



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:04 PM UTC

PDB ID : 1ZJ9 / pdb_00001zj9
Title : Structure of Mycobacterium tuberculosis NirA protein
Authors : Schnell, R.; Sandalova, T.; Hellman, U.; Lindqvist, Y.; Schneider, G.
Deposited on : 2005-04-28
Resolution : 2.90 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

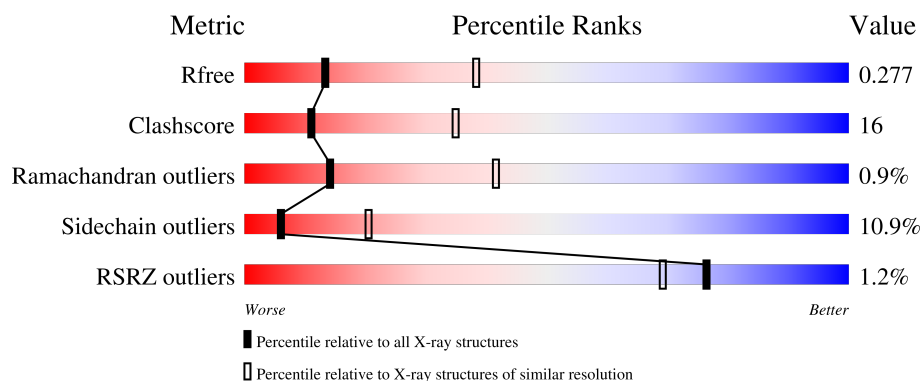
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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}	180053	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	566	2% 62% 30% . .
1	B	566	% 60% 32% . .

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	SF4	A	556	-	-	X	-
3	SF4	B	1556	-	-	X	-
4	SRM	B	1569	X	-	-	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 8792 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 Probable ferredoxin-dependent nitrite reductase NirA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	546	Total	C	N	O	S	0	0	0
			4316	2714	792	797	13			
1	B	546	Total	C	N	O	S	0	0	0
			4316	2714	792	797	13			

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	MET	-	expression tag	UNP P71753
A	-9	ASP	-	expression tag	UNP P71753
A	-8	VAL	-	expression tag	UNP P71753
A	-7	SER	-	expression tag	UNP P71753
A	-6	HIS	-	expression tag	UNP P71753
A	-5	HIS	-	expression tag	UNP P71753
A	-4	HIS	-	expression tag	UNP P71753
A	-3	HIS	-	expression tag	UNP P71753
A	-2	HIS	-	expression tag	UNP P71753
A	-1	HIS	-	expression tag	UNP P71753
A	0	GLY	-	expression tag	UNP P71753
A	1	MET	-	expression tag	UNP P71753
A	2	ALA	-	expression tag	UNP P71753
B	-10	MET	-	expression tag	UNP P71753
B	-9	ASP	-	expression tag	UNP P71753
B	-8	VAL	-	expression tag	UNP P71753
B	-7	SER	-	expression tag	UNP P71753
B	-6	HIS	-	expression tag	UNP P71753
B	-5	HIS	-	expression tag	UNP P71753
B	-4	HIS	-	expression tag	UNP P71753
B	-3	HIS	-	expression tag	UNP P71753
B	-2	HIS	-	expression tag	UNP P71753
B	-1	HIS	-	expression tag	UNP P71753
B	0	GLY	-	expression tag	UNP P71753
B	1	MET	-	expression tag	UNP P71753

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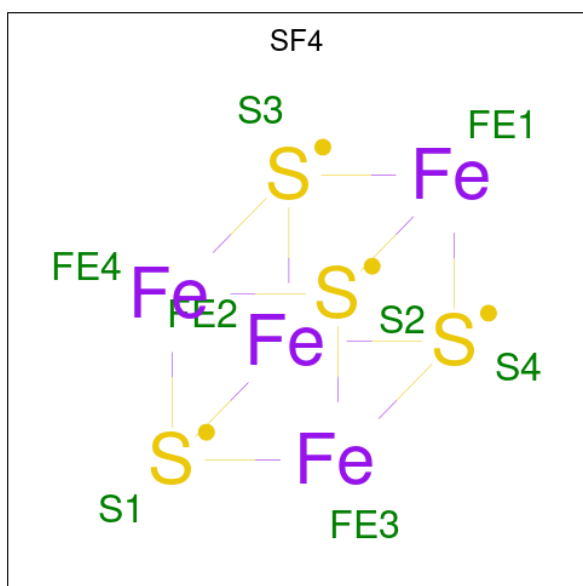
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Chain	Residue	Modelled	Actual	Comment	Reference
B	2	ALA	-	expression tag	UNP P71753

- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

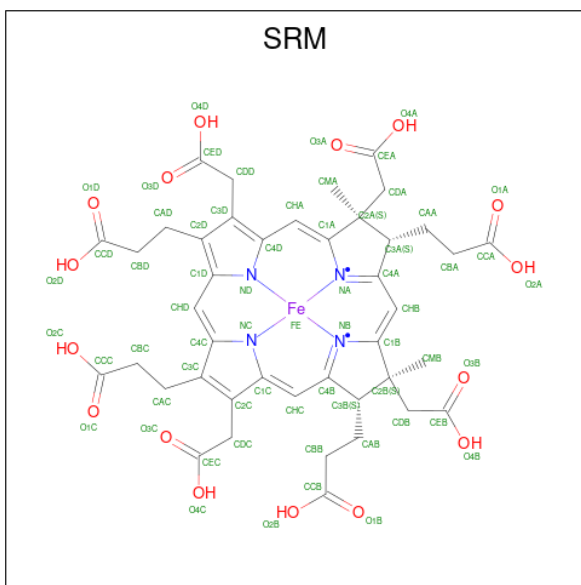
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Cl 1 1	0	0
2	B	1	Total Cl 1 1	0	0

- Molecule 3 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe₄S₄).



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

- Molecule 4 is SIROHEME (CCD ID: SRM) (formula: C₄₂H₄₄FeN₄O₁₆).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 63	C 42	Fe 1	N 4	O 16	0	0
4	B	1	Total 63	C 42	Fe 1	N 4	O 16	0	0

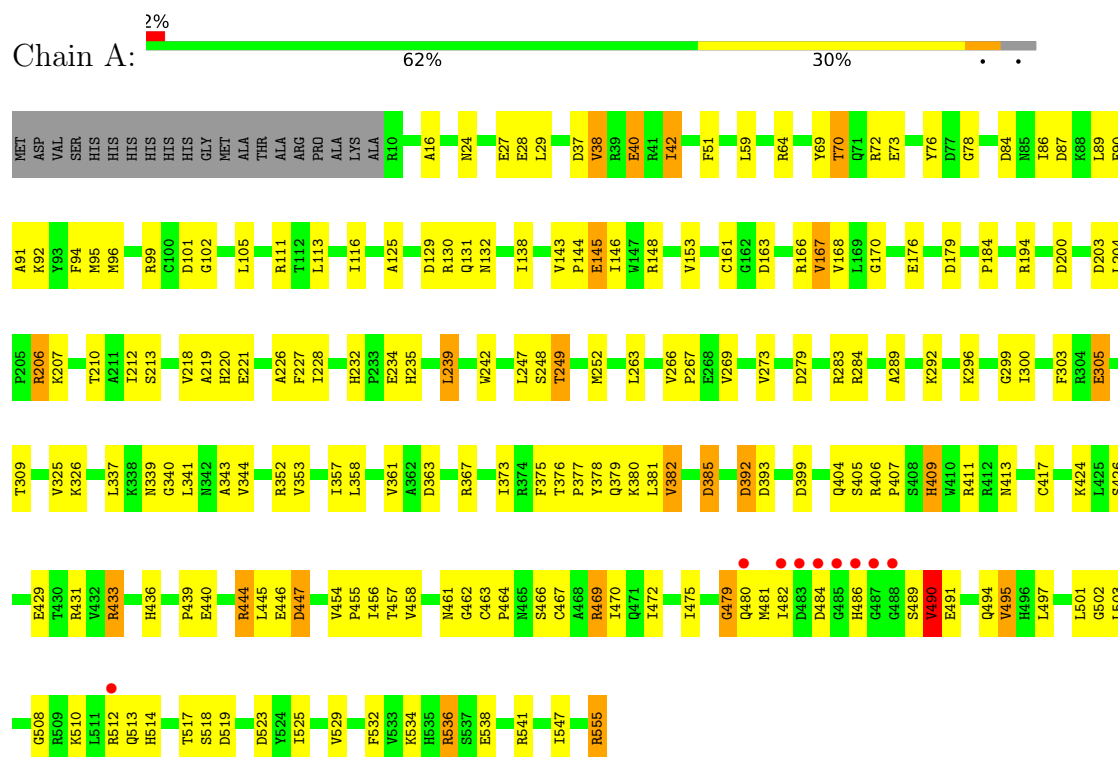
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	13	Total O 13 13	0	0
5	B	3	Total O 3 3	0	0

