



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 16, 2017 – 12:38 AM EDT

PDB ID : 3DEI
Title : Crystal Structures of Caspase-3 with Bound Isoquinoline-1,3,4-trione Derivative Inhibitors
Authors : Wu, J.; Du, J.; Li, J.; Ding, J.
Deposited on : unknown
Resolution : 2.80 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20030345
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20030345

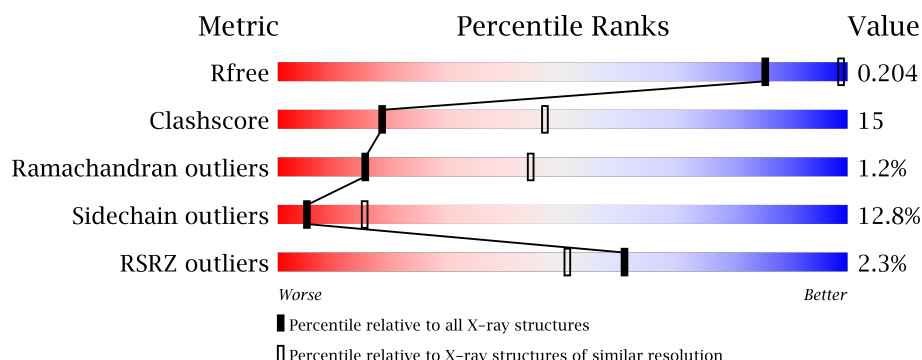
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	100719	2583 (2.80-2.80)
Clashscore	112137	3033 (2.80-2.80)
Ramachandran outliers	110173	2983 (2.80-2.80)
Sidechain outliers	110143	2985 (2.80-2.80)
RSRZ outliers	101464	2610 (2.80-2.80)

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	249	<div> <div>0.2%</div> <div>60% 28% 5% 7%</div> </div>
1	B	249	<div> <div>0.2%</div> <div>63% 26% 5% 6%</div> </div>
1	C	249	<div> <div>2%</div> <div>55% 33% 5% 6%</div> </div>
1	D	249	<div> <div>5%</div> <div>50% 37% 6% 6%</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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	RXB	C	300	-	-	X	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7672 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.

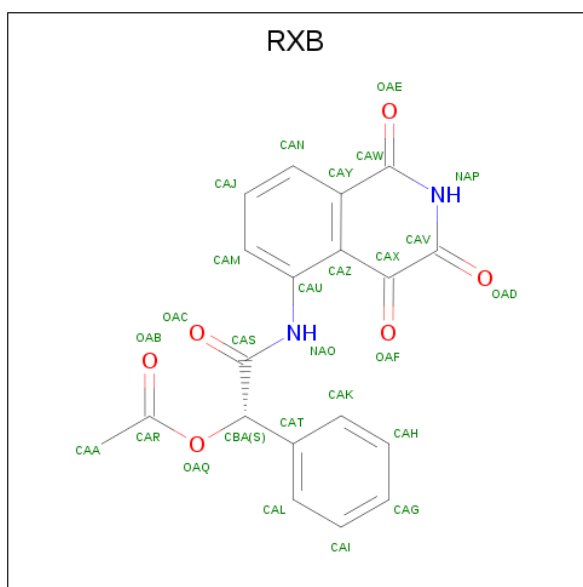
- Molecule 1 is a protein called Caspase-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	232	Total	C	N	O	S	0	0	0
			1883	1192	323	354	14			
1	B	233	Total	C	N	O	S	0	0	0
			1891	1198	324	355	14			
1	C	234	Total	C	N	O	S	0	0	0
			1901	1204	327	356	14			
1	D	234	Total	C	N	O	S	0	0	0
			1896	1201	325	356	14			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	190	ASP	GLU	SEE REMARK 999	UNP P42574
B	190	ASP	GLU	SEE REMARK 999	UNP P42574
C	190	ASP	GLU	SEE REMARK 999	UNP P42574
D	190	ASP	GLU	SEE REMARK 999	UNP P42574

- Molecule 2 is (1S)-2-oxo-1-phenyl-2-[(1,3,4-trioxo-1,2,3,4-tetrahydroisoquinolin-5-yl)amino]ethyl acetate (three-letter code: RXB) (formula: C₁₉H₁₄N₂O₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	1	Total	C	N	O	0	0
			27	19	2	6		

