



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:39 AM GMT

PDB ID : 3MM5
Title : Dissimilatory sulfite reductase in complex with the substrate sulfite
Authors : Parey, K.; Warkentin, E.; Kroneck, P.M.H.; Ermler, U.
Deposited on : 2010-04-19
Resolution : 1.80 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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

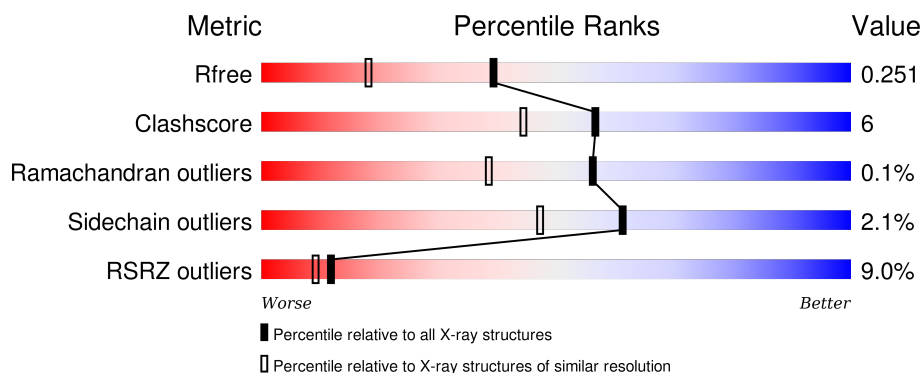
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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}	91344	4533 (1.80-1.80)
Clashscore	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)
RSRZ outliers	91569	4547 (1.80-1.80)

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>2%</div> <div>88%</div> <div>11%</div> </div>
1	D	418	<div> <div>16%</div> <div>87%</div> <div>12%</div> <div>.</div> </div>
2	B	366	<div> <div>%</div> <div>90%</div> <div>8%</div> <div>..</div> </div>
2	E	366	<div> <div>16%</div> <div>84%</div> <div>13%</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:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	SRM	D	580	-	-	-	X
3	SRM	E	570	-	-	-	X
4	SF4	D	575	-	-	-	X
4	SF4	D	576	-	-	-	X
4	SF4	E	585	-	-	X	X
4	SF4	E	586	-	-	-	X
5	SO3	D	590	-	-	-	X

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 13810 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.

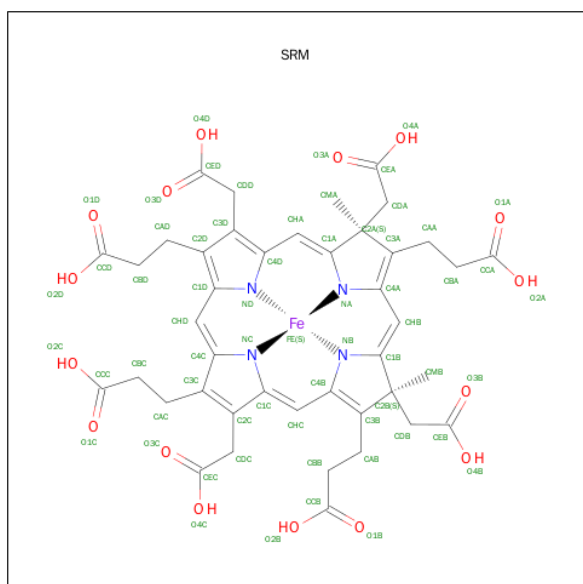
- Molecule 1 is a protein called Sulfite reductase, dissimilatory-type subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	417	Total	C	N	O	S	0	3	0
			3348	2146	557	619	26			
1	D	417	Total	C	N	O	S	0	1	0
			3336	2139	558	613	26			

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

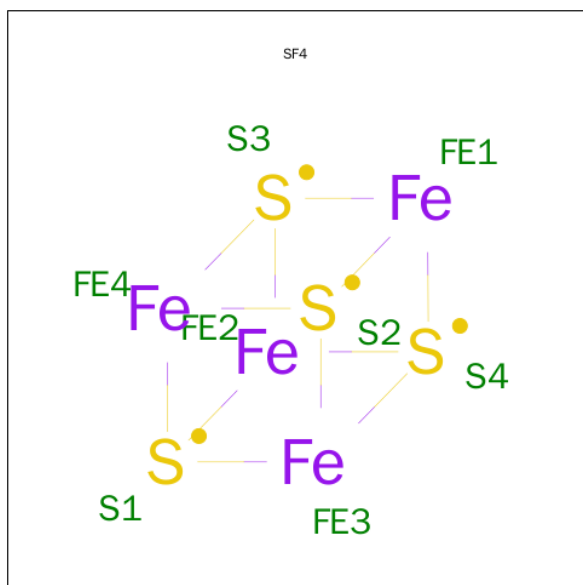
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	363	Total	C	N	O	S	5	5	0
			2930	1883	492	533	22			
2	E	363	Total	C	N	O	S	0	0	0
			2901	1862	491	526	22			