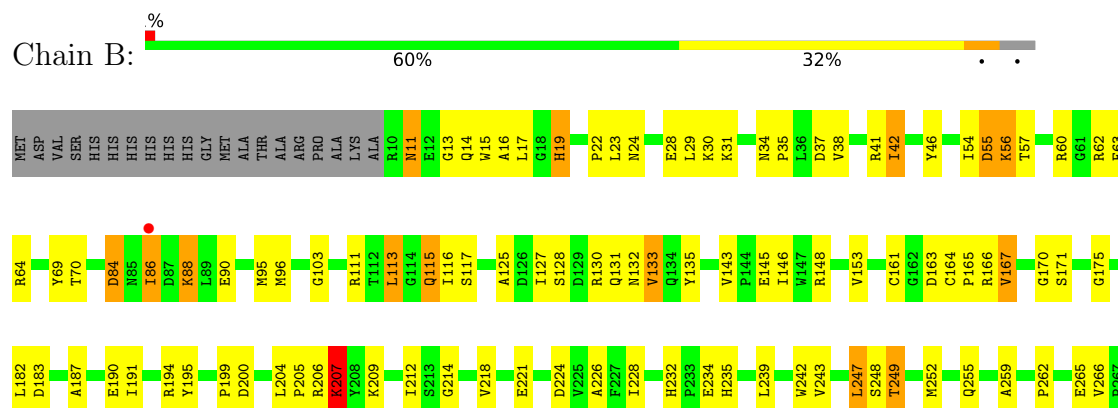
3 Residue-property plots

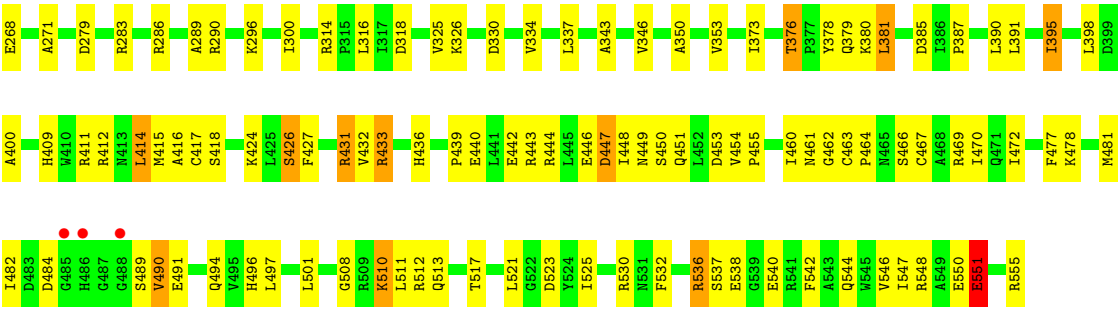
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Probable ferredoxin-dependent nitrite reductase NirA



- Molecule 1: Probable ferredoxin-dependent nitrite reductase NirA





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	59.98Å 83.31Å 108.80Å 90.00° 102.22° 90.00°	Depositor
Resolution (Å)	55.00 – 2.90 55.00 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.0 (55.00-2.90) 99.0 (55.00-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.07 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.202 , 0.280 0.204 , 0.277	Depositor DCC
R_{free} test set	1203 reflections (5.13%)	wwPDB-VP
Wilson B-factor (Å ²)	42.3	Xtriage
Anisotropy	0.335	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 25.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8792	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.53% 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, CL, 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	0.82	3/4409 (0.1%)	0.99	8/5974 (0.1%)
1	B	0.68	1/4409 (0.0%)	0.98	2/5974 (0.0%)
All	All	0.75	4/8818 (0.0%)	0.99	10/11948 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	479	GLY	C-N	28.64	1.73	1.33
1	A	490	VAL	C-N	-9.49	1.20	1.33
1	A	239	LEU	C-N	6.56	1.43	1.33
1	B	143	VAL	CA-CB	5.24	1.56	1.54

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	479	GLY	O-C-N	-8.91	111.11	122.70
1	A	385	ASP	CA-C-N	-6.67	117.78	122.59
1	A	385	ASP	C-N-CA	-6.67	117.78	122.59
1	A	249	THR	N-CA-C	-6.64	103.65	113.61
1	B	249	THR	N-CA-C	-6.50	104.58	113.30
1	A	213	SER	N-CA-C	5.83	117.72	108.79
1	A	490	VAL	O-C-N	5.49	129.43	122.57
1	B	449	ASN	N-CA-C	5.38	116.83	111.07
1	A	299	GLY	N-CA-C	-5.09	106.32	112.48
1	A	340	GLY	N-CA-C	-5.08	108.61	115.21

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	4316	0	4255	129	0
1	B	4316	0	4256	153	0
2	A	1	0	0	1	0
2	B	1	0	0	1	0
3	A	8	0	0	6	0
3	B	8	0	0	5	0
4	A	63	0	34	8	0
4	B	63	0	34	10	0
5	A	13	0	0	0	0
5	B	3	0	0	0	0
All	All	8792	0	8579	285	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 16.

