



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 31, 2017 – 05:42 PM EDT

PDB ID : 3MM6
Title : Dissimilatory sulfite reductase cyanide complex
Authors : Parey, K.; Warkentin, E.; Kroneck, P.M.H.; Ermler, U.
Deposited on : unknown
Resolution : 1.90 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 : rb-20030345
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) : rb-20030345

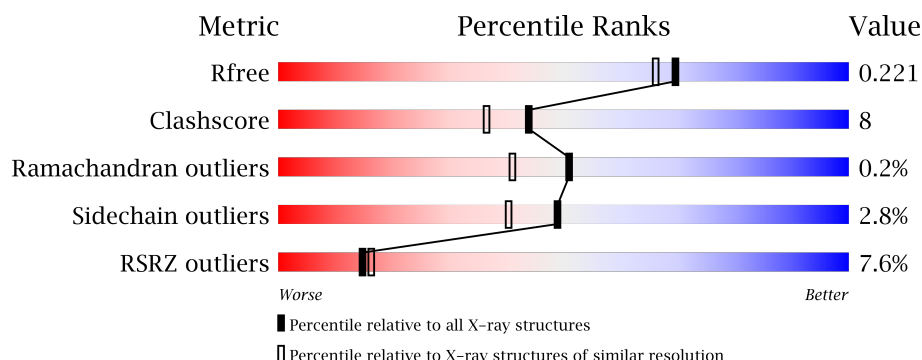
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	5047 (1.90-1.90)
Clashscore	112137	5731 (1.90-1.90)
Ramachandran outliers	110173	5669 (1.90-1.90)
Sidechain outliers	110143	5670 (1.90-1.90)
RSRZ outliers	101464	5100 (1.90-1.90)

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	418	<div> <div>4%</div> <div>86%</div> <div>14%</div> </div>
1	D	418	<div> <div>11%</div> <div>82%</div> <div>17%</div> <div>.</div> </div>
2	B	366	<div> <div>2%</div> <div>87%</div> <div>12%</div> <div>..</div> </div>
2	E	366	<div> <div>14%</div> <div>77%</div> <div>19%</div> <div>..</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria.

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	SF4	E	585	-	-	X	-
5	CYN	A	591	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 13513 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.

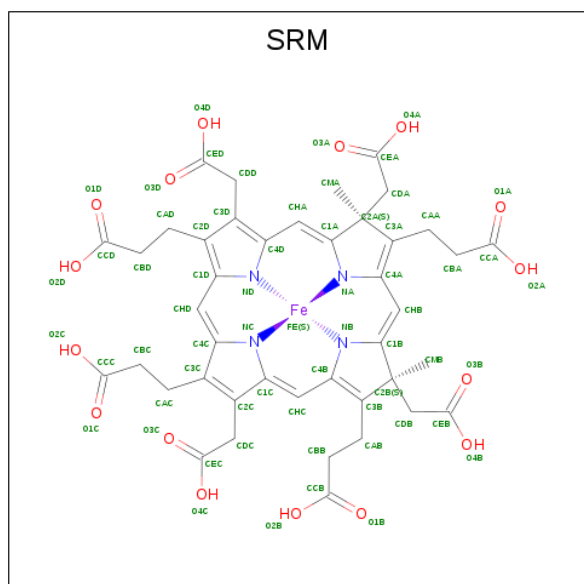
- Molecule 1 is a protein called Sulfite reductase, dissimilatory-type subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	417	Total	C	N	O	S	0	0	0
			3330	2134	557	613	26			
1	D	417	Total	C	N	O	S	0	0	0
			3330	2134	557	613	26			

- Molecule 2 is a protein called Sulfite reductase, dissimilatory-type subunit beta.

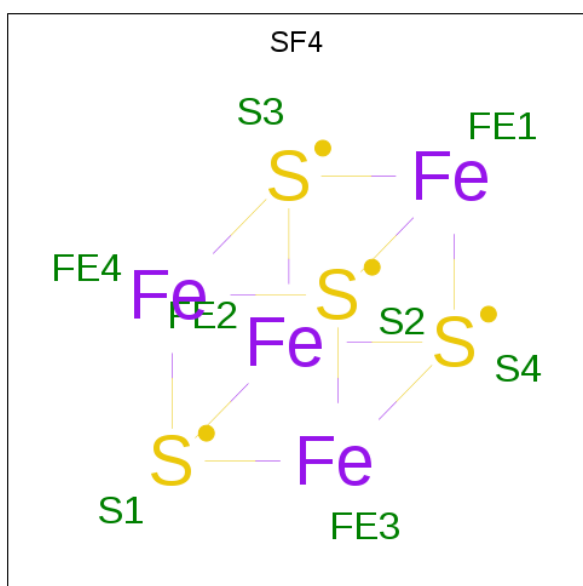
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	363	Total	C	N	O	S	0	0	0
			2901	1862	491	526	22			
2	E	363	Total	C	N	O	S	0	0	0
			2901	1862	491	526	22			

- Molecule 3 is SIROHEME (three-letter code: SRM) (formula: $C_{42}H_{42}FeN_4O_{16}$).



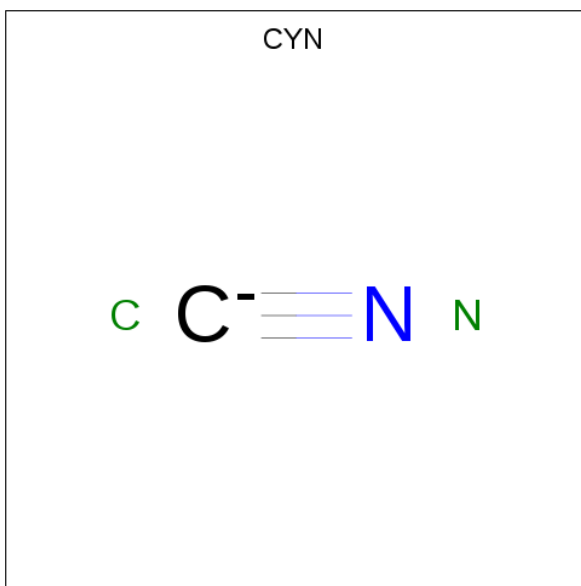
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	Fe	N	O	
			63	42	1	4	16	
3	B	1	Total	C	Fe	N	O	
			63	42	1	4	16	
3	D	1	Total	C	Fe	N	O	
			63	42	1	4	16	
3	E	1	Total	C	Fe	N	O	
			63	42	1	4	16	

- Molecule 4 is IRON/SULFUR CLUSTER (three-letter code: SF4) (formula: Fe₄S₄).



