



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:52 AM UTC

PDB ID : 6ELQ / pdb_00006elq
Title : Carbon Monoxide Dehydrogenase IV from Carboxydotherrnus hydrogenofor-
mans
Authors : Domnik, L.; Goetzl, S.; Jeoung, J.H.; Dobbek, H.
Deposited on : 2017-09-29
Resolution : 2.52 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
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

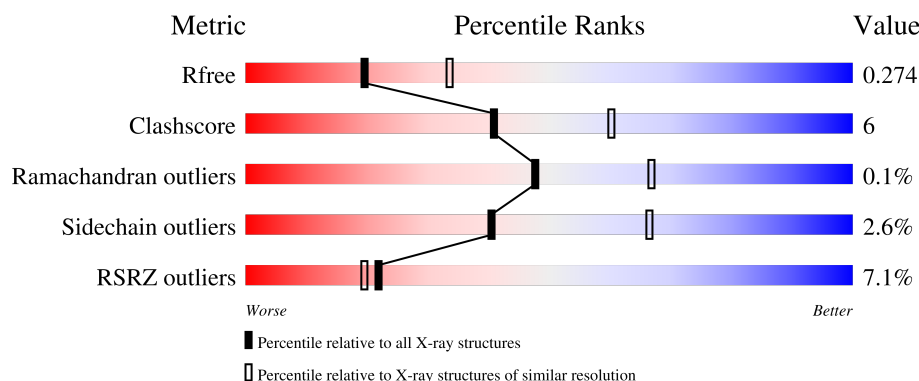
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.52 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	7383 (2.54-2.50)
Clashscore	190562	8079 (2.54-2.50)
Ramachandran outliers	187476	7944 (2.54-2.50)
Sidechain outliers	187428	7946 (2.54-2.50)
RSRZ outliers	180081	7387 (2.54-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	633	85% 14%
1	B	633	20% 75% 23% .
1	X	633	88% 11% .

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	BF8	B	5002	-	-	X	-

2 Entry composition [i](#)

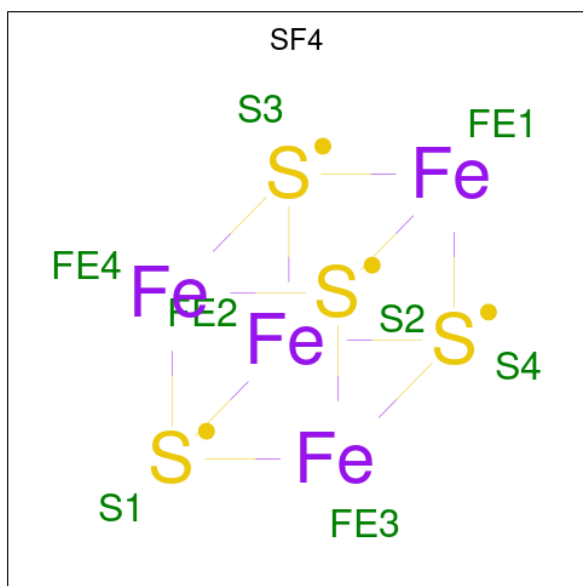
There are 4 unique types of molecules in this entry. The entry contains 14594 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Carbon Monoxide Dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	X	630	Total	C	N	O	S	0	0	0
			4762	3020	821	894	27			
1	A	630	Total	C	N	O	S	0	0	0
			4762	3020	821	894	27			
1	B	630	Total	C	N	O	S	0	0	0
			4762	3020	821	894	27			

- Molecule 2 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe_4S_4).



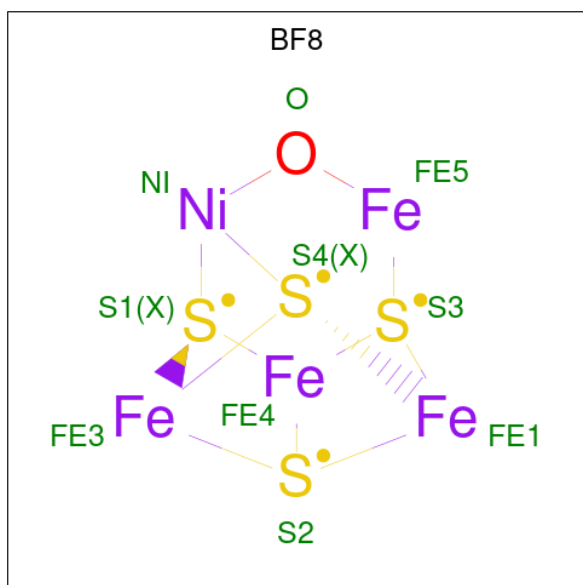
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	X	1	Total	Fe	S	0	0
			8	4	4		
2	A	1	Total	Fe	S	0	0
			8	4	4		
2	A	1	Total	Fe	S	0	0
			8	4	4		