All (285) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:479:GLY:C	1:A:480:GLN:N	1.73	1.46
1:A:96:MET:HE1	1:A:146:ILE:HG21	1.31	1.11
1:B:469:ARG:HG2	1:B:472:ILE:HD11	1.42	1.00
1:A:132:ASN:ND2	4:A:569:SRM:O1B	1.99	0.96
1:A:463:CYS:HB2	1:A:464:PRO:HD2	1.45	0.96
1:A:426:SER:CB	3:A:556:SF4:S2	2.60	0.89
1:A:426:SER:HB2	3:A:556:SF4:S2	2.12	0.89
1:A:239:LEU:HD13	1:A:266:VAL:HG21	1.54	0.88
1:B:376:THR:HG22	1:B:378:TYR:H	1.35	0.88
1:B:462:GLY:HA3	3:B:1556:SF4:S2	2.13	0.88
1:A:376:THR:HG22	1:A:378:TYR:H	1.40	0.87
1:B:431:ARG:HG3	1:B:431:ARG:HH11	1.40	0.85
1:B:207:LYS:NZ	2:B:1570:CL:CL	2.47	0.84
1:A:206:ARG:NH2	4:A:569:SRM:O4D	2.15	0.80
1:A:490:VAL:HG12	1:A:491:GLU:H	1.47	0.79
1:A:70:THR:HG23	1:A:424:LYS:HZ2	1.47	0.78
1:B:22:PRO:HG3	1:B:31:LYS:HG3	1.65	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:426:SER:HB3	3:A:556:SF4:S2	2.25	0.77
1:A:440:GLU:O	1:A:444:ARG:HB2	1.86	0.75
1:B:132:ASN:HD21	1:B:167:VAL:CG2	1.99	0.75
1:A:232:HIS:CD2	1:A:235:HIS:H	2.05	0.75
1:B:409:HIS:ND1	1:B:455:PRO:HB3	2.03	0.74
1:B:426:SER:HB3	3:B:1556:SF4:S2	2.27	0.74
1:A:70:THR:HG23	1:A:424:LYS:NZ	2.06	0.71
1:A:357:ILE:O	1:A:361:VAL:HG23	1.91	0.71
4:A:569:SRM:CED	4:A:569:SRM:HBD2	2.22	0.70
1:B:337:LEU:HD21	1:B:343:ALA:HB2	1.73	0.70
1:B:224:ASP:OD2	1:B:290:ARG:NH2	2.24	0.70
1:B:206:ARG:HD2	1:B:289:ALA:HB3	1.74	0.70
1:B:11:ASN:OD1	1:B:19:HIS:HB3	1.91	0.69
1:A:405:SER:C	1:A:407:PRO:HD3	2.18	0.69
1:A:462:GLY:HA3	3:A:556:SF4:S2	2.33	0.69
1:A:96:MET:HE1	1:A:146:ILE:CG2	2.17	0.69
1:B:431:ARG:HG3	1:B:431:ARG:NH1	2.03	0.69
1:A:132:ASN:OD1	1:A:167:VAL:HG11	1.93	0.69
1:B:113:LEU:HD12	1:B:135:TYR:OH	1.93	0.69
1:B:30:LYS:NZ	1:B:204:LEU:O	2.26	0.69
1:B:175:GLY:HA2	1:B:183:ASP:OD2	1.92	0.69
1:B:34:ASN:HB3	1:B:37:ASP:OD2	1.94	0.68
1:B:521:LEU:O	1:B:525:ILE:HG12	1.93	0.68
1:B:132:ASN:ND2	1:B:167:VAL:HG21	2.08	0.67
1:A:221:GLU:HB2	1:A:501:LEU:HD22	1.75	0.67
1:A:232:HIS:HD2	1:A:235:HIS:H	1.41	0.67
1:B:373:ILE:HD11	1:B:381:LEU:HD12	1.76	0.67
1:B:444:ARG:HH22	1:B:523:ASP:HA	1.60	0.67
1:B:417:CYS:HB2	3:B:1556:SF4:S1	2.35	0.66
1:A:463:CYS:CB	1:A:464:PRO:HD2	2.24	0.66
1:B:46:TYR:HB3	1:B:54:ILE:HG12	1.78	0.66
1:B:463:CYS:HB2	1:B:464:PRO:HD2	1.78	0.65
1:B:544:GLN:O	1:B:547:ILE:HG13	1.97	0.65
1:B:221:GLU:HB2	1:B:501:LEU:HD22	1.78	0.64
1:B:255:GLN:HB2	1:B:300:ILE:HD11	1.80	0.64
1:A:417:CYS:HB2	3:A:556:SF4:S1	2.38	0.64
1:A:203:ASP:OD1	1:A:283:ARG:NH2	2.32	0.63
1:A:96:MET:CE	1:A:146:ILE:HG21	2.19	0.63
1:B:132:ASN:HD21	1:B:167:VAL:HG22	1.64	0.63
1:B:38:VAL:O	1:B:42:ILE:HG12	1.99	0.62
1:B:132:ASN:HD21	1:B:167:VAL:HG21	1.65	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:166:ARG:NH2	1:B:207:LYS:HB3	2.13	0.62
1:A:378:TYR:O	1:A:380:LYS:HD2	1.98	0.62
1:B:481:MET:HE2	1:B:491:GLU:HA	1.80	0.62
1:B:470:ILE:HD13	1:B:478:LYS:HG3	1.81	0.62
1:B:111:ARG:HG2	1:B:115:GLN:NE2	2.16	0.60
1:A:252:MET:HE1	1:A:292:LYS:O	2.01	0.60
1:A:105:LEU:HB2	1:A:373:ILE:HG23	1.83	0.60
1:B:164:CYS:HB2	1:B:165:PRO:CD	2.31	0.60
1:B:207:LYS:HE3	4:B:1569:SRM:O3D	2.01	0.60
1:A:145:GLU:OE2	1:A:148:ARG:NH2	2.35	0.60
1:A:64:ARG:O	1:A:161:CYS:HB2	2.01	0.59
1:B:268:GLU:OE2	1:B:314:ARG:NH2	2.27	0.59
1:B:34:ASN:OD1	1:B:35:PRO:HD2	2.03	0.59
1:B:46:TYR:CD1	1:B:54:ILE:HA	2.37	0.59
1:A:227:PHE:C	1:A:228:ILE:HD12	2.27	0.59
1:B:381:LEU:N	1:B:381:LEU:HD23	2.18	0.59
1:A:494:GLN:HE22	1:A:510:LYS:HE3	1.66	0.59
1:B:38:VAL:HG21	1:B:62:ARG:CB	2.33	0.59
1:A:469:ARG:HG2	1:A:472:ILE:HD11	1.85	0.58
1:B:232:HIS:HD2	1:B:234:GLU:H	1.48	0.58
1:A:392:ASP:OD1	1:A:392:ASP:N	2.36	0.58
1:A:454:VAL:CG2	1:A:455:PRO:HD2	2.34	0.57
1:B:462:GLY:CA	3:B:1556:SF4:S2	2.91	0.57
1:A:513:GLN:H	1:A:555:ARG:NH2	2.03	0.57
1:B:96:MET:HE1	1:B:146:ILE:HD13	1.84	0.57
1:B:190:GLU:O	1:B:194:ARG:HB2	2.04	0.57
1:A:532:PHE:O	1:A:536:ARG:HB3	2.05	0.57
1:B:447:ASP:CG	1:B:530:ARG:HH12	2.13	0.57
1:A:131:GLN:HG2	1:A:375:PHE:O	2.05	0.56
1:A:373:ILE:HD11	1:A:381:LEU:HD12	1.86	0.56
1:B:46:TYR:CE1	1:B:54:ILE:HA	2.40	0.56
1:A:168:VAL:HG22	1:A:210:THR:HB	1.87	0.56
1:B:14:GLN:OE1	1:B:17:LEU:HD12	2.06	0.56
1:B:239:LEU:HD13	1:B:266:VAL:HG21	1.86	0.56
1:B:442:GLU:O	1:B:446:GLU:HB2	2.07	0.55
1:B:467:CYS:HA	4:B:1569:SRM:C4C	2.37	0.55
1:B:11:ASN:OD1	1:B:19:HIS:CB	2.54	0.55
1:B:444:ARG:NH2	1:B:523:ASP:HA	2.21	0.55
1:B:247:LEU:HD12	1:B:469:ARG:HG3	1.88	0.55
1:A:176:GLU:HB3	1:A:337:LEU:HB3	1.89	0.55
1:B:128:SER:OG	1:B:132:ASN:HB2	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:206:ARG:HH22	4:A:569:SRM:CED	2.18	0.54
1:B:482:ILE:HD11	1:B:517:THR:HG22	1.89	0.54
4:B:1569:SRM:CED	4:B:1569:SRM:HBD2	2.38	0.54
1:A:73:GLU:HA	1:A:91:ALA:HB2	1.89	0.54
1:B:187:ALA:O	1:B:191:ILE:HG13	2.08	0.54
1:A:206:ARG:HD2	1:A:289:ALA:HB3	1.90	0.54
1:B:532:PHE:O	1:B:536:ARG:HB2	2.07	0.53
1:A:247:LEU:HD12	1:A:469:ARG:HG3	1.90	0.53
1:A:494:GLN:HE22	1:A:510:LYS:CE	2.22	0.53
1:B:496:HIS:CE1	1:B:510:LYS:HD2	2.43	0.53
1:A:433:ARG:NH1	1:A:518:SER:O	2.42	0.53
1:A:490:VAL:CG1	1:A:491:GLU:H	2.19	0.53
1:A:444:ARG:NH2	1:A:523:ASP:OD1	2.38	0.53
1:B:544:GLN:O	1:B:547:ILE:CG1	2.57	0.53
1:A:232:HIS:CD2	1:A:234:GLU:H	2.27	0.53
1:A:454:VAL:HG23	1:A:455:PRO:HD2	1.91	0.53
1:B:467:CYS:HA	4:B:1569:SRM:C3C	2.39	0.53
1:A:339:ASN:ND2	1:A:341:LEU:HD12	2.24	0.52
1:B:461:ASN:OD1	1:B:466:SER:HA	2.10	0.52
1:B:64:ARG:O	1:B:161:CYS:HB2	2.09	0.52
1:B:497:LEU:O	1:B:508:GLY:HA2	2.10	0.52
1:A:163:ASP:HB3	1:A:204:LEU:O	2.09	0.52
1:B:376:THR:HG22	1:B:378:TYR:N	2.14	0.52
1:B:511:LEU:HD22	1:B:551:GLU:OE2	2.08	0.52
1:A:69:TYR:CE2	4:A:569:SRM:HMA2	2.45	0.52
1:B:38:VAL:HG21	1:B:62:ARG:HB2	1.90	0.52
1:A:37:ASP:O	1:A:40:GLU:HG2	2.09	0.52
1:B:127:ILE:HG22	1:B:379:GLN:HA	1.91	0.52
1:A:252:MET:HE1	1:A:296:LYS:HB2	1.90	0.51
1:B:41:ARG:HD2	1:B:46:TYR:OH	2.10	0.51
1:B:113:LEU:HG	1:B:133:VAL:HG21	1.92	0.51
1:B:209:LYS:NZ	4:B:1569:SRM:O1C	2.33	0.51
1:B:232:HIS:HE1	1:B:318:ASP:O	1.93	0.51
1:B:255:GLN:CB	1:B:300:ILE:HD11	2.40	0.51
1:A:429:GLU:HG3	1:A:433:ARG:HB2	1.92	0.51
1:A:513:GLN:H	1:A:555:ARG:HH22	1.57	0.51
1:B:103:GLY:HA2	1:B:131:GLN:HB3	1.92	0.51
1:A:24:ASN:HB2	1:A:27:GLU:H	1.75	0.51
1:A:116:ILE:HD13	1:A:146:ILE:HG23	1.91	0.51
1:B:432:VAL:O	1:B:433:ARG:C	2.54	0.51
1:B:232:HIS:CD2	1:B:234:GLU:H	2.28	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:143:VAL:HB	1:A:144:PRO:HD3	1.93	0.50
1:B:41:ARG:HB3	1:B:46:TYR:HE2	1.76	0.50
1:A:59:LEU:CD1	1:A:94:PHE:HZ	2.24	0.50