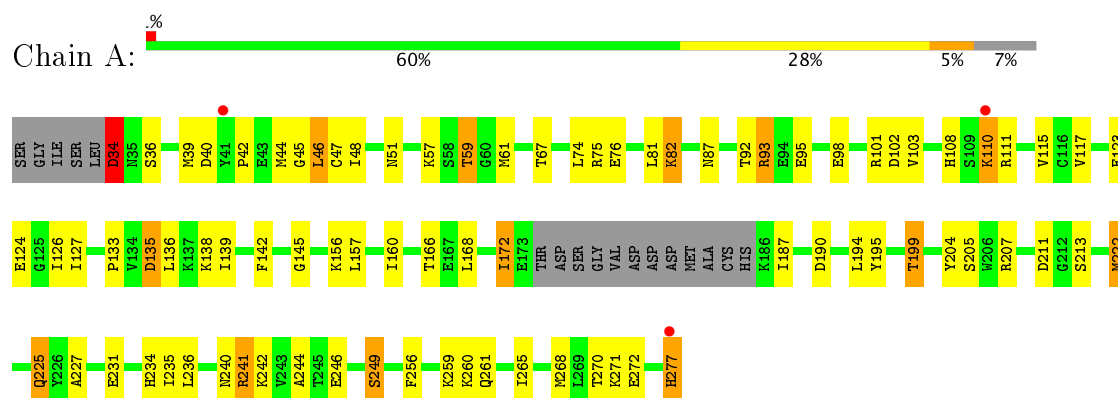
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	29	Total O 29 29	0	0
3	B	24	Total O 24 24	0	0
3	C	18	Total O 18 18	0	0
3	D	3	Total O 3 3	0	0

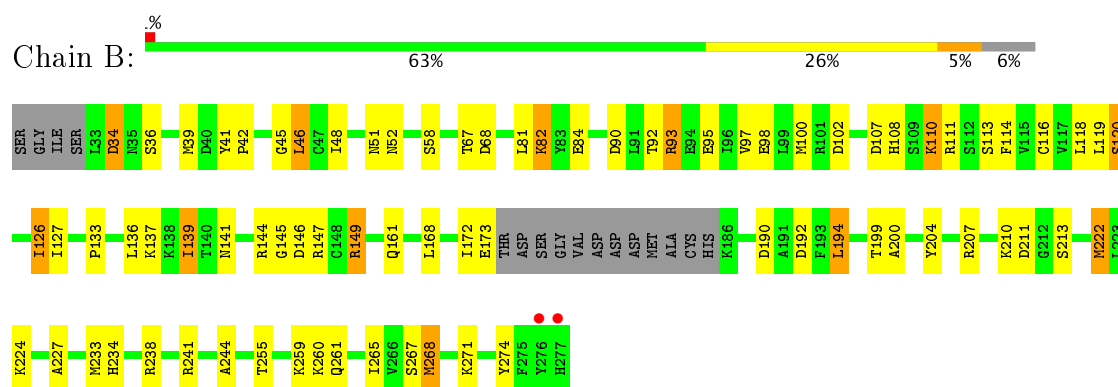
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