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



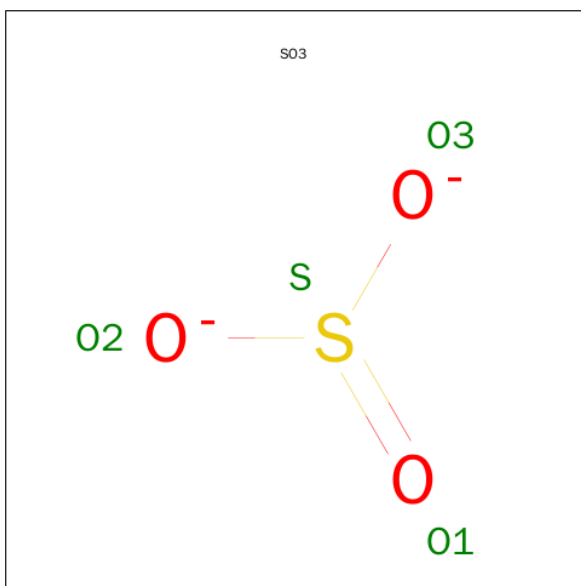
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	Fe	N	O	
			63	42	1	4	16	
3	B	1	Total	C	Fe	N	O	
			63	42	1	4	16	
3	D	1	Total	C	Fe	N	O	
			63	42	1	4	16	
3	E	1	Total	C	Fe	N	O	
			63	42	1	4	16	

- Molecule 4 is IRON/SULFUR CLUSTER (three-letter code: SF4) (formula: Fe_4S_4).



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 SULFITE ION (three-letter code: SO3) (formula: O₃S).



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

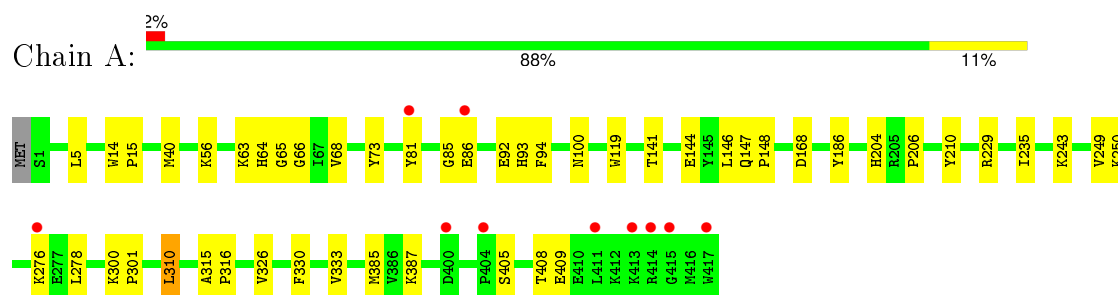
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	229	Total	O	0	0
			229	229		
6	B	565	Total	O	0	0
			565	565		
6	D	92	Total	O	0	0
			92	92		
6	E	85	Total	O	0	0
			85	85		

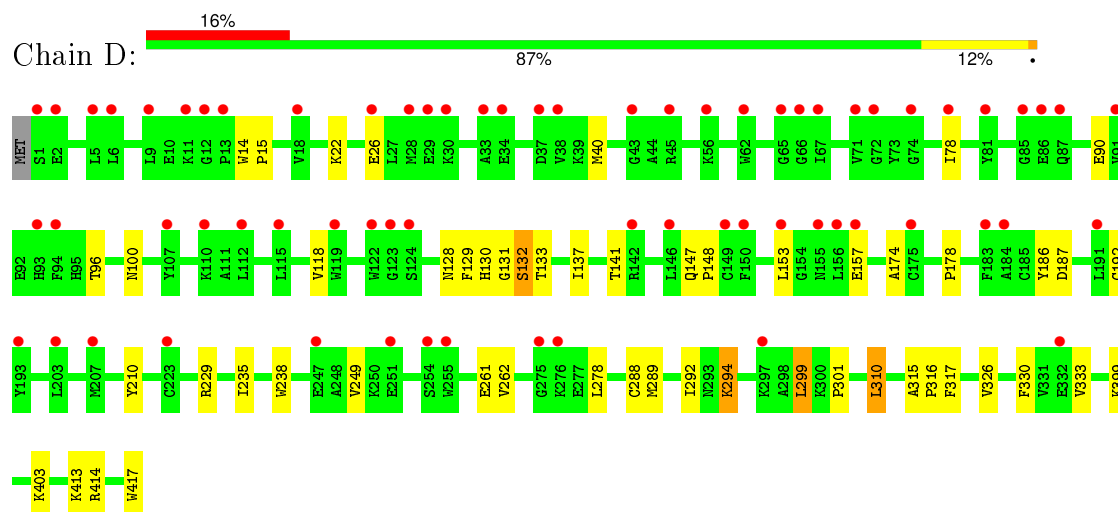
3 Residue-property plots

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.

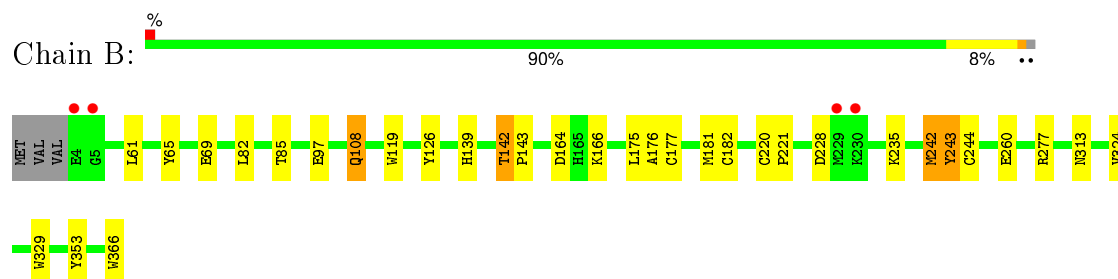
- Molecule 1: Sulfite reductase, dissimilatory-type subunit alpha