1:A:102:GLY:C	1:A:131:GLN:HE21	2.18	0.50
1:A:111:ARG:NH1	1:A:363:ASP:OD1	2.27	0.50
1:A:252:MET:CE	1:A:296:LYS:HB2	2.40	0.50
1:B:175:GLY:CA	1:B:183:ASP:OD2	2.58	0.50
1:A:339:ASN:HB3	1:A:341:LEU:H	1.77	0.50
1:B:55:ASP:OD1	1:B:55:ASP:N	2.41	0.50
1:A:239:LEU:HD13	1:A:266:VAL:CG2	2.35	0.50
1:B:41:ARG:HB3	1:B:46:TYR:CE2	2.47	0.50
1:B:414:LEU:HD23	1:B:415:MET:N	2.26	0.50
1:B:252:MET:HE1	1:B:296:LYS:HB2	1.94	0.49
1:B:38:VAL:HG21	1:B:62:ARG:HB3	1.94	0.49
1:B:13:GLY:HA2	1:B:19:HIS:HB2	1.94	0.49
1:B:35:PRO:HA	1:B:62:ARG:HD3	1.94	0.49
1:A:529:VAL:O	1:A:532:PHE:HB3	2.13	0.49
1:A:458:VAL:HA	1:A:475:ILE:O	2.13	0.49
1:B:427:PHE:O	1:B:481:MET:HE3	2.13	0.49
1:B:463:CYS:HB2	1:B:464:PRO:CD	2.43	0.49
1:A:353:VAL:HG21	1:A:358:LEU:HD21	1.95	0.48
1:B:226:ALA:HB3	1:B:242:TRP:HB2	1.94	0.48
1:A:69:TYR:HB2	1:A:95:MET:HE3	1.95	0.48
1:A:413:ASN:HB3	1:A:457:THR:HG22	1.96	0.48
1:B:116:ILE:HD13	1:B:146:ILE:HG23	1.95	0.48
1:A:436:HIS:O	1:A:439:PRO:HD2	2.14	0.48
1:B:378:TYR:O	1:B:380:LYS:HD2	2.14	0.48
1:A:170:GLY:HA2	1:A:212:ILE:O	2.14	0.47
1:B:538:GLU:H	1:B:538:GLU:CD	2.22	0.47
1:A:273:VAL:HG13	1:A:303:PHE:HZ	1.80	0.47
1:A:490:VAL:HG12	1:A:491:GLU:N	2.24	0.47
1:B:86:ILE:O	1:B:90:GLU:HB2	2.14	0.47
1:A:207:LYS:HZ1	4:A:569:SRM:C1D	2.28	0.47
1:B:195:TYR:CZ	1:B:271:ALA:HB2	2.50	0.47
1:B:346:VAL:HG11	1:B:398:LEU:HD13	1.96	0.47
1:A:16:ALA:HB2	1:A:279:ASP:HB3	1.95	0.47
1:A:24:ASN:O	1:A:28:GLU:HB2	2.15	0.47
1:A:226:ALA:HB3	1:A:242:TRP:HB2	1.96	0.47
1:A:461:ASN:CG	1:A:470:ILE:HG12	2.40	0.47
1:B:163:ASP:OD2	1:B:286:ARG:NH2	2.38	0.47
1:B:111:ARG:HG2	1:B:115:GLN:HE22	1.77	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:130:ARG:HE	1:A:167:VAL:HG23	1.80	0.46
1:A:305:GLU:O	1:A:309:THR:HG23	2.15	0.46
1:B:206:ARG:O	1:B:207:LYS:C	2.57	0.46
1:A:339:ASN:CB	1:A:341:LEU:H	2.29	0.46
1:B:232:HIS:CD2	1:B:235:HIS:H	2.33	0.46
1:A:129:ASP:HB3	1:A:379:GLN:HG3	1.98	0.46
1:A:130:ARG:HB2	4:A:569:SRM:CBB	2.46	0.46
1:B:69:TYR:HB2	1:B:95:MET:HE3	1.98	0.46
1:B:164:CYS:HB2	1:B:165:PRO:HD2	1.97	0.45
1:A:51:PHE:CE1	1:A:92:LYS:HA	2.52	0.45
1:A:166:ARG:NH1	2:A:570:CL:CL	2.85	0.45
1:A:482:ILE:HD11	1:A:517:THR:HG22	1.99	0.45
1:B:70:THR:HG22	1:B:424:LYS:NZ	2.32	0.45
1:B:163:ASP:HB3	1:B:204:LEU:O	2.16	0.45
1:A:73:GLU:N	1:A:89:LEU:O	2.44	0.45
1:A:220:HIS:CB	1:A:228:ILE:HD11	2.45	0.45
1:A:249:THR:HG22	1:A:464:PRO:HB2	1.99	0.45
1:A:232:HIS:HD2	1:A:234:GLU:H	1.63	0.45
1:B:145:GLU:OE2	1:B:148:ARG:NH2	2.50	0.45
1:B:469:ARG:NH1	4:B:1569:SRM:O4C	2.45	0.44
1:A:337:LEU:HD11	1:A:343:ALA:HB2	1.99	0.44
1:B:170:GLY:HA2	1:B:212:ILE:O	2.17	0.44
1:A:466:SER:HA	1:A:470:ILE:HG13	1.99	0.44
1:B:290:ARG:HD3	4:B:1569:SRM:O1D	2.17	0.44
1:B:440:GLU:HG2	1:B:444:ARG:HG3	1.98	0.44
1:B:42:ILE:CD1	1:B:63:PHE:CE1	3.00	0.44
1:B:376:THR:C	1:B:378:TYR:H	2.26	0.44
1:B:84:ASP:OD1	1:B:84:ASP:N	2.49	0.44
1:B:350:ALA:O	1:B:431:ARG:NH1	2.50	0.44
1:B:95:MET:HA	1:B:135:TYR:O	2.17	0.44
1:A:72:ARG:NH2	1:A:76:TYR:O	2.39	0.44
1:B:166:ARG:HH12	1:B:207:LYS:HZ3	1.66	0.44
1:B:431:ARG:HH11	1:B:431:ARG:CG	2.20	0.44
1:B:88:LYS:HA	1:B:88:LYS:HD3	1.42	0.44
1:B:166:ARG:CZ	1:B:207:LYS:HB3	2.47	0.44
1:A:461:ASN:CG	1:A:470:ILE:CG1	2.91	0.43
1:B:426:SER:CB	3:B:1556:SF4:S2	3.02	0.43
1:B:477:PHE:CZ	1:B:525:ILE:HD11	2.53	0.43
1:B:542:PHE:O	1:B:546:VAL:HG23	2.18	0.43
1:B:182:LEU:O	1:B:214:GLY:HA3	2.18	0.43
4:B:1569:SRM:HMB3	4:B:1569:SRM:HAB1	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:406:ARG:N	1:A:407:PRO:HD3	2.33	0.43
1:A:454:VAL:HG21	1:A:541:ARG:HG2	2.01	0.43
1:B:416:ALA:HA	1:B:460:ILE:O	2.18	0.43
1:A:269:VAL:O	1:A:273:VAL:HG23	2.19	0.43
1:B:199:PRO:O	1:B:200:ASP:C	2.62	0.43
1:A:495:VAL:HG23	1:A:514:HIS:HD2	1.83	0.43
1:A:404:GLN:O	1:A:407:PRO:HG3	2.19	0.43
1:B:183:ASP:C	1:B:183:ASP:OD1	2.62	0.43
1:B:334:VAL:HA	1:B:343:ALA:O	2.19	0.43
1:B:387:PRO:HG2	1:B:390:LEU:HD12	2.01	0.43
1:B:540:GLU:OE2	1:B:548:ARG:NH1	2.51	0.43
1:A:125:ALA:O	1:A:353:VAL:O	2.37	0.42
1:B:376:THR:HB	1:B:380:LYS:H	1.83	0.42
1:B:16:ALA:HB2	1:B:279:ASP:HB3	2.00	0.42
1:A:99:ARG:HG3	1:A:167:VAL:HG12	2.01	0.42
1:B:436:HIS:C	1:B:439:PRO:HD2	2.45	0.42
1:A:24:ASN:OD1	1:A:284:ARG:O	2.37	0.42
1:A:78:GLY:HA3	1:A:481:MET:HE3	2.01	0.42
1:A:218:VAL:O	1:A:219:ALA:C	2.62	0.42
1:A:337:LEU:HB2	1:A:339:ASN:HB2	2.00	0.42
1:B:11:ASN:OD1	1:B:19:HIS:CG	2.73	0.42
1:A:445:LEU:C	1:A:447:ASP:N	2.77	0.42
1:A:463:CYS:HB2	1:A:464:PRO:CD	2.30	0.42
1:B:259:ALA:HB1	1:B:316:LEU:HB3	2.02	0.42
1:A:407:PRO:HB3	1:A:411:ARG:HD3	2.02	0.42
1:A:467:CYS:HB2	3:A:556:SF4:S3	2.59	0.42
1:B:391:LEU:O	1:B:395:ILE:HG12	2.19	0.42
1:A:481:MET:HE2	1:A:491:GLU:HA	2.01	0.41
1:B:130:ARG:HB2	4:B:1569:SRM:CBB	2.49	0.41
1:A:131:GLN:OE1	1:A:377:PRO:HG3	2.20	0.41
1:A:352:ARG:NH1	1:A:431:ARG:HG2	2.35	0.41
1:B:409:HIS:CE1	1:B:455:PRO:HB3	2.54	0.41
1:A:497:LEU:O	1:A:508:GLY:HA2	2.20	0.41
1:B:330:ASP:OD2	1:B:412:ARG:NE	2.45	0.41
1:B:56:LYS:HD3	1:B:60:ARG:HH21	1.86	0.41
1:B:96:MET:HE1	1:B:146:ILE:CD1	2.49	0.41
1:B:166:ARG:HH22	1:B:207:LYS:HE2	1.86	0.41
1:B:536:ARG:HG2	1:B:537:SER:N	2.36	0.41
1:A:24:ASN:HD22	1:A:27:GLU:HG2	1.86	0.41
1:A:184:PRO:HA	1:A:263:LEU:HD11	2.03	0.41
1:A:344:VAL:O	1:A:382:VAL:HA	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:95:MET:HE2	4:B:1569:SRM:HAA1	2.02	0.41
1:A:51:PHE:CZ	1:A:92:LYS:HA	2.56	0.41
1:A:266:VAL:HB	1:A:267:PRO:HD3	2.03	0.41
1:A:502:GLY:O	1:A:503:LEU:C	2.62	0.41
1:B:454:VAL:HG22	1:B:455:PRO:HD2	2.03	0.41
1:A:38:VAL:O	1:A:42:ILE:HG12	2.21	0.41
1:A:232:HIS:HD2	1:A:235:HIS:N	2.13	0.41
1:A:467:CYS:HB3	4:A:569:SRM:HDB1	2.03	0.41
1:A:494:GLN:NE2	1:A:510:LYS:HE3	2.33	0.41
1:A:494:GLN:HG2	1:A:495:VAL:N	2.36	0.41
1:B:46:TYR:CD1	1:B:54:ILE:HG12	2.56	0.41
1:B:125:ALA:O	1:B:353:VAL:O	2.39	0.41
1:B:205:PRO:HD3	1:B:283:ARG:NH1	2.35	0.41
1:B:433:ARG:HD3	1:B:521:LEU:HD12	2.03	0.41
1:B:470:ILE:CD1	1:B:478:LYS:HG3	2.49	0.41
1:B:117:SER:HB2	1:B:135:TYR:CE1	2.56	0.41
1:B:447:ASP:OD1	1:B:447:ASP:C	2.64	0.41
1:B:490:VAL:HG12	1:B:491:GLU:H	1.86	0.40
1:B:130:ARG:O	1:B:131:GLN:HB2	2.21	0.40
1:A:463:CYS:CB	1:A:464:PRO:CD	2.94	0.40
1:B:262:PRO:HG2	1:B:265:GLU:HB2	2.03	0.40
1:B:24:ASN:O	1:B:28:GLU:HB2	2.22	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	544/566 (96%)	501 (92%)	39 (7%)	4 (1%)	18	48
1	B	544/566 (96%)	494 (91%)	44 (8%)	6 (1%)	11	36
All	All	1088/1132 (96%)	995 (92%)	83 (8%)	10 (1%)	14	41