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

- Molecule 5 is CYANIDE ION (three-letter code: CYN) (formula: CN).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	N	0	0
			2	1	1		
5	A	1	Total	C	N	0	0
			2	1	1		
5	D	1	Total	C	N	0	0
			2	1	1		

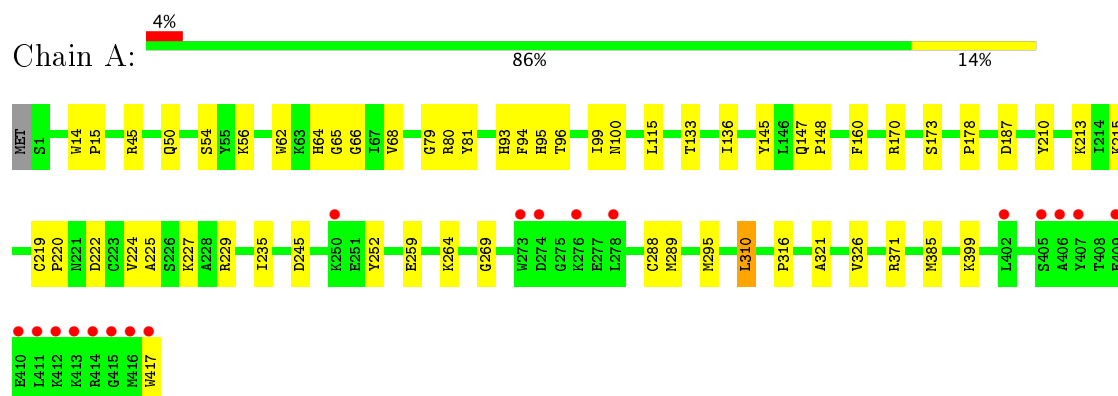
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	261	Total	O	0	0
			261	261		
6	B	299	Total	O	0	0
			299	299		
6	D	105	Total	O	0	0
			105	105		
6	E	64	Total	O	0	0
			64	64		

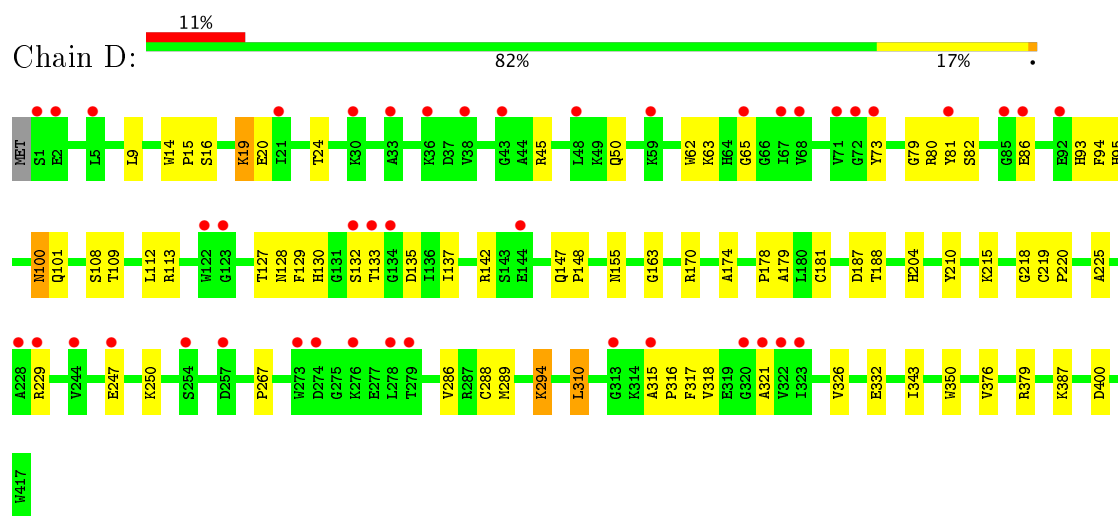
3 Residue-property plots

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

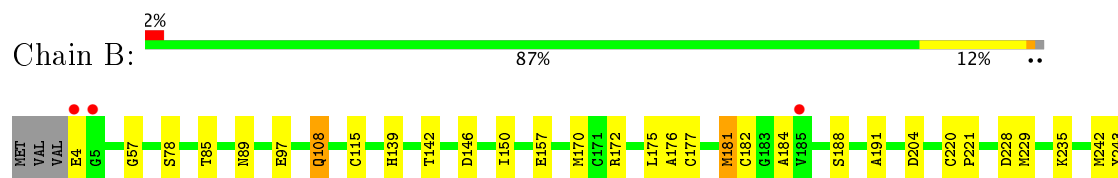
- Molecule 1: Sulfite reductase, dissimilatory-type subunit alpha



- Molecule 1: Sulfite reductase, dissimilatory-type subunit alpha

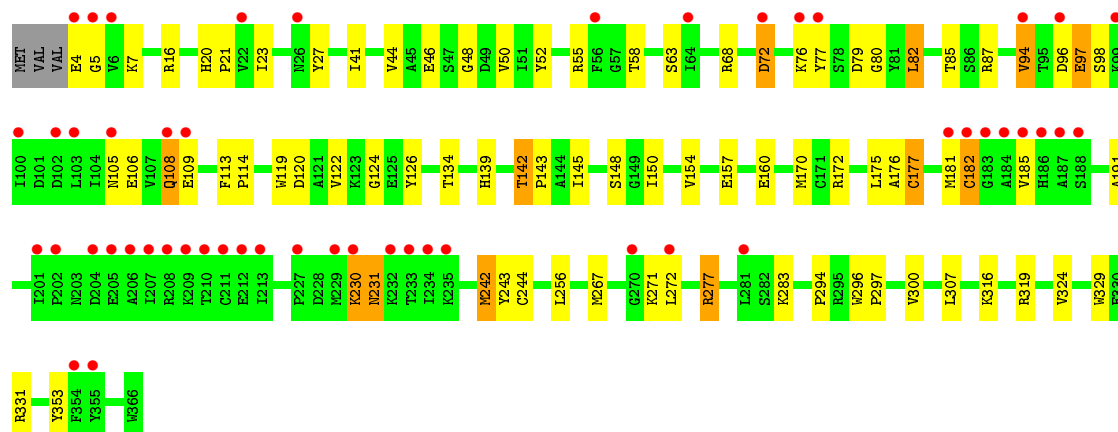
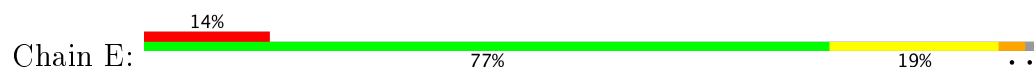