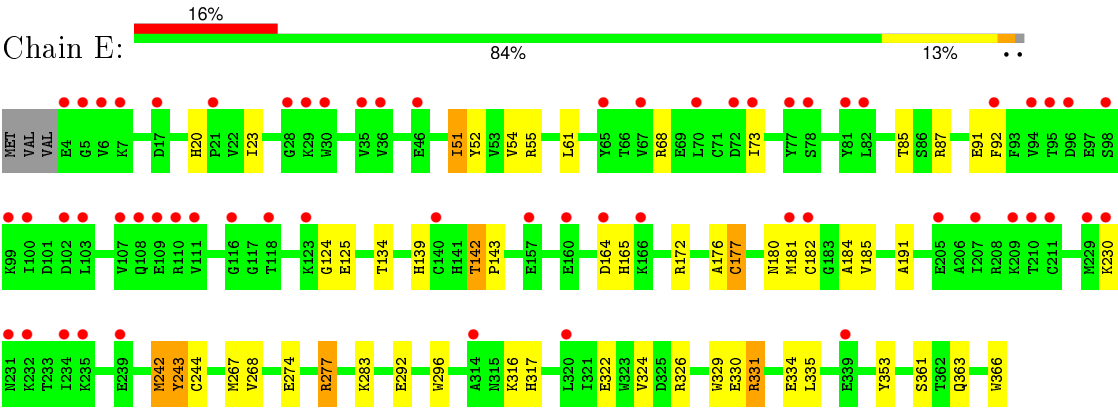
- Molecule 1: Sulfite reductase, dissimilatory-type subunit alpha



- 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	94.60Å 68.90Å 145.10Å 90.00° 107.50° 90.00°	Depositor
Resolution (Å)	15.00 – 1.80 15.00 – 1.80	Depositor EDS
% Data completeness (in resolution range)	100.0 (15.00-1.80) 99.3 (15.00-1.80)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.95 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.6.0046	Depositor
R, R_{free}	0.158 , 0.188 0.230 , 0.251	Depositor DCC
R_{free} test set	8181 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	19.4	Xtriage
Anisotropy	0.158	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 38.3	EDS
Estimated twinning fraction	0.016 for h,-k,-h-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 163613 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	13810	wwPDB-VP
Average B, all atoms (Å ²)	11.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.98% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SF4, SO3, 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	0.79	0/3444	0.75	0/4646
1	D	0.47	0/3426	0.56	0/4622
2	B	0.81	1/3025 (0.0%)	0.79	1/4114 (0.0%)
2	E	0.47	1/2984 (0.0%)	0.60	3/4058 (0.1%)
All	All	0.66	2/12879 (0.0%)	0.68	4/17440 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	68	ARG	CZ-NH1	5.59	1.40	1.33
2	B	108	GLN	CB-CG	5.20	1.66	1.52

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	108	GLN	CA-CB-CG	6.42	127.52	113.40
2	E	277	ARG	NE-CZ-NH1	6.40	123.50	120.30
2	E	331	ARG	NE-CZ-NH1	5.65	123.13	120.30
2	E	277	ARG	NE-CZ-NH2	-5.10	117.75	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3348	0	3294	42	0
1	D	3336	0	3289	39	0
2	B	2930	0	2872	31	0
2	E	2901	0	2837	40	0
3	A	63	0	34	8	0
3	B	63	0	34	1	0
3	D	63	0	34	12	0
3	E	63	0	34	8	0
4	A	16	0	0	0	0
4	B	16	0	0	0	0
4	D	16	0	0	0	0
4	E	16	0	0	3	0
5	A	4	0	0	0	0
5	D	4	0	0	0	0
6	A	229	0	0	5	0
6	B	565	0	0	5	0
6	D	92	0	0	1	0
6	E	85	0	0	1	0
All	All	13810	0	12428	153	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 6.