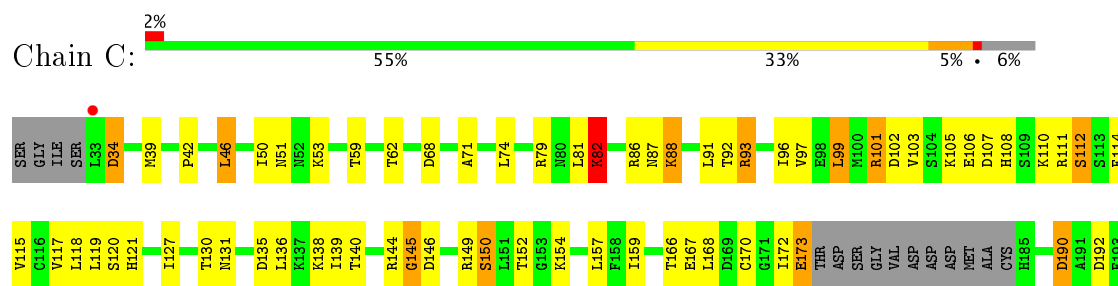
• Molecule 1: Caspase-3

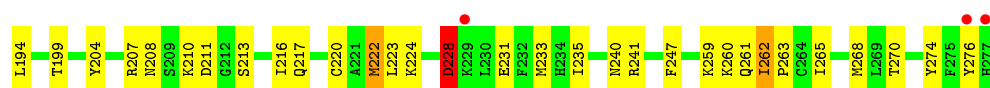


• Molecule 1: Caspase-3

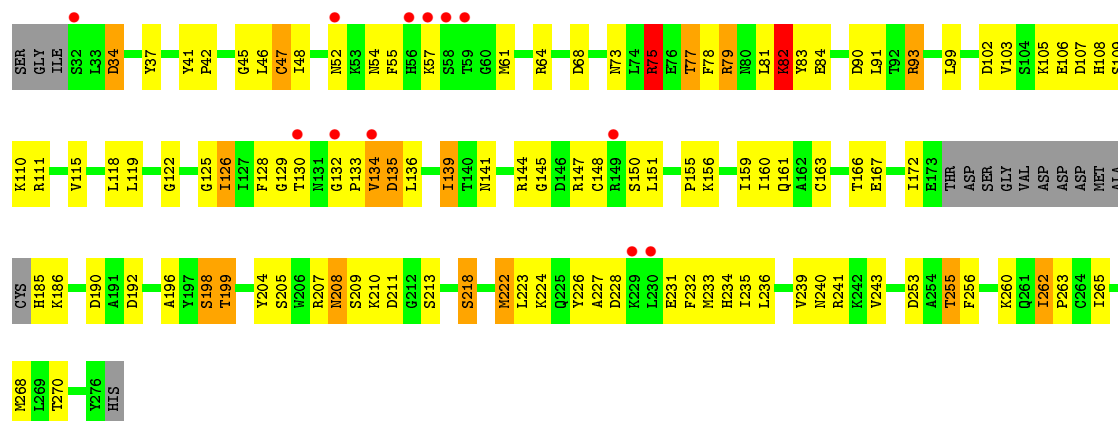


• Molecule 1: Caspase-3





● Molecule 1: Caspase-3



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	65.52Å 95.98Å 180.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.13 – 2.80 45.13 – 2.80	Depositor EDS
% Data completeness (in resolution range)	94.1 (45.13-2.80) 93.8 (45.13-2.80)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.94 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.211 , 0.291 0.211 , 0.204	Depositor DCC
R_{free} test set	1383 reflections (5.37%)	DCC
Wilson B-factor (Å ²)	68.4	Xtriage
Anisotropy	0.092	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 56.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7672	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.78% 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: OCS, RXB

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	0.63	0/1912	0.87	6/2566 (0.2%)
1	B	0.62	0/1920	0.86	9/2577 (0.3%)
1	C	0.55	0/1931	0.80	6/2592 (0.2%)
1	D	0.51	0/1925	0.79	8/2585 (0.3%)
All	All	0.58	0/7688	0.83	29/10320 (0.3%)

There are no bond length outliers.

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	102	ASP	CB-CG-OD2	7.47	125.03	118.30
1	B	190	ASP	CB-CG-OD2	7.06	124.66	118.30
1	A	190	ASP	CB-CG-OD2	6.73	124.36	118.30
1	B	90	ASP	CB-CG-OD2	6.69	124.32	118.30
1	A	34	ASP	CB-CG-OD2	6.33	124.00	118.30
1	C	190	ASP	CB-CG-OD2	6.21	123.88	118.30
1	A	135	ASP	CB-CG-OD2	6.07	123.76	118.30
1	C	135	ASP	CB-CG-OD2	6.07	123.76	118.30
1	A	40	ASP	CB-CG-OD2	6.00	123.70	118.30
1	B	68	ASP	CB-CG-OD2	6.00	123.70	118.30
1	B	102	ASP	CB-CG-OD2	5.96	123.66	118.30
1	D	102	ASP	CB-CG-OD2	5.93	123.64	118.30
1	A	211	ASP	CB-CG-OD2	5.87	123.58	118.30
1	D	34	ASP	CB-CG-OD2	5.85	123.57	118.30
1	B	34	ASP	CB-CG-OD2	5.75	123.47	118.30
1	C	102	ASP	CB-CG-OD2	5.74	123.47	118.30
1	C	107	ASP	CB-CG-OD2	5.74	123.47	118.30
1	D	135	ASP	CB-CG-OD2	5.68	123.41	118.30
1	D	190	ASP	CB-CG-OD2	5.60	123.34	118.30
1	D	107	ASP	CB-CG-OD2	5.52	123.27	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	228	ASP	CB-CG-OD2	5.33	123.10	118.30
1	D	253	ASP	CB-CG-OD2	5.32	123.09	118.30
1	B	107	ASP	CB-CG-OD2	5.30	123.07	118.30
1	B	211	ASP	CB-CG-OD2	5.29	123.06	118.30
1	B	146	ASP	CB-CG-OD2	5.26	123.03	118.30
1	B	192	ASP	CB-CG-OD2	5.25	123.03	118.30
1	C	34	ASP	CB-CG-OD2	5.23	123.01	118.30
1	D	68	ASP	CB-CG-OD2	5.12	122.91	118.30
1	D	211	ASP	CB-CG-OD2	5.09	122.88	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1883	0	1852	59	0
1	B	1891	0	1863	47	0
1	C	1901	0	1870	69	0
1	D	1896	0	1868	55	0
2	C	27	0	14	9	0
3	A	29	0	0	1	0
3	B	24	0	0	1	0
3	C	18	0	0	0	0
3	D	3	0	0	1	0
All	All	7672	0	7467	226	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 15.