- Molecule 2: Sulfite reductase, dissimilatory-type subunit beta





- Molecule 2: Sulfite reductase, dissimilatory-type subunit beta



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	94.54Å 68.68Å 145.63Å 90.00° 107.57° 90.00°	Depositor
Resolution (Å)	20.00 – 1.90 48.39 – 1.90	Depositor EDS
% Data completeness (in resolution range)	95.4 (20.00-1.90) 95.4 (48.39-1.90)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.75 (at 1.90Å)	Xtriage
Refinement program	REFMAC 5.6.0046	Depositor
R, R_{free}	0.180 , 0.217 0.181 , 0.221	Depositor DCC
R_{free} test set	6694 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	27.8	Xtriage
Anisotropy	0.115	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 50.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.020 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	13513	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.04% 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: SF4, CYN, SRM

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	0.96	0/3417	0.86	3/4610 (0.1%)
1	D	0.71	5/3417 (0.1%)	0.73	2/4610 (0.0%)
2	B	1.08	3/2984 (0.1%)	0.91	3/4058 (0.1%)
2	E	0.80	9/2984 (0.3%)	0.77	5/4058 (0.1%)
All	All	0.89	17/12802 (0.1%)	0.82	13/17336 (0.1%)

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	72	ASP	CB-CG	14.12	1.81	1.51
2	E	68	ARG	CZ-NH1	9.98	1.46	1.33
1	D	20	GLU	C-O	8.98	1.40	1.23
2	E	106	GLU	CD-OE1	7.30	1.33	1.25
2	E	96	ASP	C-N	6.91	1.50	1.34
2	E	94	VAL	CB-CG1	6.59	1.66	1.52
2	E	106	GLU	CD-OE2	6.37	1.32	1.25
2	E	96	ASP	C-O	6.27	1.35	1.23
2	E	98	SER	CB-OG	6.03	1.50	1.42
1	D	20	GLU	CG-CD	5.94	1.60	1.51
1	D	113	ARG	CD-NE	5.80	1.56	1.46
1	D	181	CYS	CB-SG	-5.54	1.72	1.81
2	B	108	GLN	CB-CG	5.39	1.67	1.52
2	B	286	VAL	CB-CG1	5.38	1.64	1.52
1	D	20	GLU	C-N	5.28	1.46	1.34
2	E	72	ASP	CG-OD2	5.27	1.37	1.25
2	B	181	MET	C-O	5.15	1.33	1.23

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	72	ASP	CB-CG-OD2	9.31	126.68	118.30
1	D	113	ARG	NE-CZ-NH1	7.20	123.90	120.30
2	B	172	ARG	NE-CZ-NH2	-6.46	117.07	120.30
2	E	68	ARG	NE-CZ-NH1	-6.39	117.11	120.30
1	D	181	CYS	CA-CB-SG	-6.27	102.71	114.00
1	A	45	ARG	NE-CZ-NH2	-5.72	117.44	120.30
2	E	277	ARG	NE-CZ-NH1	5.68	123.14	120.30
1	A	45	ARG	NE-CZ-NH1	5.49	123.04	120.30
2	E	68	ARG	NE-CZ-NH2	5.45	123.03	120.30
2	B	204	ASP	CB-CG-OD1	5.35	123.12	118.30
2	E	277	ARG	NE-CZ-NH2	-5.26	117.67	120.30
2	B	258	ASP	CB-CG-OD2	-5.21	113.61	118.30
1	A	245	ASP	CB-CG-OD1	5.12	122.91	118.30

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	3330	0	3276	46	0
1	D	3330	0	3276	56	0
2	B	2901	0	2837	29	0
2	E	2901	0	2837	67	0
3	A	63	0	34	10	0
3	B	63	0	34	1	0
3	D	63	0	34	7	0
3	E	63	0	34	5	0
4	A	16	0	0	0	0
4	B	16	0	0	0	0
4	D	16	0	0	2	0
4	E	16	0	0	2	0
5	A	4	0	0	5	0
5	D	2	0	0	1	0
6	A	261	0	0	5	0
6	B	299	0	0	3	0
6	D	105	0	0	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	E	64	0	0	1	0
All	All	13513	0	12362	191	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 8.