All (153) 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)
6:B:588:HOH:O	1:D:403[B]:LYS:HE3	1.20	1.30
3:D:580:SRM:HBB2	3:D:580:SRM:HMB1	1.39	1.04
1:A:85:GLY:HA3	1:A:92[A]:GLU:OE2	1.65	0.96
3:D:580:SRM:HBB2	3:D:580:SRM:CMB	1.95	0.95
3:A:580:SRM:O1A	2:B:139:HIS:HD2	1.55	0.88
2:E:134:THR:OG1	2:E:177:CYS:SG	2.32	0.88
1:A:81:TYR:CE2	1:A:93:HIS:CD2	2.63	0.85
1:D:229:ARG:HD2	2:E:184:ALA:HB2	1.61	0.82
3:D:580:SRM:HBD2	3:D:580:SRM:HDD1	1.67	0.77
1:D:261:GLU:HB3	1:D:294:LYS:HE2	1.65	0.76
1:D:399:LYS:HD3	1:D:417:TRP:O	1.86	0.76
1:A:5:LEU:HD11	1:A:56:LYS:HE2	1.68	0.76
1:A:81:TYR:HE2	1:A:93:HIS:CD2	2.03	0.75
1:A:81:TYR:CD2	1:A:93:HIS:CD2	2.77	0.71
3:E:570:SRM:HDD2	3:E:570:SRM:HBD1	1.72	0.71
2:E:85:THR:HB	3:E:570:SRM:HAB2	1.73	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:134:THR:HG21	2:E:182:CYS:HB2	1.74	0.69
1:A:56:LYS:HE3	6:A:709:HOH:O	1.93	0.69
3:D:580:SRM:CBB	3:D:580:SRM:HMB1	2.20	0.68
2:B:260:GLU:HB3	1:D:403[B]:LYS:HD2	1.74	0.68
1:D:131:GLY:HA3	3:D:580:SRM:HBB1	1.77	0.67
1:A:66:GLY:H	1:A:81:TYR:HE1	1.42	0.67
2:E:125:GLU:HB3	2:E:165:HIS:HB3	1.77	0.66
3:A:580:SRM:C4C	2:B:182:CYS:HA	2.26	0.65
3:A:580:SRM:O1A	2:B:139:HIS:CD2	2.45	0.65
1:A:64:HIS:HD2	6:A:525:HOH:O	1.80	0.65
1:A:94:PHE:O	2:B:139:HIS:HE1	1.80	0.65
2:B:164[B]:ASP:HB3	2:B:166:LYS:HG3	1.80	0.62
1:D:235:ILE:HD12	1:D:310:LEU:HD22	1.81	0.61
1:A:94:PHE:O	2:B:139:HIS:CE1	2.55	0.60
1:A:14:TRP:CD2	1:A:15:PRO:HD2	2.37	0.60
1:D:96:THR:HG23	2:E:139:HIS:CE1	2.38	0.58
1:D:315:ALA:HB1	1:D:316:PRO:CD	2.33	0.58
1:D:288:CYS:O	1:D:289:MET:HB2	2.02	0.58
2:E:172:ARG:HH21	3:E:570:SRM:C2C	2.17	0.58
6:A:511:HOH:O	2:E:331:ARG:HD2	2.03	0.57
2:E:20:HIS:HB3	2:E:23:ILE:HD12	1.86	0.57
2:B:313:ASN:HB3	6:B:420:HOH:O	2.04	0.57
2:B:175:LEU:CD2	6:B:508:HOH:O	2.53	0.56
1:A:65:GLY:HA2	1:A:81:TYR:HD1	1.70	0.56
2:E:176:ALA:HB2	2:E:185:VAL:HG21	1.87	0.55
2:E:134:THR:CG2	2:E:182:CYS:HB2	2.37	0.55
1:A:235:ILE:HD12	1:A:310:LEU:HD22	1.88	0.54
1:A:243:LYS:HE3	1:A:301:PRO:O	2.07	0.53
1:A:315:ALA:HB1	1:A:316:PRO:CD	2.38	0.53
1:D:22:LYS:O	1:D:26:GLU:HG3	2.08	0.53
1:D:315:ALA:HB1	1:D:316:PRO:HD3	1.90	0.53
3:D:580:SRM:HMB3	3:D:580:SRM:HBB2	1.89	0.53
2:E:134:THR:HB	4:E:585:SF4:S4	2.50	0.52
3:E:570:SRM:HDD2	3:E:570:SRM:CBD	2.40	0.52
2:B:175:LEU:HD23	2:B:175:LEU:C	2.30	0.51
1:D:249:VAL:HG11	1:D:278:LEU:HB2	1.93	0.51
1:A:65:GLY:CA	1:A:81:TYR:HD1	2.23	0.51
1:D:261:GLU:CB	1:D:294:LYS:HE2	2.37	0.51
2:B:85:THR:HB	3:B:570:SRM:HAB1	1.93	0.51
2:E:324:VAL:HG11	2:E:329:TRP:CE2	2.45	0.51
1:D:238:TRP:CE3	1:D:301:PRO:HB3	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:40:MET:SD	1:D:141:THR:HA	2.52	0.50
1:A:40:MET:SD	1:A:141:THR:HA	2.51	0.49
1:D:316:PRO:HG2	2:E:181:MET:CE	2.42	0.49
1:A:85:GLY:HA3	1:A:92[B]:GLU:HG3	1.95	0.49
1:D:262:VAL:HG22	1:D:294:LYS:HG2	1.93	0.49
1:A:68:VAL:HG22	1:A:146:LEU:CD1	2.42	0.49
1:A:81:TYR:CD2	1:A:93:HIS:HD2	2.30	0.49
2:E:326:ARG:HD2	6:E:793:HOH:O	2.12	0.49
1:D:78:ILE:HB	1:D:96:THR:OG1	2.13	0.48
2:B:324:VAL:HG11	2:B:329:TRP:CE2	2.48	0.48
3:A:580:SRM:C3C	2:B:182:CYS:HA	2.43	0.48
1:D:128:ASN:HD22	2:E:61:LEU:HA	1.78	0.48
1:D:316:PRO:HG2	2:E:181:MET:HE3	1.95	0.48
1:A:243:LYS:HD2	1:A:300:LYS:HE3	1.94	0.47
1:A:186:TYR:CD1	1:A:333:VAL:HG11	2.50	0.47
1:A:81:TYR:HD2	1:A:93:HIS:HD2	1.61	0.47
2:E:142:THR:N	2:E:143:PRO:CD	2.77	0.47
2:E:51:ILE:HG13	2:E:52:TYR:N	2.30	0.47
1:A:147:GLN:HB3	1:A:148:PRO:HD3	1.97	0.47
2:E:87:ARG:HG3	3:E:570:SRM:HBB2	1.97	0.47
1:A:81:TYR:CE2	1:A:93:HIS:CG	3.01	0.46