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

- Molecule 3 is FE(4)-NI(1)-S(5) CLUSTER with Oxygen (CCD ID: BF8) (formula: Fe_4NiOS_4).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	X	1	Total	Fe	Ni	O	S	0	0
			10	4	1	1	4		
3	A	1	Total	Fe	Ni	O	S	0	0
			10	4	1	1	4		
3	B	1	Total	Fe	Ni	O	S	0	0
			10	4	1	1	4		

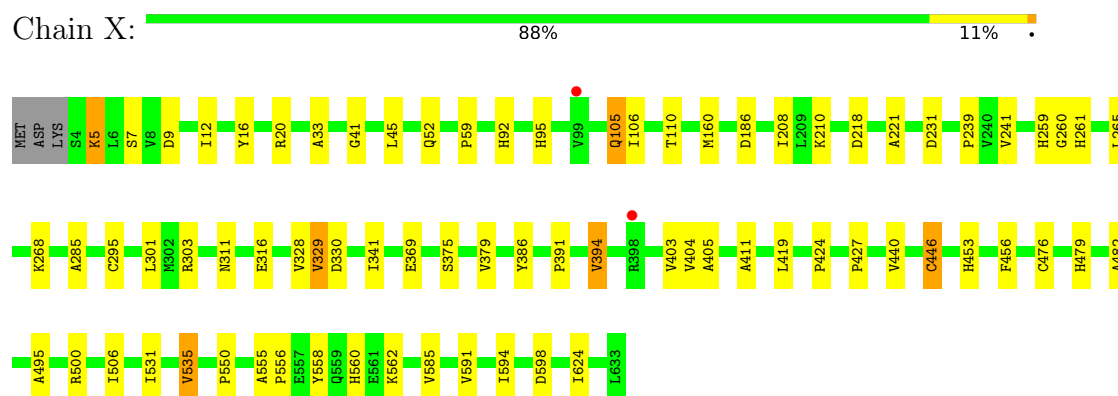
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	X	123	Total	O	0	0
			123	123		
4	A	77	Total	O	0	0
			77	77		
4	B	38	Total	O	0	0
			38	38		

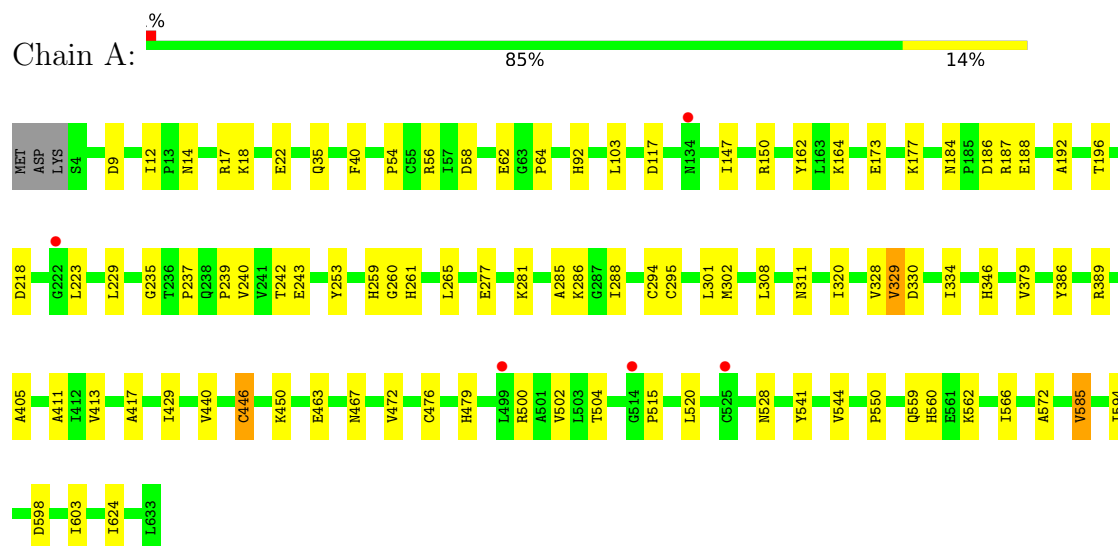
3 Residue-property plots

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.

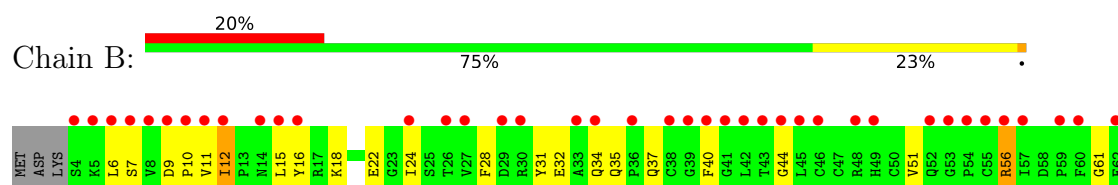
• Molecule 1: Carbon Monoxide Dehydrogenase

