



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 09:45 AM UTC

PDB ID : 3S3K / pdb_00003s3k
Title : Crystal structure of the catalytic domain of PTP10D from *Drosophila melanogaster* with a small molecular inhibitor para-NitroCatechol Sulphate
Authors : Madan, L.L.; Gopal, B.
Deposited on : 2011-05-18
Resolution : 3.20 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

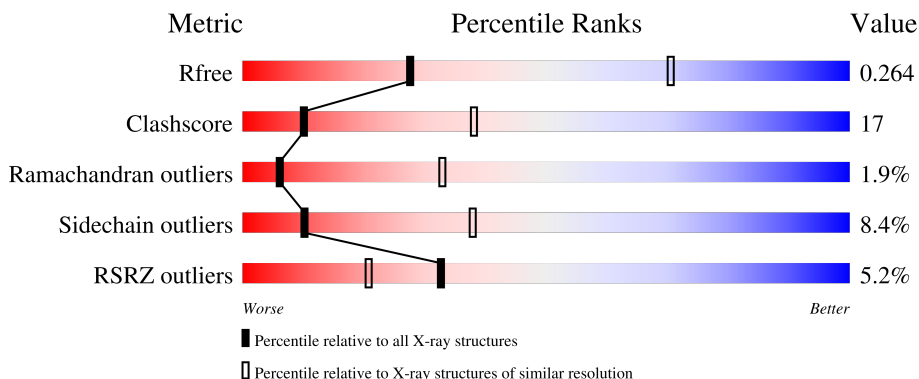
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	307	
1	B	307	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4704 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 Tyrosine-protein phosphatase 10D.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	287	Total	C	N	O	S	0	0	0
			2345	1490	420	418	17			
1	B	285	Total	C	N	O	S	0	0	0
			2318	1469	414	418	17			

There are 46 discrepancies between the modelled and reference sequences:

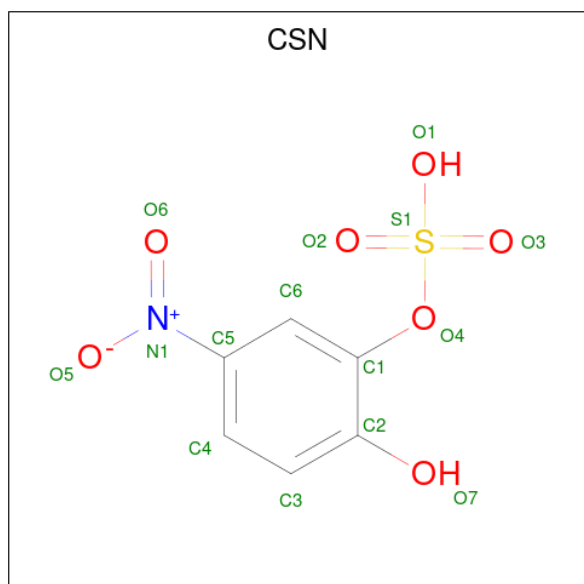
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP P35992
A	2	GLY	-	expression tag	UNP P35992
A	3	SER	-	expression tag	UNP P35992
A	4	SER	-	expression tag	UNP P35992
A	5	HIS	-	expression tag	UNP P35992
A	6	HIS	-	expression tag	UNP P35992
A	7	HIS	-	expression tag	UNP P35992
A	8	HIS	-	expression tag	UNP P35992
A	9	HIS	-	expression tag	UNP P35992
A	10	HIS	-	expression tag	UNP P35992
A	11	SER	-	expression tag	UNP P35992
A	12	SER	-	expression tag	UNP P35992
A	13	GLY	-	expression tag	UNP P35992
A	14	LEU	-	expression tag	UNP P35992
A	15	VAL	-	expression tag	UNP P35992
A	16	PRO	-	expression tag	UNP P35992
A	17	ARG	-	expression tag	UNP P35992
A	18	GLY	-	expression tag	UNP P35992
A	19	SER	-	expression tag	UNP P35992
A	20	HIS	-	expression tag	UNP P35992
A	21	MET	-	expression tag	UNP P35992
A	22	ALA	-	expression tag	UNP P35992
A	23	SER	-	expression tag	UNP P35992
B	1	MET	-	expression tag	UNP P35992
B	2	GLY	-	expression tag	UNP P35992