All (191) 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)
2:E:72:ASP:CG	2:E:72:ASP:CB	1.81	1.49
1:A:81:TYR:CE2	1:A:93:HIS:CD2	2.42	1.08
1:A:81:TYR:HE2	1:A:93:HIS:CD2	1.72	1.08
2:E:85:THR:HB	3:E:570:SRM:HAB2	1.39	1.04
2:E:185:VAL:HG23	6:E:683:HOH:O	1.62	0.97
2:B:108:GLN:HE21	2:B:115:CYS:H	1.15	0.92
2:E:230:LYS:CB	2:E:231:ASN:HB2	1.99	0.92
1:A:81:TYR:CD2	1:A:93:HIS:HD2	1.89	0.89
1:A:371:ARG:NH1	6:A:771:HOH:O	2.05	0.88
2:E:230:LYS:HB2	2:E:231:ASN:HB2	1.55	0.86
1:D:170:ARG:HH12	5:D:592:CYN:C	1.90	0.84
4:D:576:SF4:S1	6:D:774:HOH:O	2.37	0.81
1:A:81:TYR:CE2	1:A:93:HIS:HD2	1.95	0.78
2:E:134:THR:HG21	2:E:182:CYS:HB3	1.65	0.77
2:E:134:THR:HG21	2:E:182:CYS:CB	2.15	0.76
1:A:66:GLY:H	1:A:81:TYR:HE1	1.32	0.73
1:D:94:PHE:O	2:E:139:HIS:HE1	1.71	0.72
3:A:580:SRM:C4C	2:B:182:CYS:HA	2.21	0.70
1:D:288:CYS:O	1:D:289:MET:HB2	1.92	0.69
1:A:94:PHE:O	2:B:139:HIS:HE1	1.76	0.68
1:A:133:THR:OG1	5:A:591:CYN:C	2.41	0.68
1:A:229:ARG:HD3	2:B:184:ALA:HB2	1.77	0.67
2:E:55:ARG:HH22	3:E:570:SRM:HBA2	1.59	0.66
1:A:93:HIS:ND1	6:A:652:HOH:O	2.29	0.65
2:B:85:THR:HB	3:B:570:SRM:HAB1	1.82	0.62
1:A:170:ARG:HH11	5:A:591:CYN:C	2.12	0.62
1:A:14:TRP:CD2	1:A:15:PRO:HD2	2.35	0.62
1:A:81:TYR:CD2	1:A:93:HIS:CD2	2.74	0.62
1:D:63:LYS:HE2	1:D:81:TYR:HB3	1.81	0.62
1:A:229:ARG:NH1	3:A:580:SRM:O4C	2.34	0.61
1:D:16:SER:HB3	1:D:19:LYS:HB3	1.82	0.61
1:D:400:ASP:OD1	6:D:579:HOH:O	2.16	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:128:ASN:HB2	1:D:137:ILE:HB	1.84	0.60
1:D:14:TRP:CD2	1:D:15:PRO:HD2	2.38	0.59
3:A:580:SRM:NC	2:B:182:CYS:HA	2.18	0.59
2:E:272:LEU:HB2	3:E:570:SRM:CCC	2.33	0.59
2:E:296:TRP:N	2:E:297:PRO:HD3	2.17	0.58
3:A:580:SRM:CMB	3:A:580:SRM:CBB	2.82	0.58
2:E:150:ILE:O	2:E:154:VAL:HG23	2.04	0.58
2:E:230:LYS:CA	2:E:231:ASN:HB2	2.33	0.58
2:E:72:ASP:CG	2:E:72:ASP:CA	2.69	0.58
1:D:112:LEU:HD11	2:E:82:LEU:HD11	1.86	0.58
2:E:170:MET:O	2:E:319:ARG:HG2	2.03	0.57
1:D:73:TYR:CD1	1:D:204:HIS:HB3	2.40	0.57
1:A:94:PHE:O	2:B:139:HIS:CE1	2.56	0.56
1:D:81:TYR:CE2	1:D:93:HIS:HD2	2.24	0.56
1:A:65:GLY:HA2	1:A:81:TYR:HD1	1.70	0.56
2:E:134:THR:HG21	2:E:182:CYS:HB2	1.87	0.56
2:E:142:THR:N	2:E:143:PRO:CD	2.70	0.55
1:D:62:TRP:HB3	1:D:80:ARG:HD2	1.88	0.55
3:A:580:SRM:C1C	2:B:182:CYS:HA	2.37	0.55
1:D:247:GLU:HA	1:D:250:LYS:HD2	1.88	0.55
3:E:570:SRM:HBD2	3:E:570:SRM:HDD2	1.88	0.54
1:A:219:CYS:HB2	1:A:220:PRO:CD	2.37	0.54
1:D:45:ARG:HG3	1:D:45:ARG:HH11	1.71	0.54
3:D:580:SRM:CMB	3:D:580:SRM:HBB2	2.38	0.54
3:D:580:SRM:HMB1	3:D:580:SRM:HBB2	1.89	0.53
2:E:267:MET:HA	2:E:283:LYS:O	2.08	0.53
2:B:108:GLN:NE2	2:B:115:CYS:H	1.95	0.53
2:B:299:LEU:C	2:B:299:LEU:HD23	2.29	0.53
3:A:580:SRM:O1A	2:B:139:HIS:HD2	1.91	0.53
2:E:182:CYS:HB2	4:E:585:SF4:S2	2.47	0.53
2:E:296:TRP:N	2:E:297:PRO:CD	2.72	0.53
1:D:215:LYS:HE2	1:D:229:ARG:HG3	1.90	0.53
1:A:81:TYR:HD2	1:A:93:HIS:HD2	1.48	0.52
1:D:94:PHE:O	2:E:139:HIS:CE1	2.59	0.52
2:B:170:MET:O	2:B:319:ARG:HG2	2.10	0.52
1:D:387:LYS:NZ	6:D:737:HOH:O	2.43	0.52
1:A:219:CYS:HB2	1:A:220:PRO:HD2	1.91	0.52
1:A:66:GLY:N	1:A:81:TYR:HE1	2.04	0.52
2:B:175:LEU:HD23	2:B:175:LEU:C	2.30	0.52
2:B:184:ALA:HB3	6:B:499:HOH:O	2.10	0.51
2:E:120:ASP:OD1	2:E:122:VAL:HB	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:225:ALA:HB3	1:D:229:ARG:HE	1.75	0.51
1:D:286:VAL:HG22	6:D:774:HOH:O	2.10	0.51
3:A:580:SRM:CMB	3:A:580:SRM:HBB2	2.41	0.51
1:D:315:ALA:HB1	1:D:316:PRO:CD	2.41	0.51
1:A:222:ASP:OD2	1:A:227:LYS:NZ	2.42	0.51
1:D:343:ILE:HD11	1:D:379:ARG:NH2	2.26	0.51
1:D:127:THR:HA	1:D:137:ILE:O	2.11	0.50
2:E:176:ALA:HB2	2:E:185:VAL:HG21	1.92	0.50
1:A:170:ARG:HH12	5:A:590:CYN:C	2.25	0.50
1:D:63:LYS:HE2	1:D:81:TYR:CB	2.42	0.50
1:A:79:GLY:HA2	1:A:95:HIS:ND1	2.27	0.50
2:E:271:LYS:NZ	2:E:277:ARG:O	2.41	0.50
2:E:105:ASN:O	2:E:109:GLU:HB2	2.12	0.49
1:D:14:TRP:CG	1:D:15:PRO:HD2	2.47	0.49
1:D:100:ASN:HD22	1:D:101:GLN:N	2.11	0.49
1:D:174:ALA:HB1	1:D:188:THR:HB	1.94	0.48
1:A:225:ALA:O	1:A:229:ARG:HG2	2.14	0.48
1:A:235:ILE:HD12	1:A:310:LEU:HD22	1.94	0.48
1:A:399:LYS:HB2	1:A:417:TRP:CZ2	2.48	0.48
1:A:147:GLN:HB3	1:A:148:PRO:HD3	1.96	0.48
1:D:187:ASP:OD2	2:E:16:ARG:NH2	2.46	0.48
3:D:580:SRM:O1A	2:E:139:HIS:HD2	1.96	0.48
1:D:155:ASN:OD1	2:E:7:LYS:HE3	2.14	0.48
2:B:157:GLU:HG3	2:B:300:VAL:HG11	1.95	0.48
2:E:126:TYR:CD1	2:E:126:TYR:N	2.81	0.48