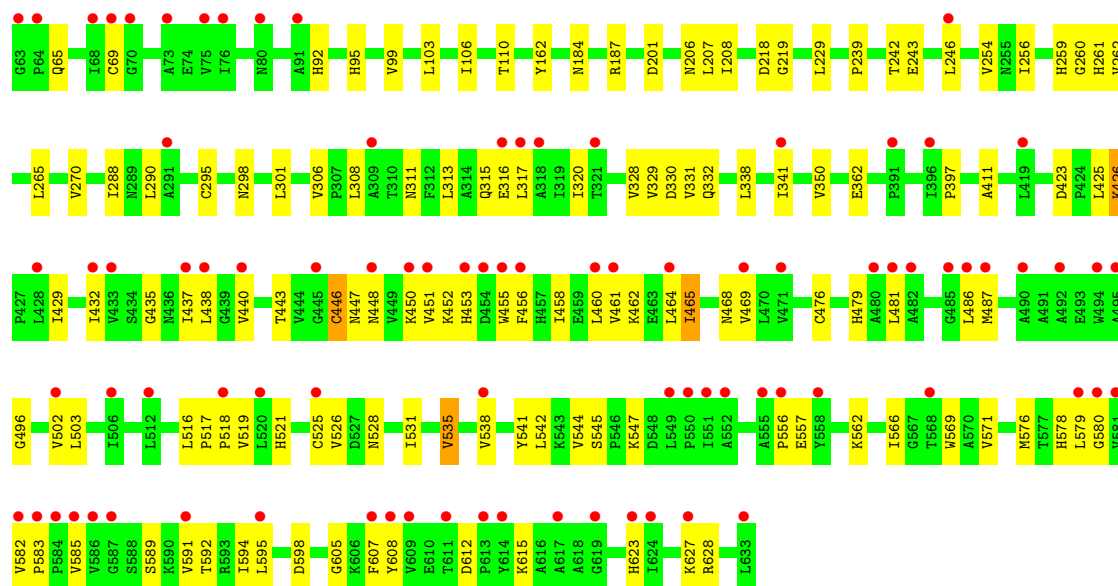


• Molecule 1: Carbon Monoxide Dehydrogenase



• Molecule 1: Carbon Monoxide Dehydrogenase





4 Data and refinement statistics

Property	Value	Source
Space group	P 3 2 1	Depositor
Cell constants a, b, c, α , β , γ	209.15Å 209.15Å 93.41Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	45.63 – 2.52 45.63 – 2.52	Depositor EDS
% Data completeness (in resolution range)	99.9 (45.63-2.52) 99.9 (45.63-2.52)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.53 (at 2.51Å)	Xtriage
Refinement program	PHENIX (1.11.1-2575_1168: ???)	Depositor
R, R_{free}	0.235 , 0.275 0.236 , 0.274	Depositor DCC
R_{free} test set	2113 reflections (2.66%)	wwPDB-VP
Wilson B-factor (Å ²)	49.0	Xtriage
Anisotropy	0.578	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 32.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.092 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	14594	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.62% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SF4, BF8

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	0.11	0/4846	0.31	0/6591
1	B	0.12	0/4846	0.32	0/6591
1	X	0.11	0/4846	0.32	0/6591
All	All	0.11	0/14538	0.32	0/19773

There are no bond length outliers.

There are no bond angle outliers.

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	4762	0	4867	50	0
1	B	4762	0	4866	99	1
1	X	4762	0	4866	42	0
2	A	16	0	0	0	0
2	B	16	0	0	2	0
2	X	8	0	0	0	0
3	A	10	0	0	0	0
3	B	10	0	0	4	0
3	X	10	0	0	1	0
4	A	77	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	38	0	0	0	0
4	X	123	0	0	0	0
All	All	14594	0	14599	187	1