1:D:14:TRP:CD2	1:D:15:PRO:HD2	2.50	0.46
2:E:177:CYS:HB2	4:E:585:SF4:S3	2.55	0.46
1:A:63:LYS:HG2	1:A:81:TYR:O	2.15	0.46
1:D:330:PHE:HB2	2:E:366:TRP:CH2	2.51	0.46
1:D:310:LEU:HD12	1:D:326:VAL:HA	1.98	0.46
1:A:315:ALA:HB1	1:A:316:PRO:HD2	1.98	0.45
3:A:580:SRM:C1C	2:B:182:CYS:HA	2.46	0.45
3:A:580:SRM:NC	2:B:182:CYS:HA	2.31	0.45
1:D:128:ASN:HB2	1:D:137:ILE:HB	1.97	0.45
1:A:330:PHE:HB2	2:B:366:TRP:CH2	2.52	0.45
1:D:414:ARG:HD3	6:D:771:HOH:O	2.16	0.45
2:B:175:LEU:HD22	6:B:508:HOH:O	2.14	0.45
2:E:73:ILE:HG22	2:E:92:PHE:CE2	2.51	0.45
3:D:580:SRM:HDA2	4:E:585:SF4:S1	2.57	0.45
2:E:164:ASP:CG	2:E:165:HIS:H	2.20	0.45
1:A:81:TYR:HD2	1:A:93:HIS:CD2	2.31	0.45
2:B:175:LEU:HD23	6:B:508:HOH:O	2.14	0.45
3:A:580:SRM:CMB	3:A:580:SRM:CBB	2.95	0.45
2:B:228:ASP:HB2	2:B:235:LYS:HG3	1.99	0.45
3:E:570:SRM:HHA	3:E:570:SRM:O4D	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:570:SRM:HAC2	3:E:570:SRM:HCD1	1.76	0.44
2:B:82:LEU:HD23	2:B:82:LEU:N	2.33	0.44
1:A:81:TYR:HE2	1:A:93:HIS:CG	2.33	0.44
1:D:317:PHE:HD2	2:E:180:ASN:HD22	1.65	0.44
2:E:242:MET:C	2:E:243:TYR:CG	2.91	0.44
1:D:90:GLU:CD	1:D:90:GLU:H	2.20	0.44
1:A:65:GLY:HA2	1:A:81:TYR:CD1	2.51	0.44
1:A:119:TRP:CH2	1:A:141:THR:HB	2.53	0.44
1:A:168:ASP:HB3	1:A:206:PRO:HG3	2.00	0.44
2:E:330:GLU:O	2:E:334:GLU:HG3	2.17	0.44
1:D:129:PHE:HA	1:D:130:HIS:HA	1.81	0.43
2:B:142:THR:N	2:B:143:PRO:CD	2.81	0.43
2:B:242:MET:C	2:B:243:TYR:CG	2.91	0.43
3:D:580:SRM:HDD1	3:D:580:SRM:CBD	2.45	0.43
1:D:131:GLY:CA	3:D:580:SRM:HBB1	2.45	0.43
3:A:580:SRM:HHB	3:A:580:SRM:HBA1	2.00	0.43
1:D:133:THR:H	3:D:580:SRM:HAB2	1.84	0.43
2:B:242:MET:SD	2:B:244:CYS:HB3	2.59	0.43
3:D:580:SRM:HHB	3:D:580:SRM:HDB1	1.84	0.42
2:E:274:GLU:OE1	2:E:361:SER:OG	2.33	0.42
2:E:316:LYS:HG2	2:E:317:HIS:CD2	2.54	0.42
2:B:353:TYR:HA	2:E:353:TYR:HA	2.00	0.42
1:A:144[A]:GLU:CD	1:A:144[A]:GLU:H	2.22	0.42
2:E:54:VAL:O	2:E:91:GLU:HA	2.19	0.42
2:B:164[A]:ASP:HB3	2:B:166:LYS:HG3	2.01	0.42
1:A:249:VAL:HG11	1:A:278:LEU:HB2	2.01	0.42
2:B:65:TYR:O	2:B:69[A]:GLU:HG3	2.20	0.41
1:D:186:TYR:CD1	1:D:333:VAL:HG11	2.55	0.41
1:D:178:PRO:HG3	1:D:187:ASP:HA	2.02	0.41
1:A:250:LYS:NZ	1:A:276:LYS:O	2.46	0.41
1:D:147:GLN:HB3	1:D:148:PRO:HD3	2.02	0.41
1:A:405:SER:OG	1:A:408:THR:HG23	2.19	0.41
1:D:118:VAL:HG11	1:D:153:LEU:HG	2.01	0.41
1:A:56:LYS:CE	6:A:709:HOH:O	2.62	0.41
2:E:55:ARG:HH22	3:E:570:SRM:HBA2	1.86	0.41
2:E:274:GLU:O	2:E:363:GLN:NE2	2.53	0.41
2:B:119:TRP:HB2	2:B:126:TYR:CD2	2.56	0.41
1:D:292:ILE:HG12	1:D:299:LEU:HB3	2.03	0.41
1:A:63:LYS:HE2	1:A:81:TYR:HB3	2.03	0.41
2:B:176:ALA:HB1	2:B:181:MET:HA	2.02	0.41
1:A:326:VAL:HB	1:A:385:MET:HA	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:191:ALA:HB3	2:E:267:MET:HB2	2.03	0.41
1:D:174:ALA:HB2	1:D:192:CYS:HB2	2.03	0.41
2:E:124:GLY:HA3	2:E:316:LYS:CD	2.51	0.41
1:A:73:TYR:CD1	1:A:204:HIS:HB3	2.56	0.41
2:E:277:ARG:HG2	2:E:322:GLU:HG2	2.03	0.40
2:B:260:GLU:HB3	1:D:403[B]:LYS:CD	2.46	0.40
1:D:132:SER:HB3	3:D:580:SRM:HCD2	2.02	0.40
2:B:220:CYS:HA	2:B:221:PRO:HD3	1.96	0.40
1:A:387:LYS:NZ	6:A:1082:HOH:O	2.53	0.40
2:E:242:MET:SD	2:E:244:CYS:HB3	2.62	0.40
2:E:176:ALA:HB1	2:E:181:MET:HA	2.04	0.40
2:E:142:THR:N	2:E:143:PRO:HD3	2.37	0.40
2:E:292:GLU:HB2	2:E:296:TRP:HA	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	418/418 (100%)	408 (98%)	10 (2%)	0	100	100
1	D	416/418 (100%)	404 (97%)	12 (3%)	0	100	100
2	B	366/366 (100%)	353 (96%)	12 (3%)	1 (0%)	46	29
2	E	361/366 (99%)	347 (96%)	13 (4%)	1 (0%)	46	29
All	All	1561/1568 (100%)	1512 (97%)	47 (3%)	2 (0%)	56	38