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	248	SER
1	B	248	SER
1	A	101	ASP
1	B	56	LYS
1	B	207	LYS
1	A	512	ARG
1	B	15	TRP
1	B	400	ALA
1	A	409	HIS
1	B	551	GLU

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	450/465 (97%)	403 (90%)	47 (10%)	7	23
1	B	450/465 (97%)	399 (89%)	51 (11%)	5	19
All	All	900/930 (97%)	802 (89%)	98 (11%)	6	20

All (98) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	LEU
1	A	38	VAL
1	A	40	GLU
1	A	42	ILE
1	A	70	THR
1	A	84	ASP
1	A	86	ILE
1	A	87	ASP
1	A	90	GLU
1	A	113	LEU
1	A	138	ILE
1	A	145	GLU
1	A	153	VAL

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Mol	Chain	Res	Type
1	A	167	VAL
1	A	179	ASP
1	A	194	ARG
1	A	200	ASP
1	A	206	ARG
1	A	300	ILE
1	A	305	GLU
1	A	325	VAL
1	A	326	LYS
1	A	367	ARG
1	A	382	VAL
1	A	385	ASP
1	A	392	ASP
1	A	393	ASP
1	A	399	ASP
1	A	409	HIS
1	A	433	ARG
1	A	444	ARG
1	A	446	GLU
1	A	447	ASP
1	A	456	ILE
1	A	469	ARG
1	A	484	ASP
1	A	486	HIS
1	A	489	SER
1	A	490	VAL
1	A	495	VAL
1	A	519	ASP
1	A	525	ILE
1	A	534	LYS
1	A	536	ARG
1	A	538	GLU
1	A	547	ILE
1	A	555	ARG
1	B	11	ASN
1	B	19	HIS
1	B	23	LEU
1	B	29	LEU
1	B	42	ILE
1	B	55	ASP
1	B	57	THR
1	B	84	ASP

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Mol	Chain	Res	Type
1	B	86	ILE
1	B	88	LYS
1	B	113	LEU
1	B	115	GLN
1	B	133	VAL
1	B	153	VAL
1	B	167	VAL
1	B	171	SER
1	B	207	LYS
1	B	218	VAL
1	B	228	ILE
1	B	243	VAL
1	B	247	LEU
1	B	249	THR
1	B	325	VAL
1	B	326	LYS
1	B	376	THR
1	B	381	LEU
1	B	385	ASP
1	B	395	ILE
1	B	411	ARG
1	B	414	LEU
1	B	418	SER
1	B	426	SER
1	B	431	ARG
1	B	433	ARG
1	B	443	ARG
1	B	447	ASP
1	B	448	ILE
1	B	450	SER
1	B	451	GLN
1	B	453	ASP
1	B	484	ASP
1	B	489	SER
1	B	490	VAL
1	B	494	GLN
1	B	510	LYS
1	B	512	ARG
1	B	513	GLN
1	B	536	ARG
1	B	550	GLU
1	B	551	GLU