All (226) 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:46:LEU:H	1:A:46:LEU:HD12	1.04	1.09
2:C:300:RXB:OAF	2:C:300:RXB:HAK	1.65	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:262:ILE:HD13	1:D:263:PRO:HD2	1.47	0.96
1:A:93:ARG:HH21	1:A:133:PRO:HD2	1.31	0.95
1:A:46:LEU:CD1	1:A:46:LEU:H	1.81	0.93
1:A:51:ASN:HD21	1:A:67:THR:HG23	1.35	0.89
1:D:208:ASN:HD22	1:D:209:SER:H	1.20	0.86
1:A:46:LEU:N	1:A:46:LEU:HD12	1.89	0.86
1:D:208:ASN:HD22	1:D:209:SER:N	1.75	0.85
1:C:99:LEU:O	1:C:103:VAL:HG23	1.78	0.84
1:C:222:MET:HA	1:C:222:MET:HE2	1.60	0.82
1:A:225:GLN:HG3	1:A:225:GLN:O	1.82	0.79
1:C:222:MET:CE	1:C:222:MET:HA	2.13	0.78
1:D:122:GLY:O	1:D:163:OCS:HB3	1.83	0.78
1:D:45:GLY:HA2	1:D:111:ARG:HD3	1.63	0.78
1:C:42:PRO:HG2	1:C:110:LYS:HB3	1.66	0.77
1:C:46:LEU:HD11	1:C:108:HIS:HE1	1.50	0.77
1:A:51:ASN:HD21	1:A:67:THR:CG2	1.98	0.76
1:A:101:ARG:HG3	1:A:142:PHE:CE1	2.23	0.73
1:D:55:PHE:HE1	1:D:64:ARG:HG3	1.51	0.73
1:A:199:THR:HG21	1:A:204:TYR:O	1.88	0.73
1:B:52:ASN:HB2	1:B:120:SER:HB2	1.71	0.73
1:C:208:ASN:HD22	1:C:211:ASP:H	1.37	0.73
1:B:51:ASN:ND2	1:B:67:THR:OG1	2.22	0.72
1:C:199:THR:HG21	1:C:204:TYR:O	1.90	0.72
1:C:51:ASN:HD22	1:C:71:ALA:HB2	1.54	0.71
1:A:222:MET:CE	1:A:222:MET:HA	2.21	0.70
1:D:119:LEU:HD23	1:D:161:GLN:HB3	1.73	0.70
1:D:78:PHE:O	1:D:83:TYR:HB2	1.91	0.70
1:B:92:THR:HG23	1:B:95:GLU:OE1	1.92	0.69
1:A:207:ARG:HG2	1:A:213:SER:HB3	1.75	0.68
1:C:208:ASN:ND2	1:C:211:ASP:H	1.92	0.68
1:C:88:LYS:HG2	1:C:91:LEU:HD11	1.77	0.66
1:C:92:THR:HB	1:C:131:ASN:ND2	2.10	0.65
1:C:51:ASN:ND2	1:C:71:ALA:HB2	2.12	0.65
1:A:222:MET:HE2	1:A:222:MET:HA	1.79	0.65
1:C:101:ARG:HH11	1:C:101:ARG:HB3	1.61	0.64
1:C:170:CYS:O	1:D:186:LYS:NZ	2.31	0.64
1:C:51:ASN:HD22	1:C:119:LEU:HD12	1.64	0.62
1:B:168:LEU:HB3	1:B:259:LYS:HD2	1.81	0.62
1:B:108:HIS:HA	1:B:111:ARG:HD2	1.81	0.62
1:A:74:LEU:HD13	1:A:117:VAL:HG11	1.82	0.61
1:A:34:ASP:HA	3:A:288:HOH:O	2.01	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:92:THR:OG1	1:A:95:GLU:HG3	2.01	0.61
1:B:46:LEU:HD12	1:B:46:LEU:H	1.66	0.60
1:C:51:ASN:ND2	1:C:119:LEU:HD12	2.17	0.60
1:A:241:ARG:HH11	1:A:241:ARG:CG	2.15	0.59
1:C:168:LEU:HD13	2:C:300:RXB:HNAO	1.67	0.59
1:B:92:THR:OG1	1:B:95:GLU:HG3	2.01	0.59
1:D:126:ILE:HG22	1:D:135:ASP:HA	1.85	0.58
1:D:148:CYS:SG	1:D:151:LEU:HB2	2.44	0.58
1:C:241:ARG:HH11	1:C:241:ARG:HG3	1.69	0.58
1:C:74:LEU:CD1	1:C:117:VAL:HG11	2.34	0.57
1:D:262:ILE:HD13	1:D:263:PRO:CD	2.28	0.57
1:C:199:THR:HG22	1:C:261:GLN:HG3	1.87	0.57
1:A:101:ARG:HG3	1:A:142:PHE:HE1	1.69	0.57