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
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	356/354 (101%)	350 (98%)	6 (2%)	68	57
1	D	354/354 (100%)	346 (98%)	8 (2%)	58	42
2	B	319/317 (101%)	312 (98%)	7 (2%)	60	45
2	E	314/317 (99%)	306 (98%)	8 (2%)	55	39
All	All	1343/1342 (100%)	1314 (98%)	29 (2%)	61	45

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	86	GLU
1	A	100	ASN
1	A	210	TYR
1	A	229	ARG
1	A	310	LEU
1	A	409	GLU
2	B	97[A]	GLU
2	B	97[B]	GLU
2	B	108	GLN
2	B	177	CYS
2	B	242	MET
2	B	243	TYR
2	B	277	ARG
1	D	100	ASN
1	D	132	SER
1	D	157	GLU
1	D	210	TYR
1	D	294	LYS
1	D	299	LEU
1	D	310	LEU
1	D	413	LYS
2	E	51	ILE
2	E	177	CYS
2	E	230	LYS

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Mol	Chain	Res	Type
2	E	242	MET
2	E	243	TYR
2	E	268	VAL
2	E	283	LYS
2	E	335	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	50	GLN
1	A	64	HIS
2	B	139	HIS
1	D	50	GLN
1	D	61	HIS
1	D	100	ASN
2	E	180	ASN
2	E	363	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

14 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	29,70,70	1.68	8 (27%)	29,112,112	4.49	9 (31%)
5	SO3	A	590	3	1,3,3	0.17	0	0,3,3	0.00	-
3	SRM	B	570	1	29,70,70	2.11	10 (34%)	29,112,112	4.29	15 (51%)
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	0,12,12	0.00	-	0,24,24	0.00	-
3	SRM	D	580	2,5	29,70,70	2.08	9 (31%)	29,112,112	4.12	15 (51%)
5	SO3	D	590	3	1,3,3	0.55	0	0,3,3	0.00	-
3	SRM	E	570	1	29,70,70	2.83	9 (31%)	29,112,112	4.46	16 (55%)
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	SO3	A	590	3	-	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	-	0/0/48/48	0/6/5/5
3	SRM	D	580	2,5	-	0/22/126/126	0/0/8/8
5	SO3	D	590	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 (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	570	SRM	C4A-NA	-5.97	1.27	1.39
3	E	570	SRM	C1C-NC	-5.69	1.29	1.36
3	E	570	SRM	C4C-NC	-5.65	1.29	1.36
3	E	570	SRM	C1B-NB	-5.63	1.29	1.38
3	B	570	SRM	C1B-NB	-4.43	1.31	1.38
3	B	570	SRM	C4A-NA	-4.22	1.31	1.39
3	D	580	SRM	C4A-NA	-3.39	1.32	1.39
3	B	570	SRM	C4C-NC	-3.39	1.32	1.36
3	A	580	SRM	C1B-NB	-2.74	1.34	1.38
3	A	580	SRM	C4B-NB	-2.63	1.34	1.39
3	E	570	SRM	C1A-NA	-2.55	1.34	1.38
3	D	580	SRM	C1A-NA	-2.43	1.34	1.38
3	A	580	SRM	C1A-NA	-2.41	1.34	1.38
3	D	580	SRM	C4B-NB	-2.25	1.34	1.39
3	D	580	SRM	C1B-NB	-2.22	1.34	1.38
3	B	570	SRM	C1C-NC	-2.21	1.33	1.36
3	E	570	SRM	C4B-NB	-2.09	1.35	1.39
3	E	570	SRM	CHC-C4B	-2.08	1.34	1.39
3	B	570	SRM	C1A-NA	-2.05	1.35	1.38
3	B	570	SRM	CAD-C2D	2.27	1.55	1.52
3	A	580	SRM	FE-NA	2.35	2.05	1.95
3	B	570	SRM	CAB-C3B	2.51	1.55	1.51
3	D	580	SRM	CAD-C2D	2.54	1.56	1.52
3	A	580	SRM	C1C-NC	2.58	1.40	1.36
3	A	580	SRM	FE-NB	2.60	2.06	1.95
3	D	580	SRM	FE-NB	3.11	2.08	1.95
3	A	580	SRM	C3C-C2C	3.26	1.47	1.37
3	D	580	SRM	FE-NA	3.30	2.08	1.95
3	B	570	SRM	C3C-C2C	3.62	1.48	1.37
3	D	580	SRM	C3C-C2C	3.94	1.49	1.37
3	A	580	SRM	C3D-C2D	4.18	1.48	1.39
3	B	570	SRM	C3D-C2D	4.38	1.49	1.39
3	B	570	SRM	CAA-C3A	4.42	1.59	1.51
3	E	570	SRM	C3C-C2C	4.79	1.51	1.37
3	E	570	SRM	C3D-C2D	6.31	1.53	1.39
3	D	580	SRM	C3D-C2D	6.60	1.54	1.39