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Chain	Residue	Modelled	Actual	Comment	Reference
B	3	SER	-	expression tag	UNP P35992
B	4	SER	-	expression tag	UNP P35992
B	5	HIS	-	expression tag	UNP P35992
B	6	HIS	-	expression tag	UNP P35992
B	7	HIS	-	expression tag	UNP P35992
B	8	HIS	-	expression tag	UNP P35992
B	9	HIS	-	expression tag	UNP P35992
B	10	HIS	-	expression tag	UNP P35992
B	11	SER	-	expression tag	UNP P35992
B	12	SER	-	expression tag	UNP P35992
B	13	GLY	-	expression tag	UNP P35992
B	14	LEU	-	expression tag	UNP P35992
B	15	VAL	-	expression tag	UNP P35992
B	16	PRO	-	expression tag	UNP P35992
B	17	ARG	-	expression tag	UNP P35992
B	18	GLY	-	expression tag	UNP P35992
B	19	SER	-	expression tag	UNP P35992
B	20	HIS	-	expression tag	UNP P35992
B	21	MET	-	expression tag	UNP P35992
B	22	ALA	-	expression tag	UNP P35992
B	23	SER	-	expression tag	UNP P35992

- Molecule 2 is N,4-DIHYDROXY-N-OXO-3-(SULFOOXY)BENZENAMINIUM (CCD ID: CSN) (formula: $C_6H_5NO_7S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	S	0	0
			15	6	1	7	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	7	Total	O	0	0
			7	7		
3	B	19	Total	O	0	0
			19	19		

4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	102.51Å 102.51Å 172.80Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	44.08 – 3.20 44.08 – 3.20	Depositor EDS
% Data completeness (in resolution range)	96.2 (44.08-3.20) 96.2 (44.08-3.20)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.37 (at 3.19Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.253 , 0.309 0.222 , 0.264	Depositor DCC
R_{free} test set	880 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	39.1	Xtriage
Anisotropy	0.027	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 55.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.047 for -h,-k,l	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	4704	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.56% 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: CSN

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	0.81	6/2412 (0.2%)	0.94	4/3272 (0.1%)
1	B	0.90	7/2383 (0.3%)	1.00	9/3236 (0.3%)
All	All	0.85	13/4795 (0.3%)	0.97	13/6508 (0.2%)

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	22	ALA	CA-CB	-6.90	1.43	1.53
1	B	244	ALA	CA-CB	-6.64	1.42	1.53
1	A	283	TRP	C-O	-6.13	1.16	1.24
1	A	243	SER	C-O	-6.10	1.17	1.24
1	A	281	ARG	C-O	-6.04	1.16	1.23
1	B	242	CYS	CB-SG	-5.82	1.62	1.81
1	A	156	ASP	C-O	-5.76	1.16	1.23
1	B	175	ILE	CA-CB	-5.59	1.47	1.54
1	A	20	HIS	C-O	-5.48	1.18	1.24
1	B	176	LEU	C-O	-5.10	1.16	1.24
1	B	243	SER	C-O	-5.10	1.17	1.23
1	B	242	CYS	C-O	-5.05	1.17	1.24
1	B	243	SER	CA-C	-5.04	1.49	1.52

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	210	ASP	N-CA-C	-10.44	101.04	113.88
1	B	34	HIS	N-CA-C	-7.26	102.36	112.45
1	A	96	GLU	N-CA-C	-6.35	99.35	109.76
1	A	104	ASN	N-CA-C	6.05	119.27	109.40
1	B	35	TYR	N-CA-C	-5.92	104.52	110.97
1	B	36	ARG	N-CA-C	5.91	118.20	111.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	284	MET	N-CA-C	5.82	117.32	110.97
1	B	215	ASN	CA-C-N	-5.78	114.42	120.38
1	B	215	ASN	C-N-CA	-5.78	114.42	120.38
1	B	235	GLN	N-CA-C	5.58	117.52	108.26
1	A	105	TYR	N-CA-C	-5.26	102.06	109.96
1	B	164	VAL	CA-C-N	5.22	125.15	119.78
1	B	164	VAL	C-N-CA	5.22	125.15	119.78

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	2345	0	2225	66	0
1	B	2318	0	2195	92	0
2	B	15	0	5	2	0
3	A	7	0	0	0	0
3	B	19	0	0	0	0
All	All	4704	0	4425	157	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 17.