1:A:241:ARG:HH11	1:A:241:ARG:HG3	1.70	0.56
1:A:231:GLU:HG3	1:A:272:GLU:HB3	1.88	0.56
1:B:207:ARG:HA	1:B:213:SER:HA	1.88	0.56
2:C:300:RXB:OAF	2:C:300:RXB:CAK	2.48	0.56
1:C:53:LYS:HE3	1:C:68:ASP:OD2	2.05	0.56
1:B:194:LEU:HD12	1:B:267:SER:HB2	1.87	0.55
1:C:81:LEU:O	1:C:82:LYS:HB2	2.05	0.55
1:A:45:GLY:HA2	1:A:111:ARG:HG2	1.89	0.55
1:C:247:PHE:O	1:C:260:LYS:NZ	2.32	0.55
2:C:300:RXB:CAL	2:C:300:RXB:HAAA	2.36	0.55
1:C:118:LEU:HD13	1:C:127:ILE:HD13	1.88	0.55
1:D:55:PHE:CE1	1:D:64:ARG:HG3	2.39	0.55
1:D:141:ASN:HB3	1:D:147:ARG:HH12	1.72	0.54
1:D:75:ARG:O	1:D:79:ARG:HB2	2.07	0.54
1:C:50:ILE:HG21	1:C:96:ILE:HD11	1.90	0.54
1:D:37:TYR:CD1	1:D:155:PRO:HD3	2.42	0.53
1:B:100:MET:HG3	1:B:139:ILE:HB	1.91	0.53
1:A:234:HIS:CE1	1:B:234:HIS:CE1	2.96	0.53
1:B:81:LEU:O	1:B:82:LYS:HB2	2.09	0.53
2:C:300:RXB:OAF	2:C:300:RXB:OAC	2.26	0.52
1:D:199:THR:HG21	1:D:204:TYR:O	2.08	0.52
1:B:233:MET:HE1	1:B:265:ILE:HG21	1.91	0.52
1:C:166:THR:O	2:C:300:RXB:HAM	2.09	0.52
1:C:262:ILE:HD13	1:C:263:PRO:HD2	1.91	0.52
1:C:207:ARG:HA	1:C:213:SER:HA	1.92	0.52
1:D:73:ASN:O	1:D:77:THR:OG1	2.28	0.51
1:C:213:SER:OG	1:C:216:ILE:HB	2.09	0.51
1:A:135:ASP:O	1:A:138:LYS:HB3	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:199:THR:HG23	1:A:261:GLN:HG3	1.91	0.51
1:B:36:SER:HA	1:B:271:LYS:HB3	1.91	0.51
1:B:222:MET:HA	1:B:222:MET:HE2	1.91	0.51
1:A:271:LYS:HE2	1:B:241:ARG:NH1	2.26	0.51
1:C:119:LEU:CD2	1:C:216:ILE:HD11	2.41	0.51
1:A:204:TYR:CG	1:A:205:SER:N	2.80	0.50
1:A:39:MET:O	1:A:277:HIS:HE1	1.95	0.50
1:A:75:ARG:NH1	1:A:87:ASN:OD1	2.40	0.50
2:C:300:RXB:OAF	2:C:300:RXB:CAS	2.59	0.50
1:A:46:LEU:HD21	1:A:108:HIS:HE1	1.77	0.50
1:A:136:LEU:HD11	1:A:160:ILE:HD13	1.93	0.50
1:B:46:LEU:CD1	1:B:114:PHE:HD1	2.25	0.49
1:D:129:GLY:N	1:D:132:GLY:O	2.36	0.49
1:A:241:ARG:NH2	1:B:34:ASP:HB3	2.27	0.49
1:A:242:LYS:NZ	1:A:246:GLU:OE1	2.44	0.49
1:B:268:MET:HE2	3:B:284:HOH:O	2.12	0.49
1:B:141:ASN:HB3	1:B:147:ARG:NH1	2.28	0.49
1:D:159:ILE:HG13	1:D:196:ALA:HB3	1.93	0.49
1:C:119:LEU:HD21	1:C:216:ILE:HD11	1.95	0.49
1:C:127:ILE:HG23	1:C:136:LEU:HG	1.95	0.49
1:D:240:ASN:HA	1:D:263:PRO:HG2	1.95	0.49
1:C:213:SER:O	1:C:217:GLN:HB2	2.12	0.49
1:A:207:ARG:HA	1:A:213:SER:HA	1.94	0.49
1:B:93:ARG:NH2	1:B:133:PRO:HD2	2.28	0.49
1:C:127:ILE:HG13	1:C:127:ILE:O	2.12	0.49
1:C:233:MET:HE1	1:C:265:ILE:HG21	1.94	0.49
1:C:59:THR:HG21	1:C:130:THR:O	2.13	0.48
1:C:228:ASP:HA	1:C:276:TYR:CD2	2.49	0.48
1:C:50:ILE:HG21	1:C:96:ILE:CD1	2.44	0.48
1:C:93:ARG:O	1:C:97:VAL:HG23	2.13	0.48
1:C:115:VAL:HG22	1:C:157:LEU:HD12	1.96	0.47