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Mol	Chain	Res	Type
1	B	555	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	ASN
1	A	26	ASN
1	A	131	GLN
1	A	142	ASN
1	A	220	HIS
1	A	231	ASN
1	A	232	HIS
1	A	235	HIS
1	A	404	GLN
1	A	459	ASN
1	A	480	GLN
1	A	486	HIS
1	A	514	HIS
1	B	115	GLN
1	B	132	ASN
1	B	232	HIS
1	B	235	HIS
1	B	327	HIS
1	B	342	ASN
1	B	451	GLN
1	B	496	HIS
1	B	531	ASN
1	B	535	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
3	SF4	B	1556	1	0,12,12	-	-	-		
4	SRM	B	1569	-	68,70,70	1.53	12 (17%)	86,112,112	1.88	24 (27%)
3	SF4	A	556	1	0,12,12	-	-	-		
4	SRM	A	569	-	68,70,70	1.55	14 (20%)	86,112,112	1.84	26 (30%)

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	SRM	A	569	-	-	12/38/126/126	-
3	SF4	B	1556	1	-	-	0/6/5/5
3	SF4	A	556	1	-	-	0/6/5/5
4	SRM	B	1569	-	1/1/19/23	12/38/126/126	-

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	569	SRM	C4C-C3C	-4.37	1.37	1.45
4	B	1569	SRM	C4C-C3C	-4.34	1.37	1.45
4	B	1569	SRM	FE-ND	3.69	2.07	1.95
4	B	1569	SRM	C1C-C2C	-3.59	1.38	1.45
4	A	569	SRM	C1C-C2C	-3.32	1.39	1.45
4	A	569	SRM	C4D-C3D	-2.83	1.38	1.44
4	A	569	SRM	CMA-C2A	-2.64	1.50	1.54
4	A	569	SRM	C1D-ND	-2.62	1.34	1.39
4	B	1569	SRM	O2C-CCC	-2.54	1.22	1.30
4	B	1569	SRM	CMA-C2A	-2.52	1.50	1.54
4	A	569	SRM	C4B-NB	-2.44	1.33	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1569	SRM	C1D-ND	-2.37	1.35	1.39
4	A	569	SRM	C1C-NC	-2.35	1.35	1.39
4	A	569	SRM	O2B-CCB	-2.34	1.23	1.30
4	B	1569	SRM	O4C-CEC	-2.34	1.23	1.30
4	B	1569	SRM	C1D-C2D	-2.32	1.39	1.44
4	B	1569	SRM	C4D-C3D	-2.32	1.39	1.44
4	A	569	SRM	CMB-C2B	-2.32	1.50	1.54
4	A	569	SRM	FE-ND	2.32	2.02	1.95
4	B	1569	SRM	O4A-CEA	-2.22	1.23	1.30
4	B	1569	SRM	C2A-C3A	-2.21	1.49	1.55
4	A	569	SRM	O4B-CEB	-2.20	1.23	1.30
4	A	569	SRM	C2A-C3A	-2.17	1.49	1.55
4	B	1569	SRM	O2B-CCB	-2.14	1.23	1.30
4	A	569	SRM	CDB-C2B	-2.12	1.49	1.56
4	A	569	SRM	CAA-C3A	-2.08	1.49	1.54

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1569	SRM	CAC-C3C-C2C	-5.70	115.86	129.24
4	B	1569	SRM	CAC-C3C-C4C	4.91	134.83	124.85
4	A	569	SRM	CAC-C3C-C2C	-4.40	118.91	129.24
4	B	1569	SRM	CHD-C1D-ND	4.22	131.52	123.86
4	A	569	SRM	CBA-CAA-C3A	-4.02	103.27	114.65
4	A	569	SRM	CHD-C1D-ND	3.94	131.00	123.86
4	A	569	SRM	CAC-C3C-C4C	3.79	132.55	124.85
4	A	569	SRM	CED-CDD-C3D	3.69	123.04	113.76
4	B	1569	SRM	CED-CDD-C3D	3.67	122.99	113.76
4	A	569	SRM	CAD-CBD-CCD	3.59	123.19	113.67
4	A	569	SRM	C1C-C2C-C3C	-3.55	102.87	106.84
4	A	569	SRM	C2C-C1C-NC	3.47	113.67	110.32
4	A	569	SRM	CHA-C4D-C3D	-3.39	117.87	125.30
4	B	1569	SRM	CAD-CBD-CCD	3.36	122.59	113.67
4	B	1569	SRM	CBA-CAA-C3A	-3.26	105.41	114.65
4	B	1569	SRM	C1C-C2C-C3C	-3.16	103.31	106.84
4	B	1569	SRM	CAB-C3B-C4B	-3.13	105.72	111.19
4	A	569	SRM	CAB-C3B-C4B	-3.11	105.75	111.19
4	A	569	SRM	CEC-CDC-C2C	3.02	121.37	113.76
4	B	1569	SRM	O2D-CCD-CBD	3.01	123.52	114.00
4	B	1569	SRM	C2C-C1C-NC	2.89	113.11	110.32
4	B	1569	SRM	CBD-CAD-C2D	2.87	120.48	112.53
4	A	569	SRM	CBC-CAC-C3C	2.83	120.36	112.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1569	SRM	CBC-CAC-C3C	2.82	120.32	112.53
4	A	569	SRM	CAA-C3A-C4A	2.72	115.94	111.19
4	B	1569	SRM	C2B-C1B-CHB	2.47	125.75	123.54
4	B	1569	SRM	CHD-C1D-C2D	-2.46	119.92	125.30
4	B	1569	SRM	C1D-CHD-C4C	-2.45	117.10	124.66
4	B	1569	SRM	O2B-CCB-CBB	2.36	121.47	114.00
4	A	569	SRM	CHD-C1D-C2D	-2.36	120.13	125.30
4	B	1569	SRM	O4D-CED-O3D	-2.33	117.33	123.33
4	A	569	SRM	O2D-CCD-CBD	2.33	121.37	114.00
4	A	569	SRM	O4D-CED-O3D	-2.29	117.44	123.33
4	A	569	SRM	O2B-CCB-CBB	2.29	121.23	114.00
4	B	1569	SRM	CEC-CDC-C2C	2.26	119.45	113.76
4	B	1569	SRM	CHC-C1C-C2C	-2.25	121.32	124.86
4	B	1569	SRM	CMA-C2A-CDA	2.23	114.52	110.74
4	A	569	SRM	CHC-C1C-C2C	-2.22	121.35	124.86
4	A	569	SRM	C1D-CHD-C4C	-2.21	117.84	124.66
4	A	569	SRM	CBD-CAD-C2D	2.19	118.58	112.53
4	A	569	SRM	O4D-CED-CDD	2.14	121.14	114.51
4	A	569	SRM	O4A-CEA-CDA	2.12	121.05	114.35
4	B	1569	SRM	C4C-C3C-C2C	2.10	109.18	106.84
4	A	569	SRM	O2B-CCB-O1B	-2.09	117.95	123.33
4	B	1569	SRM	CHA-C4D-ND	2.06	127.59	123.86
4	B	1569	SRM	C4D-CHA-C1A	-2.04	122.11	125.87
4	A	569	SRM	CMB-C2B-C3B	2.03	115.80	112.05
4	A	569	SRM	C4C-NC-C1C	-2.03	102.51	105.82
4	A	569	SRM	C2A-C3A-C4A	2.03	103.46	100.89
4	B	1569	SRM	O4D-CED-CDD	2.02	120.77	114.51

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	B	1569	SRM	NC

All (24) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	569	SRM	C1A-C2A-CDA-CEA
4	A	569	SRM	CMA-C2A-CDA-CEA
4	B	1569	SRM	C1A-C2A-CDA-CEA
4	B	1569	SRM	CMA-C2A-CDA-CEA
4	A	569	SRM	C4B-C3B-CAB-CBB
4	B	1569	SRM	C4B-C3B-CAB-CBB