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 (187) 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:B:261:HIS:CE1	3:B:5002:BF8:S3	2.73	0.80
1:B:435:GLY:HA2	1:B:628:ARG:HH22	1.47	0.79
1:B:446:CYS:HA	1:B:476:CYS:HB2	1.69	0.74
1:B:594:ILE:HA	1:B:598:ASP:HB2	1.70	0.73
1:B:612:ASP:HB3	1:B:615:LYS:HB2	1.72	0.72
1:X:33:ALA:HB1	1:A:64:PRO:HG3	1.71	0.71
1:B:465:ILE:HG21	1:B:518:PRO:HB2	1.76	0.67
1:B:259:HIS:HB3	1:B:329:VAL:HG12	1.76	0.66
1:X:45:LEU:HD12	1:X:59:PRO:HG3	1.76	0.66
1:A:117:ASP:HB2	1:A:235:GLY:HA2	1.78	0.66
1:X:7:SER:HB2	1:X:453:HIS:HB3	1.81	0.62
1:B:40:PHE:HB2	2:B:5001:SF4:S3	2.39	0.62
1:B:208:ILE:HD11	1:B:591:VAL:HG13	1.79	0.62
1:B:487:MET:SD	1:B:521:HIS:HB2	2.38	0.62
1:B:481:LEU:HD13	1:B:519:VAL:HG11	1.82	0.62
1:X:186:ASP:HB3	1:A:186:ASP:HB3	1.82	0.60
1:B:571:VAL:HG13	1:B:605:GLY:HA3	1.83	0.60
1:B:7:SER:HB3	1:B:453:HIS:HB3	1.84	0.59
1:X:105:GLN:NE2	1:X:369:GLU:OE2	2.35	0.59
1:B:576:MET:HE3	1:B:608:TYR:HB2	1.85	0.59
1:B:481:LEU:HD23	1:B:486:LEU:HD13	1.83	0.59
1:B:254:VAL:HB	1:B:288:ILE:HD13	1.85	0.58
1:B:468:ASN:OD1	1:B:496:GLY:HA3	2.04	0.58
1:B:585:VAL:HG21	1:B:595:LEU:HD12	1.85	0.58
1:X:446:CYS:HA	1:X:476:CYS:HB2	1.85	0.57
1:X:594:ILE:HA	1:X:598:ASP:HB2	1.87	0.57
1:A:239:PRO:HD3	1:A:411:ALA:HB1	1.87	0.56
1:B:11:VAL:HG11	1:B:246:LEU:HD13	1.85	0.56
1:A:18:LYS:NZ	1:A:22:GLU:OE2	2.38	0.56
1:B:103:LEU:HD22	1:B:229:LEU:HD22	1.87	0.56
1:A:450:LYS:NZ	4:A:5105:HOH:O	2.38	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:311:ASN:HB3	1:X:479:HIS:HB2	1.88	0.55
1:B:330:ASP:OD1	1:B:331:VAL:N	2.37	0.55
1:B:583:PRO:HG2	1:B:585:VAL:HG12	1.87	0.55
1:A:594:ILE:HA	1:A:598:ASP:HB2	1.89	0.54
1:B:316:GLU:HB3	1:B:341:ILE:HD13	1.89	0.54
1:B:61:GLY:HA2	1:B:65:GLN:CD	2.32	0.54
1:X:495:ALA:O	1:X:500:ARG:NH1	2.40	0.54
1:B:265:LEU:HD21	1:B:328:VAL:HG22	1.90	0.54
1:A:446:CYS:HA	1:A:476:CYS:HB2	1.88	0.53
1:B:270:VAL:HG22	1:B:290:LEU:HD13	1.90	0.53
1:A:260:GLY:HA3	1:A:330:ASP:OD1	2.08	0.53
1:A:9:ASP:HB3	1:A:12:ILE:HG12	1.91	0.53
1:A:504:THR:HG22	1:A:515:PRO:HB3	1.90	0.53
1:A:92:HIS:ND1	1:A:218:ASP:OD1	2.35	0.52
1:B:313:LEU:HB3	1:B:453:HIS:CD2	2.44	0.52
1:B:259:HIS:O	1:B:329:VAL:HA	2.10	0.52
1:B:12:ILE:HG13	1:B:317:LEU:HD12	1.91	0.52
1:B:315:GLN:HG3	1:B:338:LEU:HD21	1.92	0.52
1:A:472:VAL:HG12	1:A:520:LEU:HB2	1.90	0.51
1:X:260:GLY:HA3	1:X:330:ASP:OD1	2.11	0.51
1:X:316:GLU:HB3	1:X:341:ILE:HD13	1.91	0.51
1:B:579:LEU:HD22	1:B:583:PRO:HG3	1.92	0.51
1:B:239:PRO:HD3	1:B:411:ALA:HB1	1.92	0.51
1:X:239:PRO:HD3	1:X:411:ALA:HB1	1.93	0.51
1:A:277:GLU:OE1	1:A:288:ILE:N	2.43	0.51
1:A:463:GLU:OE1	1:A:467:ASN:ND2	2.44	0.50
1:B:525:CYS:O	1:B:528:ASN:ND2	2.44	0.50
1:X:231:ASP:OD1	1:X:303:ARG:NH2	2.41	0.50
1:A:311:ASN:HB3	1:A:479:HIS:CB	2.41	0.50
1:A:311:ASN:HB3	1:A:479:HIS:HB2	1.93	0.50