1:B:39:MET:HG3	1:B:274:TYR:O	2.14	0.47
1:C:144:ARG:O	1:C:146:ASP:N	2.47	0.47
1:C:74:LEU:HD13	1:C:117:VAL:HG11	1.97	0.47
1:D:207:ARG:HA	1:D:213:SER:HA	1.96	0.47
1:B:233:MET:HE1	1:B:265:ILE:CG2	2.45	0.47
1:D:235:ILE:O	1:D:239:VAL:HG23	2.14	0.47
1:D:148:CYS:SG	1:D:151:LEU:HD12	2.55	0.47
1:D:55:PHE:HD2	1:D:130:THR:HA	1.79	0.47
1:C:274:TYR:HD1	1:C:276:TYR:HH	1.62	0.47
1:A:207:ARG:HB2	1:A:207:ARG:NH1	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:128:PHE:CE1	1:D:133:PRO:HB3	2.49	0.46
1:A:46:LEU:HD11	1:A:108:HIS:CE1	2.51	0.46
1:C:233:MET:HE1	1:C:265:ILE:CG2	2.45	0.46
1:A:44:MET:HG3	1:A:81:LEU:O	2.16	0.46
1:B:126:ILE:HD13	1:B:133:PRO:CB	2.46	0.46
1:C:168:LEU:HB3	1:C:259:LYS:HD2	1.98	0.46
1:A:39:MET:O	1:A:277:HIS:CE1	2.68	0.46
1:B:199:THR:HG22	1:B:200:ALA:N	2.30	0.46
1:B:41:TYR:HB3	1:B:110:LYS:O	2.16	0.46
1:C:39:MET:HA	1:C:112:SER:HB3	1.98	0.46
1:A:168:LEU:HG	2:C:300:RXB:HNAP	1.80	0.46
1:B:119:LEU:HD23	1:B:161:GLN:HB3	1.98	0.45
1:C:222:MET:CA	1:C:222:MET:CE	2.91	0.45
1:B:222:MET:CE	1:B:222:MET:HA	2.46	0.45
1:B:244:ALA:O	1:B:260:LYS:NZ	2.37	0.45
1:A:48:ILE:HD13	1:A:103:VAL:HG11	1.99	0.45
1:C:150:SER:O	1:C:154:LYS:NZ	2.50	0.45
1:D:42:PRO:HG2	1:D:110:LYS:HD3	1.98	0.45
1:B:199:THR:HG21	1:B:204:TYR:O	2.16	0.45
1:B:48:ILE:HG21	1:B:100:MET:CE	2.47	0.45
1:B:46:LEU:HD12	1:B:46:LEU:N	2.28	0.44
1:C:86:ARG:NH2	1:C:106:GLU:OE2	2.42	0.44
1:A:123:GLU:O	1:A:124:GLU:C	2.53	0.44
1:C:223:LEU:HD23	1:C:235:ILE:HG21	1.98	0.44
1:D:93:ARG:HB2	1:D:134:VAL:HG23	1.99	0.44
1:D:93:ARG:HE	1:D:132:GLY:HA3	1.83	0.44
1:D:231:GLU:OE1	1:D:233:MET:HB2	2.18	0.44
1:A:46:LEU:HD11	1:A:108:HIS:ND1	2.33	0.44
1:D:141:ASN:OD1	1:D:144:ARG:HD3	2.17	0.44
1:A:249:SER:HB3	1:A:259:LYS:H	1.82	0.44
1:C:108:HIS:HA	1:C:111:ARG:HD2	1.98	0.44
1:D:125:GLY:HA2	3:D:279:HOH:O	2.18	0.44
1:D:232:PHE:HA	1:D:235:ILE:HD12	2.00	0.44
1:A:194:LEU:C	1:A:194:LEU:HD23	2.39	0.43
1:C:210:LYS:HB3	1:C:210:LYS:HE2	1.78	0.43
2:C:300:RXB:OAC	2:C:300:RXB:CAX	2.66	0.43
1:A:195:TYR:N	1:A:195:TYR:CD1	2.86	0.43
1:A:270:THR:O	1:B:241:ARG:HB2	2.18	0.43
1:A:82:LYS:HD3	1:A:82:LYS:N	2.34	0.43
1:D:54:ASN:N	1:D:90:ASP:OD2	2.45	0.43
1:A:172:ILE:HD11	1:B:149:ARG:CZ	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:59:THR:C	1:A:61:MET:H	2.22	0.43
1:D:198:SER:O	1:D:263:PRO:HA	2.18	0.43
1:A:235:ILE:H	1:A:235:ILE:HG12	1.67	0.43
1:C:231:GLU:O	1:C:235:ILE:HG12	2.19	0.43
1:D:236:LEU:HB3	1:D:265:ILE:CD1	2.49	0.43
1:B:222:MET:HG2	1:B:238:ARG:HB2	2.01	0.43
1:A:241:ARG:CG	1:A:241:ARG:NH1	2.81	0.43
1:C:101:ARG:HH11	1:C:101:ARG:CB	2.30	0.43