All (55) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	580	SRM	CAB-C3B-C2B	-17.06	104.26	123.46
3	D	580	SRM	CDD-C3D-C4D	-11.29	108.99	127.34
3	B	570	SRM	CAA-C3A-C2A	-11.27	110.78	123.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	570	SRM	CAA-C3A-C2A	-11.03	111.04	123.46
3	D	580	SRM	CAB-C3B-C2B	-9.96	112.25	123.46
3	E	570	SRM	CAC-C3C-C4C	-9.57	116.62	127.01
3	B	570	SRM	CBC-CAC-C3C	-9.20	96.04	112.53
3	E	570	SRM	CDC-C2C-C1C	-8.13	114.25	127.38
3	E	570	SRM	CAD-C2D-C1D	-7.76	118.58	127.01
3	B	570	SRM	C4A-NA-C1A	-7.67	102.74	106.90
3	D	580	SRM	CAD-C2D-C1D	-7.29	119.09	127.01
3	B	570	SRM	CAB-C3B-C2B	-6.62	116.01	123.46
3	E	570	SRM	C4A-NA-C1A	-6.13	103.58	106.90
3	D	580	SRM	CAA-C3A-C2A	-5.97	116.74	123.46
3	A	580	SRM	CAA-C3A-C2A	-5.91	116.81	123.46
3	B	570	SRM	CEC-CDC-C2C	-5.01	106.73	116.31
3	B	570	SRM	C3B-C4B-NB	-4.74	104.81	110.09
3	E	570	SRM	CEC-CDC-C2C	-4.69	107.34	116.31
3	E	570	SRM	CAC-C3C-C2C	-4.62	117.88	129.38
3	E	570	SRM	CBC-CAC-C3C	-4.44	104.57	112.53
3	B	570	SRM	CAB-CBB-CCB	-4.43	104.62	112.75
3	E	570	SRM	CHB-C4A-C3A	-3.80	116.84	125.48
3	A	580	SRM	C3A-C4A-NA	-3.73	105.94	110.09
3	E	570	SRM	CDD-C3D-C4D	-3.61	121.48	127.34
3	A	580	SRM	CAD-CBD-CCD	-3.57	106.20	112.75
3	E	570	SRM	CAD-C2D-C3D	-3.38	120.98	129.38
3	B	570	SRM	CHB-C4A-C3A	-3.08	118.49	125.48
3	D	580	SRM	CAD-C2D-C3D	-3.06	121.76	129.38
3	B	570	SRM	CAD-C2D-C3D	-2.57	122.98	129.38
3	D	580	SRM	CED-CDD-C3D	-2.50	111.54	116.31
3	D	580	SRM	C3A-C4A-NA	-2.31	107.51	110.09
3	B	570	SRM	C1C-CHC-C4B	-2.30	116.88	124.39
3	E	570	SRM	CMA-C2A-CDA	-2.21	107.14	109.76
3	E	570	SRM	C4B-NB-C1B	-2.12	105.75	106.90
3	D	580	SRM	C3B-C4B-NB	2.14	112.46	110.09
3	D	580	SRM	C2A-CDA-CEA	2.20	118.86	115.45
3	D	580	SRM	CHB-C4A-C3A	2.23	130.56	125.48
3	D	580	SRM	CAC-C3C-C4C	2.34	129.54	127.01
3	B	570	SRM	CMB-C2B-CDB	2.38	112.59	109.76
3	E	570	SRM	CMB-C2B-CDB	2.51	112.75	109.76
3	D	580	SRM	CDC-C2C-C1C	2.67	131.69	127.38
3	B	570	SRM	CAD-C2D-C1D	2.80	130.05	127.01
3	A	580	SRM	C3B-C4B-NB	3.31	113.77	110.09
3	E	570	SRM	CAB-CBB-CCB	3.75	119.62	112.75
3	B	570	SRM	CHC-C4B-NB	3.84	131.20	123.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	580	SRM	CMA-C2A-CDA	5.16	115.90	109.76
3	D	580	SRM	CMA-C2A-CDA	5.40	116.19	109.76
3	A	580	SRM	CBD-CAD-C2D	5.57	122.52	112.53
3	B	570	SRM	CBD-CAD-C2D	5.61	122.58	112.53
3	D	580	SRM	CBD-CAD-C2D	5.89	123.08	112.53
3	A	580	SRM	C2A-CDA-CEA	6.55	125.60	115.45
3	D	580	SRM	C4A-NA-C1A	7.31	110.87	106.90
3	E	570	SRM	C3A-C4A-NA	7.40	118.32	110.09
3	B	570	SRM	C3A-C4A-NA	7.57	118.51	110.09
3	A	580	SRM	C4A-NA-C1A	9.98	112.32	106.90

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 31 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	580	SRM	8	0
3	B	570	SRM	1	0
3	D	580	SRM	12	0
3	E	570	SRM	8	0
4	E	585	SF4	3	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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%)	0.01	10 (2%) 62 57	6, 9, 20, 38	0
1	D	417/418 (99%)	0.93	67 (16%) 3 2	4, 10, 18, 25	0
2	B	363/366 (99%)	-0.10	4 (1%) 82 80	5, 9, 16, 46	0
2	E	363/366 (99%)	0.95	60 (16%) 2 2	6, 10, 20, 36	0
All	All	1560/1568 (99%)	0.45	141 (9%) 12 9	4, 9, 19, 46	0

All (141) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	E	4	GLU	9.3
1	D	1	SER	8.7
1	D	81	TYR	6.9
2	E	5	GLY	6.8
2	E	230	LYS	6.6
1	D	67	ILE	6.1
1	A	415	GLY	5.6
1	D	28	MET	5.3
2	E	7	LYS	5.1
1	D	93	HIS	4.9
1	D	33	ALA	4.9
1	D	71	VAL	4.8
1	A	413	LYS	4.8
2	E	46	GLU	4.7
1	D	86	GLU	4.6
2	E	229	MET	4.6
1	A	417	TRP	4.5
2	B	4	GLU	4.5
1	D	122	TRP	4.3
2	E	205	GLU	4.3
1	D	157	GLU	4.0