1:A:213:LYS:HE2	5:A:591:CYN:C	2.44	0.48
3:A:580:SRM:HMB3	3:A:580:SRM:HBB2	1.94	0.48
1:A:178:PRO:HG3	1:A:187:ASP:HA	1.96	0.48
1:D:147:GLN:HB3	1:D:148:PRO:HD3	1.95	0.48
2:B:228:ASP:HB2	2:B:235:LYS:HG3	1.96	0.48
1:D:267:PRO:HG2	6:D:774:HOH:O	2.13	0.48
1:A:133:THR:HG21	5:A:591:CYN:C	2.44	0.47
3:D:580:SRM:HHB	3:D:580:SRM:HBA1	1.95	0.47
3:D:580:SRM:HDB1	2:E:134:THR:HG22	1.96	0.47
2:E:134:THR:HB	4:E:585:SF4:S4	2.54	0.47
1:D:9:LEU:HD21	2:E:294:PRO:O	2.15	0.47
1:A:56:LYS:NZ	6:A:526:HOH:O	2.42	0.47
1:D:79:GLY:HA2	1:D:95:HIS:ND1	2.30	0.46
1:D:294:LYS:HB3	1:D:294:LYS:HE2	1.61	0.46
1:D:316:PRO:HA	1:D:321:ALA:N	2.30	0.46
2:E:175:LEU:C	2:E:175:LEU:HD23	2.35	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:315:ALA:HB1	1:D:316:PRO:HD2	1.98	0.46
1:D:179:ALA:O	2:E:23:ILE:HG23	2.16	0.46
1:D:310:LEU:HD12	1:D:326:VAL:HA	1.97	0.45
1:D:45:ARG:HG3	1:D:45:ARG:NH1	2.30	0.45
2:E:52:TYR:CE1	2:E:97:GLU:HB3	2.51	0.45
1:D:187:ASP:OD1	2:E:27:TYR:OH	2.34	0.45
2:E:124:GLY:HA3	2:E:316:LYS:HD3	1.99	0.45
2:E:185:VAL:HG13	2:E:191:ALA:HB1	1.97	0.45
2:B:146:ASP:O	2:B:150:ILE:HD12	2.17	0.45
2:B:353:TYR:HA	2:E:353:TYR:HA	1.97	0.45
1:D:129:PHE:HA	1:D:130:HIS:HA	1.66	0.45
1:D:350:TRP:CG	1:D:376:VAL:HG21	2.51	0.45
1:D:81:TYR:CE2	1:D:93:HIS:CD2	3.03	0.45
1:A:99:ILE:HB	1:A:136:ILE:HB	1.99	0.45
2:E:76:LYS:HD3	2:E:77:TYR:CZ	2.52	0.45
2:B:324:VAL:HG11	2:B:329:TRP:CE2	2.52	0.45
2:E:44:VAL:HG22	2:E:50:VAL:HG22	1.99	0.45
2:E:157:GLU:HG3	2:E:300:VAL:CG1	2.47	0.44
1:A:173:SER:OG	1:A:215:LYS:HG2	2.18	0.44
2:E:58:THR:HG22	2:E:113:PHE:CE2	2.53	0.44
1:A:145:TYR:C	1:A:148:PRO:HD2	2.37	0.44
2:B:176:ALA:HB1	2:B:181:MET:HA	1.99	0.44
1:D:219:CYS:HB2	1:D:220:PRO:CD	2.48	0.44
1:D:135:ASP:OD2	1:D:163:GLY:HA3	2.18	0.44
3:E:570:SRM:HAC2	3:E:570:SRM:HCD1	1.85	0.43
2:E:87:ARG:CZ	2:E:172:ARG:HD2	2.47	0.43
3:A:580:SRM:HDD2	3:A:580:SRM:HAD1	1.79	0.43
2:E:20:HIS:HA	2:E:21:PRO:HD2	1.64	0.43
1:A:96:THR:HG21	6:A:553:HOH:O	2.17	0.43
1:D:24:THR:OG1	2:E:63:SER:HB2	2.19	0.43
1:A:288:CYS:O	1:A:289:MET:HB2	2.18	0.43
1:D:112:LEU:CD1	2:E:82:LEU:HD11	2.48	0.43
3:D:580:SRM:CBB	3:D:580:SRM:CMB	2.96	0.43
1:A:316:PRO:HA	1:A:321:ALA:N	2.34	0.43
1:A:96:THR:HG23	2:B:139:HIS:NE2	2.33	0.43
2:E:58:THR:HG22	2:E:113:PHE:CD2	2.53	0.43
2:E:113:PHE:HA	2:E:114:PRO:HD2	1.61	0.43
2:B:359:ARG:HD3	2:B:363:GLN:O	2.19	0.43
1:A:269:GLY:HA2	6:A:589:HOH:O	2.19	0.43
2:B:191:ALA:HB3	2:B:267:MET:HB2	2.01	0.42
2:E:324:VAL:HG11	2:E:329:TRP:CE2	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:580:SRM:C3C	2:B:182:CYS:HA	2.49	0.42
1:D:86:GLU:H	1:D:86:GLU:HG2	1.69	0.42
1:D:65:GLY:HA2	1:D:81:TYR:CD1	2.53	0.42
2:E:20:HIS:HB3	2:E:23:ILE:HG13	2.02	0.42
2:E:52:TYR:HB2	2:E:94:VAL:O	2.18	0.42
1:D:50:GLN:HE22	1:D:82:SER:HB2	1.85	0.42
2:E:242:MET:SD	2:E:244:CYS:HB3	2.59	0.42
2:B:313:ASN:HB3	6:B:478:HOH:O	2.19	0.42
2:E:230:LYS:HE3	2:E:230:LYS:HA	2.02	0.42
1:D:178:PRO:HG3	1:D:187:ASP:HA	2.01	0.41
1:A:259:GLU:OE2	1:A:264:LYS:HE3	2.19	0.41
1:D:316:PRO:HG2	2:E:181:MET:HE3	2.01	0.41
1:D:317:PHE:HA	1:D:318:VAL:HA	1.80	0.41
1:D:218:GLY:HA3	4:D:575:SF4:S2	2.60	0.41
3:D:580:SRM:HDD2	3:D:580:SRM:HAD1	1.91	0.41
2:E:175:LEU:HD23	2:E:176:ALA:N	2.35	0.41
2:E:145:ILE:HG12	2:E:177:CYS:HA	2.01	0.41
2:E:20:HIS:CD2	2:E:79:ASP:HB2	2.55	0.41
1:A:115:LEU:HD21	1:A:160:PHE:CG	2.54	0.41
2:B:57:GLY:HA2	2:B:89:ASN:OD1	2.21	0.41
2:E:46:GLU:C	2:E:48:GLY:H	2.22	0.41
1:D:109:THR:N	2:E:80:GLY:O	2.52	0.41
1:A:219:CYS:CB	1:A:220:PRO:CD	2.98	0.41
1:A:62:TRP:HB3	1:A:80:ARG:HD2	2.03	0.41
1:D:108:SER:HA	2:E:80:GLY:O	2.21	0.40
1:D:93:HIS:HB2	1:D:142:ARG:HG2	2.03	0.40
1:A:252:TYR:CD2	1:A:295:MET:HG2	2.56	0.40
2:B:220:CYS:HA	2:B:221:PRO:HD3	1.76	0.40
2:B:330:GLU:HB2	6:B:369:HOH:O	2.21	0.40
2:E:134:THR:CG2	2:E:182:CYS:HB3	2.43	0.40
2:E:331:ARG:HD3	2:E:331:ARG:HA	1.94	0.40
1:A:326:VAL:HB	1:A:385:MET:HA	2.03	0.40
1:A:64:HIS:HE1	2:B:249:THR:O	2.05	0.40
2:E:108:GLN:N	2:E:108:GLN:HE21	2.20	0.40
2:E:119:TRP:CE2	2:E:170:MET:CE	3.04	0.40
2:E:4:GLU:HG3	2:E:5:GLY:H	1.87	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	415/418 (99%)	405 (98%)	10 (2%)	0	100	100
1	D	415/418 (99%)	391 (94%)	24 (6%)	0	100	100
2	B	361/366 (99%)	350 (97%)	10 (3%)	1 (0%)	44	34
2	E	361/366 (99%)	333 (92%)	26 (7%)	2 (1%)	28	16
All	All	1552/1568 (99%)	1479 (95%)	70 (4%)	3 (0%)	51	41