All (157) 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:29:LYS:C	1:B:30:ASN:HD22	1.46	1.22
1:B:304:GLY:O	1:B:305:LYS:HB2	1.41	1.12
1:B:176:LEU:O	1:B:177:ASN:HB2	1.51	1.04
1:A:95:ASP:HB3	1:A:98:SER:HB3	1.39	1.04
1:B:28:ILE:HD11	1:B:262:ILE:HG22	1.45	0.96
1:B:29:LYS:C	1:B:30:ASN:ND2	2.30	0.89
1:B:30:ASN:HD22	1:B:30:ASN:N	1.72	0.83
1:B:244:ALA:HB2	2:B:308:CSN:H6	1.61	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:179:SER:HB2	1:B:186:MET:HB3	1.67	0.77
1:B:235:GLN:O	1:B:235:GLN:HG3	1.89	0.71
1:B:176:LEU:O	1:B:177:ASN:CB	2.30	0.69
1:A:186:MET:HE1	1:A:230:ARG:HG2	1.76	0.68
1:B:215:ASN:ND2	1:B:216:PRO:HD3	2.09	0.68
1:B:304:GLY:O	1:B:305:LYS:CB	2.29	0.67
1:A:282:VAL:O	1:A:283:TRP:C	2.36	0.67
1:B:233:ALA:HA	1:B:234:GLU:HG2	1.76	0.66
1:B:35:TYR:OH	1:B:296:GLN:HG2	1.96	0.66
1:A:85:SER:C	1:A:104:ASN:HD22	2.04	0.66
1:A:199:ILE:HG22	1:A:200:LEU:N	2.11	0.66
1:B:70:ASN:HA	1:B:73:LYS:HD2	1.77	0.65
1:B:299:LEU:O	1:B:303:GLU:HG3	1.95	0.64
1:B:29:LYS:O	1:B:30:ASN:ND2	2.30	0.64
1:B:241:HIS:ND1	1:B:241:HIS:C	2.56	0.64
1:A:246:VAL:O	1:A:284:MET:O	2.17	0.63
1:A:269:ASP:O	1:A:273:ILE:HG12	1.99	0.63
1:A:228:ARG:NH2	1:A:256:ASP:OD2	2.33	0.62
1:A:246:VAL:HG23	1:A:247:GLY:H	1.65	0.61
1:A:54:LYS:HG3	1:A:55:HIS:CD2	2.35	0.61
1:A:70:ASN:HA	1:A:73:LYS:HD2	1.83	0.60
1:A:256:ASP:O	1:A:260:GLN:HG3	2.02	0.60
1:B:28:ILE:HD11	1:B:262:ILE:CG2	2.28	0.60
1:B:233:ALA:HA	1:B:235:GLN:H	1.67	0.60
1:B:233:ALA:CB	1:B:235:GLN:H	2.15	0.60
1:A:253:ILE:HG22	1:A:277:MET:HE1	1.84	0.59
1:B:277:MET:HB3	1:B:284:MET:HG3	1.84	0.59
1:A:253:ILE:HG22	1:A:277:MET:CE	2.33	0.59
1:B:278:ARG:NH2	1:B:285:VAL:O	2.35	0.59
1:A:90:GLN:O	1:A:131:ARG:NH2	2.37	0.58
1:A:246:VAL:HG23	1:A:247:GLY:N	2.19	0.58
1:B:41:ASP:O	1:B:42:SER:C	2.46	0.58
1:A:208:TRP:HB3	1:A:248:ARG:NH2	2.19	0.57
1:B:30:ASN:ND2	1:B:30:ASN:N	2.46	0.57
1:A:95:ASP:CB	1:A:98:SER:HB3	2.24	0.57
1:B:233:ALA:CA	1:B:235:GLN:H	2.18	0.57
1:B:216:PRO:O	1:B:218:GLN:N	2.37	0.56
1:B:71:ARG:N	1:B:72:PRO:HD2	2.21	0.56
1:A:199:ILE:CG2	1:A:200:LEU:N	2.69	0.56
1:A:208:TRP:CD2	1:A:248:ARG:HD3	2.41	0.55
1:B:107:PRO:HD3	1:B:280:GLU:HG3	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:259:LEU:HD21	1:A:301:VAL:HG21	1.89	0.55
1:B:95:ASP:C	1:B:95:ASP:OD1	2.50	0.55
1:B:244:ALA:CB	2:B:308:CSN:H6	2.35	0.55
1:A:273:ILE:O	1:A:277:MET:HG3	2.08	0.54