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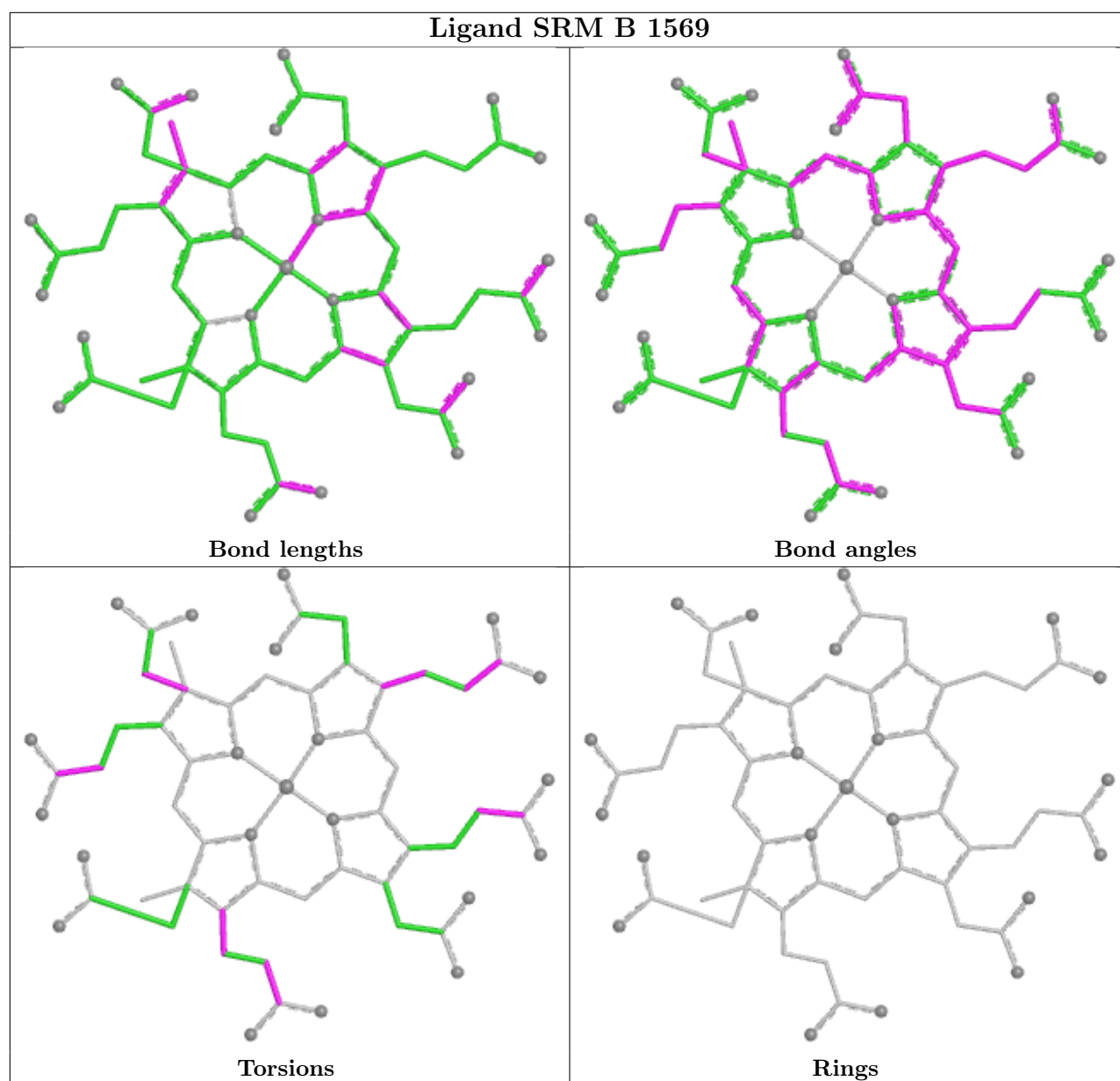
Mol	Chain	Res	Type	Atoms
4	B	1569	SRM	C3D-C2D-CAD-CBD
4	A	569	SRM	CAC-CBC-CCC-O2C
4	A	569	SRM	CAC-CBC-CCC-O1C
4	A	569	SRM	C3A-C2A-CDA-CEA
4	B	1569	SRM	CAC-CBC-CCC-O1C
4	B	1569	SRM	CAD-CBD-CCD-O2D
4	A	569	SRM	CAD-CBD-CCD-O2D
4	A	569	SRM	C3D-C2D-CAD-CBD
4	A	569	SRM	CAB-CBB-CCB-O2B
4	B	1569	SRM	CAD-CBD-CCD-O1D
4	A	569	SRM	CAB-CBB-CCB-O1B
4	A	569	SRM	CAD-CBD-CCD-O1D
4	B	1569	SRM	CAC-CBC-CCC-O2C
4	A	569	SRM	C3A-CAA-CBA-CCA
4	B	1569	SRM	CAB-CBB-CCB-O1B
4	B	1569	SRM	CAB-CBB-CCB-O2B
4	B	1569	SRM	CAA-CBA-CCA-O2A
4	B	1569	SRM	CAA-CBA-CCA-O1A

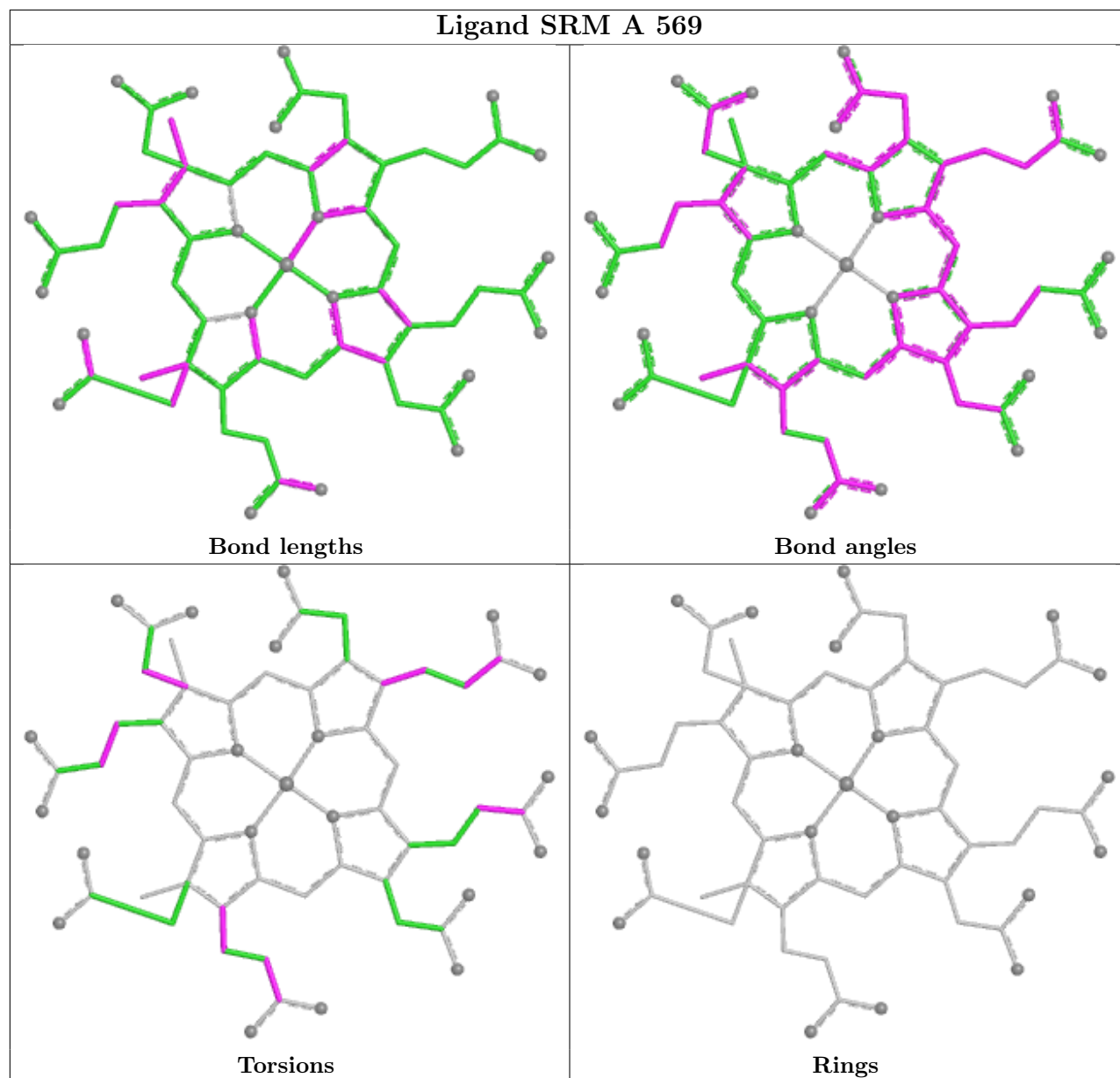
There are no ring outliers.