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	E	231	ASN
2	B	142	THR
2	E	142	THR

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	353/354 (100%)	346 (98%)	7 (2%)	60	55
1	D	353/354 (100%)	345 (98%)	8 (2%)	56	49
2	B	314/317 (99%)	305 (97%)	9 (3%)	48	39
2	E	314/317 (99%)	301 (96%)	13 (4%)	35	24
All	All	1334/1342 (99%)	1297 (97%)	37 (3%)	49	40

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	50	GLN
1	A	54	SER
1	A	68	VAL
1	A	100	ASN
1	A	210	TYR
1	A	224	VAL
1	A	310	LEU
2	B	4	GLU
2	B	78	SER
2	B	97	GLU
2	B	177	CYS
2	B	188	SER
2	B	229	MET
2	B	242	MET
2	B	243	TYR
2	B	283	LYS
1	D	19	LYS
1	D	100	ASN
1	D	132	SER
1	D	133	THR
1	D	210	TYR
1	D	294	LYS
1	D	310	LEU
1	D	332	GLU
2	E	41	ILE
2	E	82	LEU
2	E	97	GLU
2	E	108	GLN
2	E	148	SER
2	E	160	GLU
2	E	177	CYS
2	E	182	CYS
2	E	230	LYS
2	E	242	MET
2	E	243	TYR
2	E	256	LEU
2	E	307	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	50	GLN
1	A	93	HIS

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Mol	Chain	Res	Type
1	A	100	ASN
2	B	108	GLN
2	B	139	HIS
1	D	50	GLN
1	D	93	HIS
1	D	100	ASN
2	E	108	GLN
2	E	139	HIS
2	E	180	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

15 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	$\# Z > 2$
4	SF4	A	575	1	0,12,12	0.00	-	0,24,24	0.00	-
4	SF4	A	576	1	0,12,12	0.00	-	0,24,24	0.00	-
3	SRM	A	580	2,5	34,70,70	1.67	8 (23%)	34,112,112	4.13	13 (38%)
5	CYN	A	590	3	0,1,1	0.00	-	0,0,0	0.00	-
5	CYN	A	591	-	0,1,1	0.00	-	0,0,0	0.00	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SRM	B	570	1	34,70,70	2.30	7 (20%)	34,112,112	3.92	17 (50%)
4	SF4	B	585	2	0,12,12	0.00	-	0,24,24	0.00	-
4	SF4	B	586	2	0,12,12	0.00	-	0,24,24	0.00	-
4	SF4	D	575	1	0,12,12	0.00	-	0,24,24	0.00	-
4	SF4	D	576	1,6	0,12,12	0.00	-	0,24,24	0.00	-
3	SRM	D	580	2,5	34,70,70	2.06	8 (23%)	34,112,112	3.67	14 (41%)
5	CYN	D	592	3	0,1,1	0.00	-	0,0,0	0.00	-
3	SRM	E	570	1	34,70,70	2.41	12 (35%)	34,112,112	3.89	18 (52%)
4	SF4	E	585	2	0,12,12	0.00	-	0,24,24	0.00	-
4	SF4	E	586	2	0,12,12	0.00	-	0,24,24	0.00	-

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
4	SF4	A	575	1	-	0/0/48/48	0/6/5/5
4	SF4	A	576	1	-	0/0/48/48	0/6/5/5
3	SRM	A	580	2,5	-	0/22/126/126	0/0/8/8
5	CYN	A	590	3	-	0/0/0/0	0/0/0/0
5	CYN	A	591	-	-	0/0/0/0	0/0/0/0
3	SRM	B	570	1	-	0/22/126/126	0/0/8/8
4	SF4	B	585	2	-	0/0/48/48	0/6/5/5
4	SF4	B	586	2	-	0/0/48/48	0/6/5/5
4	SF4	D	575	1	-	0/0/48/48	0/6/5/5
4	SF4	D	576	1,6	-	0/0/48/48	0/6/5/5
3	SRM	D	580	2,5	-	0/22/126/126	0/0/8/8
5	CYN	D	592	3	-	0/0/0/0	0/0/0/0
3	SRM	E	570	1	-	0/22/126/126	0/0/8/8
4	SF4	E	585	2	-	0/0/48/48	0/6/5/5
4	SF4	E	586	2	-	0/0/48/48	0/6/5/5