1:B:109:HIS:HB2	1:B:257:ARG:HG3	1.88	0.54
1:B:92:VAL:H	1:B:98:SER:HB2	1.73	0.54
1:B:207:THR:O	1:B:208:TRP:C	2.51	0.54
1:A:85:SER:HB3	1:A:104:ASN:HD22	1.71	0.54
1:A:103:ALA:HB2	1:A:118:THR:HB	1.90	0.53
1:A:282:VAL:HG22	1:A:283:TRP:N	2.24	0.53
1:B:81:PRO:HA	1:B:104:ASN:HD21	1.74	0.53
1:B:118:THR:O	1:B:241:HIS:HB2	2.09	0.53
1:A:85:SER:CB	1:A:104:ASN:HD22	2.22	0.53
1:B:122:LEU:HB2	1:B:125:THR:HG23	1.91	0.53
1:B:247:GLY:N	1:B:284:MET:O	2.32	0.53
1:A:89:LEU:HB3	1:A:131:ARG:NH2	2.24	0.52
1:A:90:GLN:HB2	1:A:135:GLU:OE2	2.10	0.52
1:B:103:ALA:HB2	1:B:118:THR:HB	1.93	0.51
1:A:261:GLN:C	1:A:263:ASN:H	2.19	0.51
1:A:220:LEU:O	1:A:224:VAL:HG23	2.11	0.51
1:B:104:ASN:HD22	1:B:281:ARG:HD2	1.75	0.50
1:A:33:GLU:OE1	1:A:36:ARG:NH1	2.44	0.50
1:B:181:TYR:OH	1:B:230:ARG:NH2	2.45	0.50
1:A:34:HIS:O	1:A:38:MET:HG2	2.11	0.50
1:A:26:ILE:HG12	1:A:34:HIS:CD2	2.47	0.50
1:A:253:ILE:CG2	1:A:277:MET:CE	2.90	0.49
1:B:104:ASN:HD22	1:B:281:ARG:HH11	1.59	0.49
1:B:208:TRP:C	1:B:209:PRO:O	2.54	0.49
1:B:208:TRP:O	1:B:209:PRO:O	2.30	0.49
1:B:21:MET:CG	1:B:22:ALA:H	2.26	0.49
1:B:154:LYS:O	1:B:243:SER:OG	2.27	0.49
1:B:250:GLY:CA	1:B:284:MET:HE3	2.43	0.49
1:B:26:ILE:C	1:B:26:ILE:HD13	2.38	0.49
1:A:246:VAL:HB	1:A:283:TRP:O	2.13	0.49
1:B:69:CYS:SG	1:B:96:GLU:HB2	2.53	0.49
1:A:71:ARG:N	1:A:72:PRO:HD2	2.27	0.49
1:B:247:GLY:HA2	1:B:286:GLN:H	1.78	0.49
1:A:79:ILE:HD12	1:A:246:VAL:HG11	1.95	0.48
1:A:134:TRP:CZ2	1:A:193:ARG:HB3	2.49	0.48
1:B:38:MET:HB3	1:B:46:PHE:CE2	2.49	0.48
1:B:275:TYR:C	1:B:275:TYR:CD2	2.92	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:21:MET:HG3	1:B:22:ALA:H	1.79	0.47
1:B:233:ALA:HA	1:B:235:GLN:N	2.28	0.47
1:A:253:ILE:CG2	1:A:277:MET:HE1	2.44	0.47
1:B:106:VAL:CG2	1:B:117:VAL:HG13	2.44	0.47
1:B:171:ILE:HD13	1:B:193:ARG:HD2	1.95	0.47
1:B:261:GLN:O	1:B:263:ASN:N	2.48	0.47
1:A:24:ARG:HH22	1:A:52:GLU:CD	2.23	0.47
1:A:47:SER:HA	1:A:288:GLU:OE1	2.15	0.47
1:B:65:ALA:HB2	1:B:81:PRO:HD2	1.96	0.47
1:B:259:LEU:HD21	1:B:301:VAL:HG21	1.97	0.47
1:B:60:GLN:HB2	1:B:82:TYR:CD1	2.50	0.46
1:B:123:HIS:CE1	1:B:126:ARG:NH2	2.84	0.46
1:B:213:VAL:HG21	1:B:293:CYS:HB2	1.97	0.46
1:A:28:ILE:HG12	1:A:268:VAL:HG23	1.98	0.46
1:A:144:LEU:HD11	1:A:252:PHE:CB	2.46	0.45
1:A:281:ARG:HG3	1:A:282:VAL:N	2.31	0.45
1:B:104:ASN:ND2	1:B:281:ARG:HH11	2.14	0.45
