7,13,24,30	0
3	TPP	A	3236	26/26	0.12	-0.59	12,15,27,34	0
5	MG	A	3238	1/1	0.10	-0.63	12,12,12,12	0
6	CA	B	3239	1/1	0.08	-0.75	44,44,44,44	0
4	PYR	A	3237	6/6	0.13	-0.89	21,25,28,29	0
2	SF4	B	3233	8/8	0.07	-1.77	20,23,25,26	0
2	SF4	A	3235	8/8	0.06	-1.84	16,18,18,18	0
2	SF4	B	3235	8/8	0.06	-1.89	12,13,14,15	0
2	SF4	A	3234	8/8	0.06	-2.23	23,24,24,25	0
2	SF4	A	3233	8/8	0.06	-2.72	27,30,32,33	0
2	SF4	B	3234	8/8	0.05	-3.35	16,17,19,19	0

6.5 Other polymers ⓘ

There are no such residues in this entry.