1:B:443:THR:HG21	1:B:461:VAL:HG22	1.92	0.50
1:B:487:MET:HA	1:B:517:PRO:HB3	1.92	0.50
1:X:531:ILE:O	1:X:535:VAL:HG13	2.11	0.49
1:B:531:ILE:O	1:B:535:VAL:HG13	2.12	0.49
1:B:261:HIS:CD2	1:B:295:CYS:HB2	2.48	0.49
1:A:413:VAL:O	1:A:417:ALA:N	2.45	0.49
1:B:562:LYS:O	1:B:566:ILE:HG12	2.13	0.49
1:A:237:PRO:HB3	1:A:302:MET:HB3	1.94	0.49
1:B:201:ASP:OD2	1:B:206:ASN:ND2	2.42	0.49
1:X:261:HIS:CD2	1:X:295:CYS:HB2	2.48	0.49
1:X:419:LEU:HD11	1:X:506:ILE:HA	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:425:LEU:HB3	1:B:429:ILE:HD12	1.94	0.48
1:B:11:VAL:HG12	1:B:15:LEU:HD12	1.95	0.48
1:A:429:ILE:HG13	1:A:541:TYR:CD1	2.49	0.48
1:B:464:LEU:O	1:B:469:VAL:HB	2.14	0.47
1:B:106:ILE:HA	1:B:110:THR:HG22	1.97	0.47
1:X:241:VAL:HG22	1:X:404:VAL:HG22	1.96	0.47
1:B:313:LEU:HB3	1:B:453:HIS:HD2	1.80	0.47
1:A:18:LYS:HE3	1:A:320:ILE:O	2.15	0.47
1:B:301:LEU:HD22	1:B:308:LEU:HD21	1.97	0.47
1:B:313:LEU:HB2	1:B:447:ASN:HB2	1.96	0.47
1:X:311:ASN:HB3	1:X:479:HIS:CB	2.45	0.47
1:B:585:VAL:O	1:B:592:THR:OG1	2.33	0.46
1:X:456:PHE:CD1	1:X:556:PRO:HB3	2.50	0.46
1:A:14:ASN:O	1:A:17:ARG:HG3	2.15	0.46
1:A:550:PRO:HB2	1:A:624:ILE:HG23	1.98	0.46
1:X:265:LEU:HD21	1:X:328:VAL:HG22	1.98	0.46
1:B:184:ASN:HB3	1:B:187:ARG:HB3	1.97	0.46
1:X:9:ASP:HB3	1:X:12:ILE:HG12	1.97	0.46
1:A:173:GLU:O	1:A:177:LYS:HG2	2.15	0.46
1:A:184:ASN:HB3	1:A:187:ARG:HB3	1.98	0.46
1:B:18:LYS:NZ	1:B:22:GLU:OE2	2.40	0.46
1:B:31:TYR:CZ	1:B:450:LYS:HA	2.51	0.46
1:B:34:GLN:O	1:B:37:GLN:HG3	2.16	0.45
1:B:525:CYS:SG	3:B:5002:BF8:O	2.75	0.45
1:B:6:LEU:HD11	1:B:16:TYR:CD2	2.52	0.45
1:B:32:GLU:HA	1:B:35:GLN:HG3	1.99	0.45
1:B:92:HIS:ND1	1:B:218:ASP:OD2	2.33	0.45
1:B:95:HIS:CD2	1:B:262:VAL:HG22	2.51	0.45
1:B:623:HIS:CE1	1:B:627:LYS:HE2	2.52	0.45
1:X:160:MET:HE1	1:X:221:ALA:HB2	1.97	0.45
1:B:350:VAL:HA	1:B:362:GLU:O	2.16	0.45
1:A:265:LEU:HD21	1:A:328:VAL:HG22	1.99	0.45
1:B:11:VAL:HG13	1:B:397:PRO:HG3	1.99	0.45
1:B:256:ILE:HB	1:B:290:LEU:HD23	1.99	0.45
1:B:311:ASN:HB3	1:B:479:HIS:CB	2.47	0.45
1:B:311:ASN:HB3	1:B:479:HIS:HB3	1.97	0.45
1:X:259:HIS:HB3	1:X:329:VAL:HG23	1.99	0.45
1:X:403:VAL:HG11	1:X:482:ALA:HB1	1.98	0.45
1:B:545:SER:OG	1:B:547:LYS:HG2	2.17	0.45
1:B:69:CYS:HB3	2:B:5000:SF4:S4	2.56	0.44
1:B:328:VAL:HA	1:B:350:VAL:HG23	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:219:GLY:HA3	1:B:569:TRP:CD1	2.52	0.44
1:B:458:ILE:O	1:B:462:LYS:HG3	2.17	0.44
1:B:578:HIS:HA	1:B:608:TYR:HB3	1.99	0.44
1:A:301:LEU:HD22	1:A:308:LEU:HD11	1.99	0.44
1:B:316:GLU:O	1:B:320:ILE:HG12	2.18	0.44
1:X:550:PRO:HB2	1:X:624:ILE:HG23	2.00	0.43
1:A:192:ALA:O	1:A:196:THR:HG23	2.18	0.43
1:A:223:LEU:HD11	1:A:566:ILE:HG12	1.99	0.43
1:B:438:LEU:HD12	1:B:628:ARG:HH11	1.82	0.43
1:X:424:PRO:O	1:X:427:PRO:HD2	2.18	0.43
1:A:35:GLN:NE2	4:A:5112:HOH:O	2.51	0.43
1:B:40:PHE:O	1:B:44:GLY:N	2.50	0.43