1:B:89:LEU:HD22	1:B:131:ARG:HG2	1.99	0.45
1:B:250:GLY:HA2	1:B:284:MET:HE3	1.98	0.45
1:B:126:ARG:HD3	1:B:156:ASP:OD2	2.16	0.45
1:A:85:SER:CB	1:A:104:ASN:ND2	2.80	0.45
1:B:84:HIS:CD2	1:B:84:HIS:C	2.95	0.44
1:B:233:ALA:HA	1:B:234:GLU:CG	2.43	0.44
1:A:111:SER:HB3	1:A:114:GLU:HB2	1.99	0.44
1:B:178:ASP:HB2	1:B:180:HIS:CE1	2.52	0.44
1:B:202:HIS:CE1	1:B:204:HIS:HB2	2.53	0.44
1:A:38:MET:HE3	1:A:46:PHE:HA	2.00	0.44
1:A:85:SER:HB3	1:A:104:ASN:ND2	2.33	0.44
1:B:130:TRP:HB3	1:B:171:ILE:HG21	1.98	0.44
1:A:282:VAL:O	1:A:284:MET:N	2.50	0.44
1:B:71:ARG:N	1:B:72:PRO:CD	2.81	0.44
1:B:134:TRP:CZ2	1:B:193:ARG:HB3	2.53	0.44
1:A:217:PRO:HB2	1:A:297:CYS:SG	2.58	0.43
1:A:231:ILE:HA	1:A:231:ILE:HD12	1.80	0.43
1:B:63:THR:H	1:B:83:ASP:CG	2.26	0.43
1:B:217:PRO:HG2	1:B:296:GLN:OE1	2.18	0.43
1:B:256:ASP:O	1:B:260:GLN:HG3	2.19	0.43
1:A:208:TRP:CD1	1:A:209:PRO:HD2	2.54	0.43
1:A:27:LEU:HD23	1:A:267:TYR:CZ	2.54	0.43
1:A:77:THR:HG23	1:B:68:PRO:HG3	2.00	0.43
1:B:92:VAL:H	1:B:98:SER:CB	2.30	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:110:ASN:N	1:B:114:GLU:OE1	2.48	0.43
1:B:208:TRP:CZ2	1:B:248:ARG:HG2	2.54	0.43
1:A:115:PHE:CE2	1:A:253:ILE:HG23	2.54	0.42
1:A:64:PHE:C	1:A:66:ASP:H	2.27	0.42
1:A:143:MET:HG3	1:A:241:HIS:CE1	2.53	0.42
1:B:178:ASP:OD2	1:B:178:ASP:N	2.52	0.42
1:A:148:PHE:CE1	1:A:153:GLU:HB2	2.54	0.42
1:A:109:HIS:HB2	1:A:257:ARG:HG3	2.00	0.42
1:B:273:ILE:O	1:B:277:MET:HG3	2.19	0.42
1:A:28:ILE:HD12	1:A:266:ASP:HA	2.01	0.42
1:B:143:MET:HE2	1:B:204:HIS:NE2	2.34	0.42
1:B:277:MET:HE2	1:B:284:MET:HE2	2.01	0.42
1:A:183:ASP:HB3	1:A:207:THR:HB	2.01	0.42
1:A:70:ASN:HB3	1:A:100:TYR:CD2	2.55	0.41
1:B:305:LYS:HD3	1:B:305:LYS:HA	1.59	0.41
1:A:145:THR:OG1	1:A:146:ARG:N	2.53	0.41
1:A:228:ARG:HH21	1:A:256:ASP:CG	2.27	0.41
1:A:99:ASP:OD1	1:A:99:ASP:C	2.62	0.41
1:B:26:ILE:HD12	1:B:31:PHE:CD1	2.55	0.41
1:B:233:ALA:HA	1:B:234:GLU:CB	2.50	0.41
1:B:140:ALA:HA	1:B:201:ARG:O	2.21	0.41
1:B:234:GLU:HG2	1:B:235:GLN:N	2.36	0.41
1:A:224:VAL:HG21	1:A:255:LEU:HD23	2.02	0.41
1:B:84:HIS:CD2	1:B:84:HIS:O	2.74	0.40
1:B:26:ILE:HD12	1:B:31:PHE:HD1	1.86	0.40
1:B:62:CYS:HA	1:B:83:ASP:OD1	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	285/307 (93%)	258 (90%)	25 (9%)	2 (1%)	18	52
1	B	283/307 (92%)	248 (88%)	26 (9%)	9 (3%)	3	21
All	All	568/614 (92%)	506 (89%)	51 (9%)	11 (2%)	6	33