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Mol	Chain	Res	Type	RSRZ
2	E	28	GLY	4.0
2	E	81	TYR	4.0
2	E	70	LEU	3.8
1	D	12	GLY	3.7
1	D	65	GLY	3.7
2	E	100	ILE	3.7
2	E	123	LYS	3.6
1	D	66	GLY	3.6
2	E	21	PRO	3.5
1	D	37	ASP	3.5
1	D	26	GLU	3.4
2	E	73	ILE	3.4
2	E	109	GLU	3.3
1	D	112	LEU	3.3
2	E	181	MET	3.3
1	D	9	LEU	3.2
1	D	91	VAL	3.2
2	E	98	SER	3.2
1	D	223	CYS	3.2
2	E	103	LEU	3.1
1	D	38	VAL	3.1
2	E	210	THR	3.1
2	E	99	LYS	3.1
2	E	232	LYS	3.0
1	D	74	GLY	3.0
1	D	207	MET	3.0
2	E	182	CYS	3.0
1	D	6	LEU	3.0
1	D	2	GLU	2.9
1	D	251	GLU	2.9
1	A	411	LEU	2.9
2	E	6	VAL	2.9
1	D	30	LYS	2.9
1	D	247	GLU	2.9
1	D	85	GLY	2.9
1	D	123	GLY	2.9
1	D	153	LEU	2.9
1	D	275	GLY	2.9
1	D	276	LYS	2.8
2	B	5	GLY	2.8
2	E	231	ASN	2.8
2	E	207	ILE	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	175	CYS	2.8
2	E	111	VAL	2.7
1	D	255	TRP	2.7
2	E	234	ILE	2.7
1	D	43	GLY	2.7
1	D	13	PRO	2.6
1	D	183	PHE	2.6
1	D	124	SER	2.6
1	A	414	ARG	2.6
2	E	65	TYR	2.6
2	E	67	VAL	2.6
2	E	140	CYS	2.6
2	E	166	LYS	2.6
1	D	115	LEU	2.6
2	E	36	VAL	2.6
1	D	45	ARG	2.6
2	E	211	CYS	2.6
2	E	164	ASP	2.6
1	A	81	TYR	2.6
2	E	77	TYR	2.6
1	D	11	LYS	2.5
1	D	156	LEU	2.5
1	D	191	LEU	2.5
2	E	314	ALA	2.5
1	D	87	GLN	2.5
2	B	229	MET	2.5
2	E	102	ASP	2.5
2	E	209	LYS	2.5
2	E	118	THR	2.5
2	E	30	TRP	2.4
1	D	149	CYS	2.4
1	D	203	LEU	2.4
2	E	96	ASP	2.4
1	D	34	GLU	2.4
1	D	110	LYS	2.4
2	E	235	LYS	2.4
2	E	35	VAL	2.4
1	D	94	PHE	2.4
2	E	29	LYS	2.3
2	E	320	LEU	2.3
1	D	254	SER	2.3
1	A	276	LYS	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	404	PRO	2.3
2	E	339	GLU	2.3
1	D	146	LEU	2.3
2	E	108	GLN	2.3
1	D	297	LYS	2.3
1	D	62	TRP	2.3
2	E	82	LEU	2.3
2	E	72	ASP	2.3
1	D	5	LEU	2.3
2	E	94	VAL	2.3
1	D	142	ARG	2.2
1	D	18	VAL	2.2
1	D	56	LYS	2.2
1	D	29	GLU	2.2
1	D	332	GLU	2.2
2	E	239	GLU	2.2
1	D	155	ASN	2.2
1	D	107	TYR	2.2
1	D	193	TYR	2.2
2	E	78	SER	2.2
2	E	107	VAL	2.2
2	E	116	GLY	2.1
1	D	119	TRP	2.1
2	E	92	PHE	2.1
1	D	72	GLY	2.1
2	E	157	GLU	2.1
2	E	95	THR	2.1
1	D	78	ILE	2.1
2	E	160	GLU	2.1
1	A	400	ASP	2.1
2	E	17	ASP	2.1
1	D	150	PHE	2.1
1	D	184	ALA	2.1
1	A	86	GLU	2.1
2	E	110	ARG	2.0
2	B	230	LYS	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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	SRM	E	570	63/63	0.88	0.35	5.22	8,9,11,13	0
3	SRM	D	580	63/63	0.89	0.35	5.07	12,17,25,26	0
4	SF4	D	576	8/8	0.94	0.24	4.28	6,6,7,7	0
4	SF4	D	575	8/8	0.95	0.28	3.23	6,6,7,7	0
4	SF4	E	585	8/8	0.93	0.28	3.10	9,10,11,12	0
4	SF4	E	586	8/8	0.95	0.21	2.78	9,9,10,10	0
5	SO3	D	590	4/4	0.94	0.25	2.64	48,50,52,56	4
5	SO3	A	590	4/4	0.99	0.10	0.24	18,23,30,30	0
4	SF4	A	576	8/8	0.97	0.12	0.23	5,7,8,8	0
3	SRM	A	580	63/63	0.97	0.08	-1.10	8,10,15,18	0
3	SRM	B	570	63/63	0.97	0.08	-1.15	6,8,11,13	0
4	SF4	B	586	8/8	0.98	0.06	-1.75	7,7,8,8	0
4	SF4	A	575	8/8	0.99	0.04	-3.38	7,8,8,9	0
4	SF4	B	585	8/8	0.99	0.03	-4.21	8,10,13,13	0

6.5 Other polymers ⓘ

There are no such residues in this entry.