All (35) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	570	SRM	C4C-NC	-6.89	1.28	1.36
3	E	570	SRM	C4A-NA	-5.54	1.28	1.39
3	E	570	SRM	C1C-NC	-5.38	1.30	1.36
3	B	570	SRM	C4A-NA	-5.17	1.29	1.39
3	E	570	SRM	C4C-NC	-4.27	1.31	1.36
3	E	570	SRM	C1B-NB	-3.64	1.31	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	580	SRM	C4A-NA	-3.55	1.32	1.39
3	B	570	SRM	C1B-NB	-3.46	1.32	1.38
3	D	580	SRM	C4C-NC	-2.51	1.33	1.36
3	E	570	SRM	CHC-C4B	-2.32	1.34	1.39
3	E	570	SRM	C1A-NA	-2.13	1.34	1.38
3	E	570	SRM	C4B-NB	-2.09	1.35	1.39
3	D	580	SRM	CHA-C1A	2.24	1.39	1.36
3	A	580	SRM	CAD-C2D	2.28	1.55	1.52
3	A	580	SRM	C1D-ND	2.29	1.41	1.36
3	E	570	SRM	CAD-C2D	2.35	1.56	1.52
3	A	580	SRM	C1C-NC	2.44	1.39	1.36
3	E	570	SRM	CAA-C3A	2.54	1.55	1.51
3	B	570	SRM	CHA-C1A	2.60	1.40	1.36
3	A	580	SRM	FE-NA	2.66	2.06	1.95
3	A	580	SRM	FE-NB	2.72	2.06	1.95
3	E	570	SRM	CHA-C1A	2.79	1.40	1.36
3	D	580	SRM	CAD-C2D	2.97	1.57	1.52
3	A	580	SRM	CHA-C1A	3.10	1.40	1.36
3	D	580	SRM	FE-NA	3.57	2.09	1.95
3	B	570	SRM	C3C-C2C	3.59	1.48	1.37
3	D	580	SRM	FE-NB	3.74	2.10	1.95
3	A	580	SRM	C3C-C2C	3.82	1.49	1.37
3	D	580	SRM	C3C-C2C	4.23	1.50	1.37
3	B	570	SRM	CAA-C3A	4.44	1.58	1.51
3	A	580	SRM	C3D-C2D	4.46	1.50	1.39
3	E	570	SRM	C3C-C2C	4.50	1.51	1.37
3	B	570	SRM	C3D-C2D	5.51	1.52	1.39
3	E	570	SRM	C3D-C2D	5.88	1.53	1.39
3	D	580	SRM	C3D-C2D	6.84	1.55	1.39

All (62) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	580	SRM	CAB-C3B-C2B	-18.06	103.10	123.52
3	D	580	SRM	CAB-C3B-C2B	-14.08	107.61	123.52
3	E	570	SRM	CAA-C3A-C2A	-10.70	111.42	123.52
3	B	570	SRM	CAA-C3A-C2A	-9.89	112.34	123.52
3	B	570	SRM	CAB-C3B-C2B	-9.31	113.00	123.52
3	B	570	SRM	CBC-CAC-C3C	-8.22	96.79	112.47
3	B	570	SRM	C4A-NA-C1A	-6.86	101.75	106.37
3	A	580	SRM	CAA-C3A-C2A	-6.00	116.73	123.52
3	E	570	SRM	CBC-CAC-C3C	-5.54	101.89	112.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	580	SRM	CAA-C3A-C2A	-5.40	117.42	123.52
3	E	570	SRM	C4D-CHA-C1A	-5.39	119.45	130.12
3	E	570	SRM	CEC-CDC-C2C	-5.31	105.81	116.14
3	B	570	SRM	C3B-C2B-C1B	-5.30	91.86	101.25
3	A	580	SRM	CDD-C3D-C2D	-4.62	118.16	126.48
3	E	570	SRM	CAC-C3C-C2C	-4.47	118.26	129.38
3	B	570	SRM	C4D-CHA-C1A	-4.45	121.31	130.12
3	B	570	SRM	CEC-CDC-C2C	-4.35	107.68	116.14
3	E	570	SRM	C4A-NA-C1A	-4.34	103.44	106.37
3	B	570	SRM	C3B-C4B-NB	-4.34	105.17	110.12
3	A	580	SRM	CAC-CBC-CCC	-4.23	105.43	112.66
3	E	570	SRM	C3B-C2B-C1B	-4.17	93.86	101.25
3	E	570	SRM	CHB-C4A-C3A	-4.12	116.12	125.47
3	E	570	SRM	CDC-C2C-C3C	-4.09	119.11	126.48
3	A	580	SRM	CAD-C2D-C3D	-3.97	119.49	129.38
3	D	580	SRM	CAD-C2D-C3D	-3.74	120.08	129.38
3	D	580	SRM	CDD-C3D-C2D	-3.67	119.86	126.48
3	D	580	SRM	C4D-CHA-C1A	-3.66	122.86	130.12
3	A	580	SRM	CAD-CBD-CCD	-3.62	106.48	112.66
3	D	580	SRM	CDD-C3D-C4D	-3.58	121.93	127.36
3	D	580	SRM	C3A-C2A-C1A	-3.54	94.98	101.25
3	E	570	SRM	CDD-C3D-C4D	-3.53	122.00	127.36
3	A	580	SRM	C4D-CHA-C1A	-3.07	124.04	130.12
3	E	570	SRM	CDC-C2C-C1C	-3.03	122.81	127.39
3	A	580	SRM	C3B-C2B-C1B	-2.56	96.72	101.25
3	B	570	SRM	CAB-CBB-CCB	-2.50	108.39	112.66
3	D	580	SRM	C3A-C4A-NA	-2.46	107.30	110.12
3	B	570	SRM	C2A-CDA-CEA	-2.29	112.00	115.39
3	A	580	SRM	C3A-C2A-C1A	-2.27	97.22	101.25
3	B	570	SRM	CDC-C2C-C3C	-2.27	122.39	126.48
3	B	570	SRM	CHB-C4A-C3A	-2.21	120.45	125.47
3	D	580	SRM	C3B-C2B-C1B	-2.20	97.35	101.25
3	B	570	SRM	CAD-CBD-CCD	-2.09	109.09	112.66
3	D	580	SRM	C4B-NB-C1B	-2.08	104.97	106.37
3	E	570	SRM	CHA-C1A-NA	2.02	127.36	124.18
3	E	570	SRM	C2A-CDA-CEA	2.31	118.80	115.39
3	D	580	SRM	CED-CDD-C3D	2.49	120.99	116.14
3	A	580	SRM	CDD-C3D-C4D	2.54	131.21	127.36
3	B	570	SRM	CHC-C4B-NB	2.94	129.32	123.80
3	D	580	SRM	C3B-C4B-NB	3.02	113.57	110.12
3	B	570	SRM	C2B-CDB-CEB	3.10	119.97	115.39
3	E	570	SRM	CED-CDD-C3D	3.16	122.29	116.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	580	SRM	CBD-CAD-C2D	3.54	119.25	112.48
3	E	570	SRM	CAA-CBA-CCA	3.72	119.01	112.66
3	B	570	SRM	CBD-CAD-C2D	4.65	121.37	112.48
3	A	580	SRM	C4A-NA-C1A	4.95	109.70	106.37
3	E	570	SRM	CAB-CBB-CCB	5.88	122.71	112.66
3	B	570	SRM	C3A-C4A-NA	6.67	117.74	110.12
3	D	580	SRM	C4A-NA-C1A	7.10	111.16	106.37
3	E	570	SRM	CBD-CAD-C2D	7.39	126.61	112.48
3	D	580	SRM	CBD-CAD-C2D	7.59	126.99	112.48
3	E	570	SRM	C3A-C4A-NA	7.77	119.00	110.12
3	A	580	SRM	C2A-CDA-CEA	8.22	127.55	115.39