1:X:210:LYS:HD3	1:X:210:LYS:HA	1.81	0.43
1:A:40:PHE:CD1	1:A:56:ARG:HB2	2.54	0.43
1:B:525:CYS:HB2	3:B:5002:BF8:O	2.17	0.43
1:B:556:PRO:O	1:B:580:GLY:HA3	2.19	0.43
1:A:253:TYR:CE1	1:A:286:LYS:HD3	2.53	0.43
1:A:147:ILE:HG12	1:A:150:ARG:HH21	1.84	0.43
1:A:572:ALA:HA	1:A:603:ILE:HD13	1.99	0.43
1:B:298:ASN:HD21	1:B:479:HIS:HD2	1.65	0.43
1:X:106:ILE:HA	1:X:110:THR:HG22	2.01	0.43
1:B:579:LEU:HD11	1:B:607:PHE:HD2	1.84	0.43
1:X:208:ILE:HD11	1:X:591:VAL:HG13	2.00	0.43
1:A:294:CYS:HB3	1:A:295:CYS:H	1.69	0.43
1:A:103:LEU:HD22	1:A:229:LEU:HD22	1.99	0.43
1:B:18:LYS:HE3	1:B:320:ILE:O	2.19	0.43
1:B:201:ASP:HB3	1:B:207:LEU:HD21	2.01	0.43
1:X:41:GLY:HA3	1:A:54:PRO:HG3	2.00	0.42
1:B:579:LEU:HD11	1:B:607:PHE:CD2	2.53	0.42
1:B:503:LEU:HB3	1:B:516:LEU:HB2	2.01	0.42
1:B:451:VAL:HG11	1:B:456:PHE:HB2	2.01	0.42
1:A:528:ASN:HD22	1:A:566:ILE:HD13	1.85	0.42
1:B:542:LEU:HB2	1:B:544:VAL:HG22	2.01	0.42
1:B:242:THR:OG1	1:B:243:GLU:N	2.52	0.42
1:X:391:PRO:HA	1:X:394:VAL:HG13	2.01	0.42
1:A:259:HIS:HB3	1:A:329:VAL:HG23	2.02	0.42
1:X:239:PRO:HA	1:X:405:ALA:O	2.19	0.42
1:X:555:ALA:HB1	1:X:558:TYR:HB3	2.02	0.42
1:A:285:ALA:HB2	1:A:386:TYR:CE2	2.54	0.42
1:B:313:LEU:HD23	1:B:453:HIS:HB2	2.01	0.42
1:X:285:ALA:HB2	1:X:386:TYR:CE2	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:261:HIS:CD2	1:A:295:CYS:HB2	2.55	0.42
1:B:9:ASP:HA	1:B:10:PRO:HD3	1.92	0.42
1:B:429:ILE:HD13	1:B:541:TYR:CD1	2.54	0.41
1:X:560:HIS:CE1	1:X:562:LYS:HE2	2.55	0.41
1:B:260:GLY:HA3	1:B:330:ASP:OD1	2.20	0.41
1:B:456:PHE:CD1	1:B:556:PRO:HB3	2.55	0.41
1:A:58:ASP:OD2	1:A:62:GLU:HB2	2.19	0.41
1:B:301:LEU:HD22	1:B:308:LEU:HD11	2.02	0.41
1:B:448:ASN:H	1:B:557:GLU:HB3	1.85	0.41
1:X:92:HIS:ND1	1:X:218:ASP:OD1	2.37	0.41
1:A:346:HIS:CE1	1:A:389:ARG:HG3	2.55	0.41
1:B:24:ILE:HG12	1:B:341:ILE:HG12	2.02	0.41
1:B:525:CYS:CB	3:B:5002:BF8:O	2.69	0.41
1:X:95:HIS:CD2	1:X:95:HIS:C	2.99	0.41
1:X:562:LYS:HE3	3:X:702:BF8:S4	2.60	0.41
1:B:298:ASN:HD21	1:B:479:HIS:CD2	2.39	0.41
1:B:423:ASP:OD1	1:B:426:LYS:N	2.54	0.41
1:B:452:LYS:HB3	1:B:455:TRP:HB3	2.03	0.41
1:B:585:VAL:HG23	1:B:591:VAL:HG12	2.03	0.41
1:X:5:LYS:H	1:X:5:LYS:HG2	1.70	0.41
1:A:242:THR:OG1	1:A:243:GLU:N	2.53	0.41
1:B:503:LEU:H	1:B:503:LEU:HG	1.65	0.41
1:A:281:LYS:HA	1:A:285:ALA:O	2.20	0.41
1:X:16:TYR:CZ	1:X:20:ARG:HD2	2.57	0.40
1:A:560:HIS:CE1	1:A:562:LYS:HE2	2.56	0.40
1:X:52:GLN:HG3	1:A:559:GLN:HB2	2.03	0.40
1:A:329:VAL:HG13	1:A:334:ILE:HD13	2.02	0.40
1:A:239:PRO:HA	1:A:405:ALA:O	2.21	0.40
1:B:317:LEU:HD23	1:B:317:LEU:HA	1.89	0.40
1:A:184:ASN:O	1:A:188:GLU:HG2	2.21	0.40
1:B:35:GLN:C	1:B:37:GLN:H	2.30	0.40
1:X:268:LYS:HB3	1:X:375:SER:OG	2.22	0.40
1:B:28:PHE:HE2	1:B:453:HIS:HB2	1.86	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:34:GLN:NE2	1:B:56:ARG:O[5_675]	2.19	0.01