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	40	ALA
1	B	177	ASN
1	A	283	TRP
1	B	216	PRO
1	B	218	GLN
1	B	262	ILE
1	A	262	ILE
1	B	209	PRO
1	B	217	PRO
1	B	214	PRO
1	B	215	ASN

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	251/278 (90%)	233 (93%)	18 (7%)	13	44
1	B	250/278 (90%)	226 (90%)	24 (10%)	8	32
All	All	501/556 (90%)	459 (92%)	42 (8%)	10	38

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	HIS
1	A	21	MET
1	A	33	GLU
1	A	46	PHE
1	A	66	ASP

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Mol	Chain	Res	Type
1	A	90	GLN
1	A	95	ASP
1	A	117	VAL
1	A	131	ARG
1	A	146	ARG
1	A	207	THR
1	A	231	ILE
1	A	238	ILE
1	A	241	HIS
1	A	243	SER
1	A	266	ASP
1	A	280	GLU
1	A	305	LYS
1	B	21	MET
1	B	26	ILE
1	B	28	ILE
1	B	30	ASN
1	B	37	LEU
1	B	46	PHE
1	B	85	SER
1	B	117	VAL
1	B	157	GLN
1	B	178	ASP
1	B	207	THR
1	B	209	PRO
1	B	215	ASN
1	B	219	THR
1	B	229	ASP
1	B	234	GLU
1	B	235	GLN
1	B	241	HIS
1	B	243	SER
1	B	246	VAL
1	B	263	ASN
1	B	265	SER
1	B	268	VAL
1	B	305	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	55	HIS

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Mol	Chain	Res	Type
1	A	104	ASN
1	A	109	HIS
1	A	174	GLN
1	A	263	ASN
1	B	30	ASN
1	B	84	HIS
1	B	104	ASN
1	B	177	ASN
1	B	180	HIS
1	B	197	GLN
1	B	215	ASN
1	B	218	GLN
1	B	261	GLN
1	B	290	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is 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
2	CSN	B	308	-	15,15,15	4.82	6 (40%)	18,22,22	2.12	4 (22%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	CSN	B	308	-	-	3/7/9/9	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	308	CSN	O6-N1	11.56	1.42	1.22
2	B	308	CSN	C6-C1	8.74	1.53	1.38
2	B	308	CSN	C6-C5	8.60	1.53	1.39
2	B	308	CSN	O4-S1	-5.48	1.47	1.58
2	B	308	CSN	C5-N1	-4.12	1.35	1.45
2	B	308	CSN	O4-C1	-2.55	1.37	1.42

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	308	CSN	C5-C6-C1	-6.94	111.55	119.06
2	B	308	CSN	C6-C1-C2	4.13	124.21	120.04
2	B	308	CSN	O6-N1-C5	2.12	121.73	118.82
2	B	308	CSN	C3-C4-C5	2.12	122.88	120.08

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	308	CSN	C1-O4-S1-O2
2	B	308	CSN	C1-O4-S1-O3
2	B	308	CSN	C1-O4-S1-O1

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	308	CSN	2	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	287/307 (93%)	0.38	9 (3%) 51 32	27, 53, 81, 104	0
1	B	285/307 (92%)	0.53	21 (7%) 20 13	27, 52, 87, 120	0
All	All	572/614 (93%)	0.46	30 (5%) 33 21	27, 52, 83, 120	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	41	ASP	4.6
1	B	39	SER	3.7
1	A	93	ASP	3.5
1	B	304	GLY	3.5
1	B	233	ALA	3.3
1	B	305	LYS	3.1
1	B	211	PHE	3.0
1	A	304	GLY	3.0
1	B	44	PHE	3.0
1	A	30	ASN	2.9
1	B	234	GLU	2.7
1	B	59	ASP	2.7
1	B	37	LEU	2.7
1	B	215	ASN	2.6
1	B	33	GLU	2.6
1	A	94	ASP	2.6
1	A	305	LYS	2.6
1	B	40	ALA	2.5
1	B	235	GLN	2.5
1	A	68	PRO	2.5
1	B	162	ASP	2.5
1	B	77	THR	2.4
1	A	63	THR	2.3
1	B	96	GLU	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	42	SER	2.3
1	A	124	SER	2.2
1	B	30	ASN	2.2
1	B	217	PRO	2.2
1	A	92	VAL	2.1
1	B	151	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 [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	CSN	B	308	15/15	0.90	0.29	38,38,38,38	0

6.5 Other polymers [i](#)

There are no such residues in this entry.