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

10 monomers are involved in 33 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	580	SRM	10	0
5	A	590	CYN	1	0
5	A	591	CYN	4	0
3	B	570	SRM	1	0
4	D	575	SF4	1	0
4	D	576	SF4	1	0
3	D	580	SRM	7	0
5	D	592	CYN	1	0
3	E	570	SRM	5	0
4	E	585	SF4	2	0

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	417/418 (99%)	-0.02	18 (4%) 36 39	18, 31, 64, 112	0
1	D	417/418 (99%)	0.61	44 (10%) 7 8	22, 66, 111, 138	0
2	B	363/366 (99%)	-0.12	6 (1%) 70 73	18, 28, 43, 114	0
2	E	363/366 (99%)	0.75	51 (14%) 3 3	28, 69, 106, 148	0
All	All	1560/1568 (99%)	0.30	119 (7%) 15 16	18, 45, 99, 148	0

All (119) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	E	4	GLU	9.5
1	D	1	SER	7.8
1	A	417	TRP	7.8
2	B	4	GLU	6.3
2	E	232	LYS	5.9
2	E	229	MET	5.5
2	E	185	VAL	5.2
2	E	186	HIS	4.8
2	E	184	ALA	4.7
2	E	281	LEU	4.5
2	B	5	GLY	4.4
2	E	207	ILE	4.3
1	A	415	GLY	4.2
1	D	43	GLY	4.2
2	E	205	GLU	4.2
1	D	33	ALA	4.2
1	A	406	ALA	4.1
2	E	209	LYS	4.1
2	E	230	LYS	3.9
1	A	276	LYS	3.9
1	A	413	LYS	3.9

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Mol	Chain	Res	Type	RSRZ
1	D	228	ALA	3.7
1	D	81	TYR	3.7
1	D	68	VAL	3.7
1	D	229	ARG	3.7
2	E	201	ILE	3.7
2	E	77	TYR	3.6
1	D	67	ILE	3.6
1	D	276	LYS	3.5
1	D	85	GLY	3.4
1	D	5	LEU	3.4
1	D	30	LYS	3.3
1	A	412	LYS	3.2
1	D	36	LYS	3.2
2	E	188	SER	3.2
1	A	416	MET	3.2
2	E	6	VAL	3.2
1	D	72	GLY	3.1
2	E	210	THR	3.0
1	A	278	LEU	3.0
1	D	323	ILE	3.0
1	A	414	ARG	3.0
2	E	234	ILE	3.0
2	B	185	VAL	3.0
2	E	187	ALA	2.9
1	A	411	LEU	2.9
2	E	355	TYR	2.9
2	E	100	ILE	2.9
2	E	213	ILE	2.9
2	E	181	MET	2.8
2	E	211	CYS	2.8
1	D	38	VAL	2.8
1	D	247	GLU	2.8
1	A	407	TYR	2.7
2	E	206	ALA	2.7
2	E	233	THR	2.7
2	E	208	ARG	2.7
1	D	2	GLU	2.7
1	D	86	GLU	2.7
1	D	134	GLY	2.7
2	E	202	PRO	2.6
1	D	321	ALA	2.6
2	E	182	CYS	2.6

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Mol	Chain	Res	Type	RSRZ
2	E	204	ASP	2.6
2	E	183	GLY	2.6
1	D	273	TRP	2.6
1	D	71	VAL	2.6
1	A	273	TRP	2.5
1	D	59	LYS	2.5
1	D	132	SER	2.5
1	D	320	GLY	2.5
2	E	72	ASP	2.5
2	E	212	GLU	2.5
2	E	102	ASP	2.5
2	E	96	ASP	2.5
1	A	405	SER	2.5
1	D	279	THR	2.4
2	E	235	LYS	2.4
1	A	402	LEU	2.3
1	D	133	THR	2.3
2	E	5	GLY	2.3
2	B	272	LEU	2.3
2	B	355	TYR	2.3
2	E	108	GLN	2.3
1	D	122	TRP	2.3
1	A	410	GLU	2.3
1	D	65	GLY	2.3
1	D	123	GLY	2.3
1	A	274	ASP	2.3
1	D	244	VAL	2.2
1	D	254	SER	2.2
2	E	109	GLU	2.2
1	D	274	ASP	2.2
2	E	64	ILE	2.2
2	E	103	LEU	2.2
2	E	105	ASN	2.2
1	D	92	GLU	2.2
1	D	21	ILE	2.2
2	E	26	ASN	2.2
2	B	353	TYR	2.2
2	E	94	VAL	2.2
1	D	315	ALA	2.2
2	E	22	VAL	2.2
1	D	144	GLU	2.1
1	D	278	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
2	E	272	LEU	2.1
2	E	270	GLY	2.1
1	A	409	GLU	2.1
2	E	76	LYS	2.1
2	E	56	PHE	2.1
1	D	48	LEU	2.1
2	E	227	PRO	2.1
1	A	250	LYS	2.1
1	D	73	TYR	2.0
1	D	322	VAL	2.0
2	E	99	LYS	2.0
1	D	313	GLY	2.0
2	E	354	PHE	2.0
1	D	257	ASP	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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
4	SF4	B	586	8/8	0.99	0.09	-0.02	20,23,24,24	0
3	SRM	A	580	63/63	0.97	0.12	-0.10	16,24,37,49	0
3	SRM	B	570	63/63	0.98	0.11	-0.33	14,20,27,30	0
3	SRM	E	570	63/63	0.94	0.12	-0.38	33,45,56,63	0
4	SF4	A	575	8/8	0.99	0.10	-0.39	22,22,24,28	0
3	SRM	D	580	63/63	0.94	0.12	-0.61	40,59,84,98	0
5	CYN	A	591	2/2	0.98	0.08	-0.83	22,22,22,23	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	SF4	B	585	8/8	0.99	0.09	-1.08	18,22,23,25	0
4	SF4	E	585	8/8	0.92	0.05	-1.36	41,44,50,54	0
4	SF4	A	576	8/8	0.98	0.05	-1.38	26,28,30,32	0
4	SF4	E	586	8/8	0.98	0.03	-1.63	44,46,49,50	0
4	SF4	D	575	8/8	0.96	0.07	-1.64	38,42,46,46	0
4	SF4	D	576	8/8	0.98	0.04	-1.88	38,42,44,49	0
5	CYN	D	592	2/2	0.93	0.29	-	41,41,41,52	0
5	CYN	A	590	2/2	0.97	0.22	-	18,18,18,27	0

6.5 Other polymers [i](#)

There are no such residues in this entry.