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	628/633 (99%)	605 (96%)	22 (4%)	1 (0%)	43	62
1	B	628/633 (99%)	597 (95%)	30 (5%)	1 (0%)	43	62
1	X	628/633 (99%)	604 (96%)	24 (4%)	0	100	100
All	All	1884/1899 (99%)	1806 (96%)	76 (4%)	2 (0%)	48	67

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	585	VAL
1	B	332	GLN

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	512/515 (99%)	501 (98%)	11 (2%)	47	72
1	B	512/515 (99%)	493 (96%)	19 (4%)	30	54
1	X	512/515 (99%)	502 (98%)	10 (2%)	48	73
All	All	1536/1545 (99%)	1496 (97%)	40 (3%)	40	66

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	X	5	LYS

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Mol	Chain	Res	Type
1	X	105	GLN
1	X	301	LEU
1	X	329	VAL
1	X	379	VAL
1	X	394	VAL
1	X	440	VAL
1	X	446	CYS
1	X	535	VAL
1	X	585	VAL
1	A	162	TYR
1	A	164	LYS
1	A	240	VAL
1	A	329	VAL
1	A	379	VAL
1	A	440	VAL
1	A	446	CYS
1	A	500	ARG
1	A	502	VAL
1	A	544	VAL
1	A	585	VAL
1	B	12	ILE
1	B	51	VAL
1	B	56	ARG
1	B	99	VAL
1	B	162	TYR
1	B	306	VAL
1	B	426	LYS
1	B	432	ILE
1	B	437	ILE
1	B	440	VAL
1	B	446	CYS
1	B	460	LEU
1	B	465	ILE
1	B	502	VAL
1	B	526	VAL
1	B	535	VAL
1	B	538	VAL
1	B	582	VAL
1	B	589	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	X	118	GLN
1	X	479	HIS
1	A	623	HIS
1	B	65	GLN
1	B	84	GLN
1	B	95	HIS
1	B	230	GLN

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 ⓘ

8 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	$\# Z > 2$
3	BF8	B	5002	1	0,14,14	-	-	-		
2	SF4	X	701	1	0,12,12	-	-	-		
2	SF4	A	5001	1	0,12,12	-	-	-		
2	SF4	B	5000	1	0,12,12	-	-	-		
2	SF4	A	5002	1	0,12,12	-	-	-		
2	SF4	B	5001	1	0,12,12	-	-	-		
3	BF8	X	702	1	0,14,14	-	-	-		
3	BF8	A	5003	1	0,14,14	-	-	-		

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
3	BF8	B	5002	1	-	-	0/4/5/5
2	SF4	X	701	1	-	-	0/6/5/5
2	SF4	A	5001	1	-	-	0/6/5/5
2	SF4	B	5000	1	-	-	0/6/5/5
2	SF4	A	5002	1	-	-	0/6/5/5
2	SF4	B	5001	1	-	-	0/6/5/5
3	BF8	X	702	1	-	-	0/4/5/5
3	BF8	A	5003	1	-	-	0/4/5/5

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	5002	BF8	4	0
2	B	5000	SF4	1	0
2	B	5001	SF4	1	0
3	X	702	BF8	1	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	630/633 (99%)	0.17	5 (0%) 82 80	38, 54, 73, 88	0
1	B	630/633 (99%)	1.15	128 (20%) 3 2	43, 69, 95, 106	0
1	X	630/633 (99%)	-0.01	2 (0%) 90 88	36, 46, 60, 75	0
All	All	1890/1899 (99%)	0.44	135 (7%) 22 19	36, 54, 86, 106	0