1:B:255:THR:O	1:B:259:LYS:HE2	2.19	0.43
1:B:81:LEU:HD21	1:B:227:ALA:HB1	2.00	0.42
1:D:99:LEU:O	1:D:103:VAL:HG23	2.19	0.42
1:B:81:LEU:HD21	1:B:227:ALA:CB	2.49	0.42
1:C:173:GLU:HA	1:D:185:HIS:O	2.19	0.42
1:D:47:CYS:HA	1:D:115:VAL:O	2.19	0.42
1:D:135:ASP:O	1:D:136:LEU:C	2.58	0.42
1:D:192:ASP:OD1	1:D:270:THR:OG1	2.35	0.42
1:C:103:VAL:O	1:C:106:GLU:HB2	2.18	0.42
1:D:108:HIS:HA	1:D:111:ARG:HH11	1.84	0.42
1:A:42:PRO:HG2	1:A:110:LYS:HB3	2.01	0.42
1:C:92:THR:HB	1:C:131:ASN:HD21	1.81	0.42
1:A:81:LEU:O	1:A:82:LYS:HB2	2.20	0.42
1:C:117:VAL:HG22	1:C:159:ILE:HG13	2.02	0.42
1:C:74:LEU:HD23	1:C:220:CYS:SG	2.60	0.42
1:A:240:ASN:O	1:A:244:ALA:CB	2.67	0.42
1:B:233:MET:CE	1:B:265:ILE:HG21	2.48	0.42
1:D:47:CYS:SG	1:D:47:CYS:O	2.78	0.42
1:A:236:LEU:HB3	1:A:265:ILE:HD12	2.01	0.41
1:D:218:SER:O	1:D:222:MET:HB2	2.20	0.41
1:D:52:ASN:ND2	1:D:91:LEU:O	2.53	0.41
1:B:137:LYS:HG3	1:B:141:ASN:HD21	1.84	0.41
1:D:239:VAL:O	1:D:243:VAL:HG23	2.20	0.41
1:B:48:ILE:HG21	1:B:100:MET:HE1	2.02	0.41
1:B:41:TYR:O	1:B:42:PRO:C	2.58	0.41
1:B:45:GLY:HA3	1:B:113:SER:O	2.20	0.41
1:C:270:THR:O	1:D:241:ARG:HB2	2.19	0.41
1:A:249:SER:OG	1:A:256:PHE:O	2.38	0.41
1:B:116:CYS:SG	1:B:118:LEU:HD21	2.60	0.41
1:C:118:LEU:HD22	1:C:127:ILE:CD1	2.51	0.41
1:C:228:ASP:HA	1:C:276:TYR:HD2	1.86	0.41
1:B:126:ILE:HD13	1:B:133:PRO:HB3	2.03	0.41
1:C:120:SER:OG	1:C:121:HIS:N	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:231:GLU:CG	1:D:234:HIS:HE2	2.34	0.41
1:D:255:THR:HB	1:D:256:PHE:CD1	2.56	0.41
1:A:156:LYS:HE2	1:A:156:LYS:HB2	1.82	0.41
1:C:145:GLY:O	1:C:152:THR:OG1	2.39	0.41
1:D:118:LEU:HD11	1:D:139:ILE:HD11	2.02	0.41
1:C:199:THR:HG22	1:C:261:GLN:CG	2.49	0.41
1:D:41:TYR:HB3	1:D:42:PRO:HD2	2.03	0.41
1:A:231:GLU:O	1:A:235:ILE:HG12	2.21	0.40
1:A:47:CYS:HA	1:A:115:VAL:O	2.21	0.40
1:C:46:LEU:HD12	1:C:114:PHE:HD1	1.86	0.40
1:D:223:LEU:HD23	1:D:223:LEU:HA	2.00	0.40
1:A:277:HIS:CD2	1:A:277:HIS:N	2.88	0.40
1:B:93:ARG:O	1:B:97:VAL:HG23	2.21	0.40
1:C:240:ASN:HA	1:C:263:PRO:HD2	2.03	0.40
1:D:226:TYR:O	1:D:228:ASP:N	2.55	0.40
1:D:166:THR:O	1:D:167:GLU:C	2.59	0.40
1:A:241:ARG:HH21	1:B:34:ASP:HB3	1.84	0.40
1:D:81:LEU:O	1:D:82:LYS:CB	2.70	0.40
1:C:51:ASN:HD22	1:C:119:LEU:CD1	2.34	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	227/249 (91%)	212 (93%)	13 (6%)	2 (1%)	20	52
1	B	228/249 (92%)	209 (92%)	18 (8%)	1 (0%)	38	72
1	C	229/249 (92%)	219 (96%)	8 (4%)	2 (1%)	20	52
1	D	229/249 (92%)	199 (87%)	24 (10%)	6 (3%)	6	21
All	All	913/996 (92%)	839 (92%)	63 (7%)	11 (1%)	15	44