4 monomers are involved in 29 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1556	SF4	5	0
4	B	1569	SRM	10	0
3	A	556	SF4	6	0
4	A	569	SRM	8	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	479:GLY	C	480:GLN	N	1.73

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	546/566 (96%)	-0.13	9 (1%) 70 62	18, 33, 54, 72	0
1	B	546/566 (96%)	-0.14	4 (0%) 84 79	19, 34, 55, 72	0
All	All	1092/1132 (96%)	-0.13	13 (1%) 76 69	18, 33, 54, 72	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	488	GLY	4.7
1	A	486	HIS	4.4
1	A	480	GLN	4.3
1	A	485	GLY	3.6
1	A	484	ASP	3.5
1	A	487	GLY	3.3
1	B	486	HIS	3.2
1	A	483	ASP	2.8
1	A	482	ILE	2.7
1	A	512	ARG	2.7
1	B	485	GLY	2.5
1	B	86	ILE	2.1
1	B	488	GLY	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 oligosaccharides in this entry.

6.4 Ligands

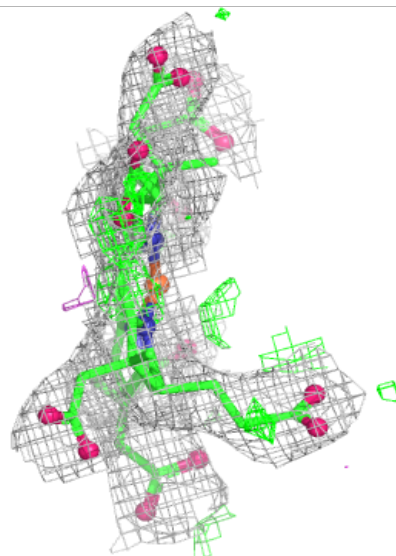
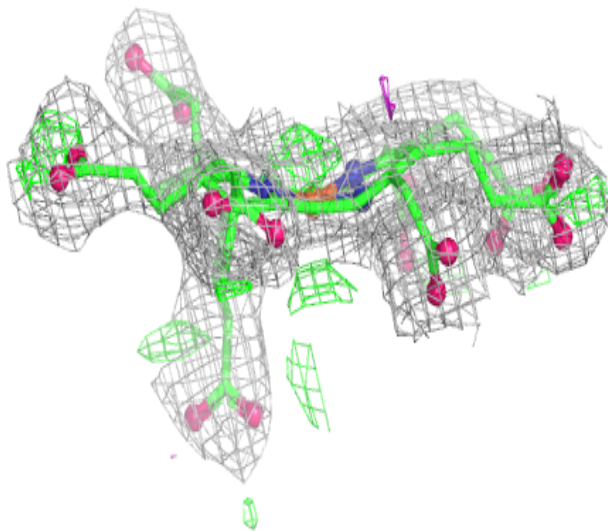
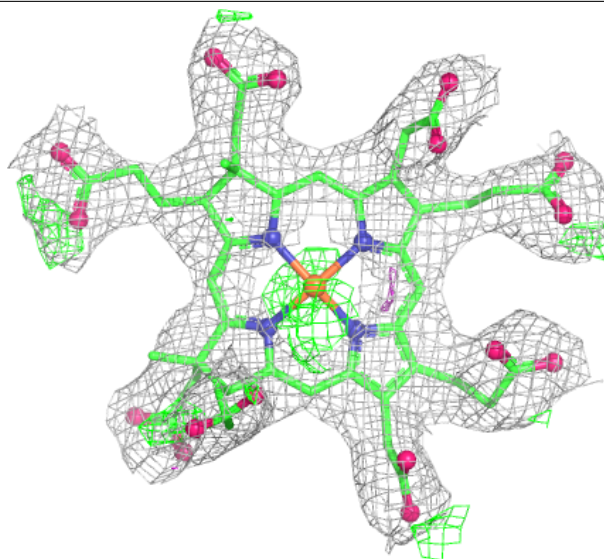
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	CL	A	570	1/1	0.73	0.26	34,34,34,34	0
2	CL	B	1570	1/1	0.91	0.12	32,32,32,32	0
4	SRM	A	569	63/63	0.93	0.13	28,36,42,47	0
4	SRM	B	1569	63/63	0.94	0.10	28,34,42,46	0
3	SF4	A	556	8/8	0.99	0.03	19,20,23,23	0
3	SF4	B	1556	8/8	0.99	0.04	24,27,28,29	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

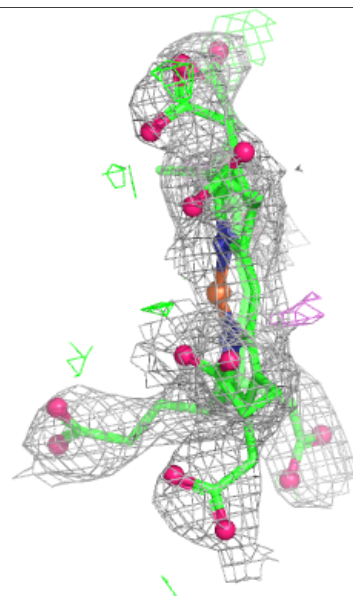
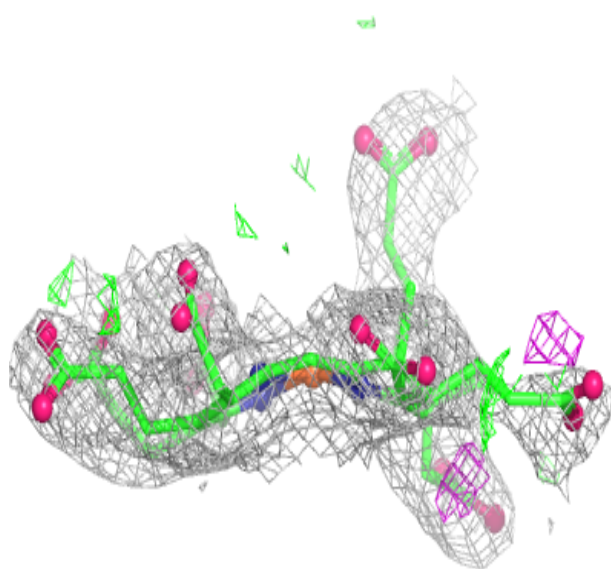
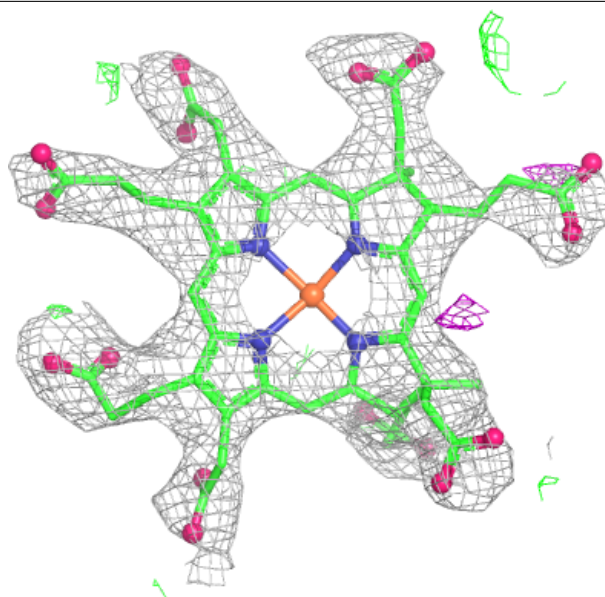
Electron density around SRM A 569:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around SRM B 1569:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers ⓘ

There are no such residues in this entry.