All (135) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	40	PHE	4.7
1	B	317	LEU	4.6
1	B	451	VAL	4.6
1	B	69	CYS	4.4
1	B	450	LYS	4.2
1	B	585	VAL	4.0
1	B	586	VAL	4.0
1	B	12	ILE	3.9
1	B	582	VAL	3.9
1	B	26	THR	3.8
1	B	57	ILE	3.8
1	B	59	PRO	3.8
1	B	587	GLY	3.8
1	B	27	VAL	3.7
1	B	611	THR	3.6
1	B	8	VAL	3.6
1	B	54	PRO	3.6
1	B	44	GLY	3.5
1	B	464	LEU	3.5
1	B	609	VAL	3.4
1	B	460	LEU	3.4
1	B	41	GLY	3.3
1	B	607	PHE	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	309	ALA	3.3
1	B	624	ILE	3.3
1	B	45	LEU	3.2
1	B	613	PRO	3.2
1	B	39	GLY	3.2
1	B	448	ASN	3.2
1	B	43	THR	3.2
1	B	73	ALA	3.1
1	B	36	PRO	3.1
1	B	42	LEU	3.1
1	B	485	GLY	3.1
1	B	494	TRP	3.1
1	B	455	TRP	3.1
1	B	584	PRO	3.0
1	B	16	TYR	3.0
1	B	15	LEU	2.9
1	B	486	LEU	2.9
1	B	456	PHE	2.9
1	B	608	TYR	2.9
1	B	614	TYR	2.9
1	B	550	PRO	2.9
1	B	60	PHE	2.9
1	B	48	ARG	2.8
1	B	461	VAL	2.8
1	B	518	PRO	2.8
1	B	438	LEU	2.8
1	B	512	LEU	2.8
1	B	5	LYS	2.8
1	B	63	GLY	2.8
1	B	445	GLY	2.8
1	B	33	ALA	2.8
1	B	471	VAL	2.7
1	B	64	PRO	2.7
1	B	14	ASN	2.7
1	B	520	LEU	2.7
1	B	52	GLN	2.7
1	B	91	ALA	2.6
1	B	76	ILE	2.6
1	B	437	ILE	2.6
1	B	583	PRO	2.6
1	B	55	CYS	2.6
1	B	318	ALA	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	555	ALA	2.6
1	X	99	VAL	2.6
1	B	68	ILE	2.6
1	A	525	CYS	2.6
1	B	591	VAL	2.5
1	A	499	LEU	2.5
1	B	49	HIS	2.5
1	B	10	PRO	2.5
1	B	29	ASP	2.5
1	B	432	ILE	2.5
1	B	469	VAL	2.5
1	B	453	HIS	2.5
1	B	480	ALA	2.5
1	B	506	ILE	2.4
1	B	633	LEU	2.4
1	B	30	ARG	2.4
1	B	75	VAL	2.4
1	B	490	ALA	2.4
1	B	556	PRO	2.4
1	A	514	GLY	2.4
1	B	595	LEU	2.4
1	B	11	VAL	2.4
1	B	80	ASN	2.4
1	B	627	LYS	2.4
1	B	291	ALA	2.4
1	B	482	ALA	2.4
1	B	38	CYS	2.3
1	B	440	VAL	2.3
1	B	46	CYS	2.3
1	B	558	TYR	2.3
1	B	568	THR	2.3
1	B	9	ASP	2.3
1	B	316	GLU	2.3
1	B	70	GLY	2.2
1	B	619	GLY	2.2
1	B	24	ILE	2.2
1	B	34	GLN	2.2
1	A	222	GLY	2.2
1	B	580	GLY	2.2
1	B	551	ILE	2.2
1	B	454	ASP	2.2
1	B	56	ARG	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	433	VAL	2.2
1	B	538	VAL	2.2
1	B	6	LEU	2.2
1	B	246	LEU	2.2
1	B	481	LEU	2.2
1	B	492	ALA	2.2
1	A	134	ASN	2.1
1	B	549	LEU	2.1
1	B	396	ILE	2.1
1	B	525	CYS	2.1
1	B	4	SER	2.1
1	B	552	ALA	2.1
1	B	617	ALA	2.1
1	B	419	LEU	2.1
1	B	579	LEU	2.1
1	B	62	GLU	2.1
1	B	487	MET	2.1
1	B	502	VAL	2.1
1	B	53	GLY	2.1
1	B	341	ILE	2.1
1	B	623	HIS	2.1
1	B	581	VAL	2.1
1	B	428	LEU	2.1
1	B	321	THR	2.1
1	X	398	ARG	2.0
1	B	391	PRO	2.0
1	B	495	ALA	2.0
1	B	7	SER	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 [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. 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	SF4	B	5001	8/8	0.91	0.10	77,83,104,104	8
3	BF8	B	5002	10/10	0.95	0.07	60,66,72,73	6
3	BF8	A	5003	10/10	0.96	0.06	46,50,59,63	7
2	SF4	B	5000	8/8	0.96	0.06	64,67,71,76	0
3	BF8	X	702	10/10	0.97	0.05	45,47,55,59	8
2	SF4	A	5001	8/8	0.98	0.04	42,44,51,55	0
2	SF4	A	5002	8/8	0.99	0.03	34,41,45,48	0
2	SF4	X	701	8/8	0.99	0.04	34,45,50,50	0

6.5 Other polymers ⓘ

There are no such residues in this entry.