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	145	GLY
1	D	34	ASP
1	D	82	LYS
1	D	145	GLY
1	B	145	GLY
1	D	75	ARG
1	D	106	GLU
1	A	227	ALA
1	C	82	LYS
1	D	227	ALA
1	A	145	GLY

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	207/221 (94%)	182 (88%)	25 (12%)	6	17
1	B	208/221 (94%)	186 (89%)	22 (11%)	8	23
1	C	209/221 (95%)	181 (87%)	28 (13%)	4	13
1	D	209/221 (95%)	177 (85%)	32 (15%)	3	9
All	All	833/884 (94%)	726 (87%)	107 (13%)	5	15

All (107) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	34	ASP
1	A	36	SER
1	A	46	LEU
1	A	57	LYS
1	A	59	THR
1	A	76	GLU
1	A	82	LYS
1	A	93	ARG
1	A	98	GLU

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Mol	Chain	Res	Type
1	A	110	LYS
1	A	126	ILE
1	A	127	ILE
1	A	139	ILE
1	A	157	LEU
1	A	166	THR
1	A	172	ILE
1	A	187	ILE
1	A	199	THR
1	A	222	MET
1	A	225	GLN
1	A	241	ARG
1	A	249	SER
1	A	260	LYS
1	A	268	MET
1	A	277	HIS
1	B	46	LEU
1	B	58	SER
1	B	82	LYS
1	B	84	GLU
1	B	93	ARG
1	B	98	GLU
1	B	110	LYS
1	B	120	SER
1	B	126	ILE
1	B	127	ILE
1	B	136	LEU
1	B	139	ILE
1	B	144	ARG
1	B	149	ARG
1	B	172	ILE
1	B	173	GLU
1	B	194	LEU
1	B	210	LYS
1	B	222	MET
1	B	224	LYS
1	B	261	GLN
1	B	268	MET
1	C	34	ASP
1	C	46	LEU
1	C	62	THR
1	C	79	ARG

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Mol	Chain	Res	Type
1	C	82	LYS
1	C	87	ASN
1	C	88	LYS
1	C	93	ARG
1	C	99	LEU
1	C	101	ARG
1	C	105	LYS
1	C	112	SER
1	C	138	LYS
1	C	139	ILE
1	C	140	THR
1	C	149	ARG
1	C	150	SER
1	C	167	GLU
1	C	172	ILE
1	C	173	GLU
1	C	190	ASP
1	C	192	ASP
1	C	194	LEU
1	C	222	MET
1	C	224	LYS
1	C	228	ASP
1	C	262	ILE
1	C	268	MET
1	D	46	LEU
1	D	47	CYS
1	D	48	ILE
1	D	57	LYS
1	D	61	MET
1	D	75	ARG
1	D	77	THR
1	D	79	ARG
1	D	82	LYS
1	D	84	GLU
1	D	93	ARG
1	D	105	LYS
1	D	109	SER
1	D	126	ILE
1	D	134	VAL
1	D	139	ILE
1	D	150	SER
1	D	156	LYS

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Mol	Chain	Res	Type
1	D	160	ILE
1	D	172	ILE
1	D	198	SER
1	D	199	THR
1	D	205	SER
1	D	208	ASN
1	D	210	LYS
1	D	218	SER
1	D	222	MET
1	D	224	LYS
1	D	255	THR
1	D	260	LYS
1	D	262	ILE
1	D	268	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	ASN
1	A	51	ASN
1	A	56	HIS
1	A	217	GLN
1	A	277	HIS
1	B	51	ASN
1	B	217	GLN
1	C	35	ASN
1	C	51	ASN
1	C	208	ASN
1	D	51	ASN
1	D	56	HIS
1	D	80	ASN
1	D	89	ASN
1	D	208	ASN
1	D	225	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues 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
1	OCS	A	163	1	8,8,9	1.52	3 (37%)	7,11,13	5.30	4 (57%)
1	OCS	B	163	1	8,8,9	1.24	2 (25%)	7,11,13	3.84	3 (42%)
1	OCS	C	163	1	8,8,9	1.49	1 (12%)	7,11,13	5.99	6 (85%)
1	OCS	D	163	1	8,8,9	1.24	1 (12%)	7,11,13	2.68	3 (42%)

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
1	OCS	A	163	1	-	0/4/7/9	0/0/0/0
1	OCS	B	163	1	-	0/4/7/9	0/0/0/0
1	OCS	C	163	1	-	0/4/7/9	0/0/0/0
1	OCS	D	163	1	-	0/4/7/9	0/0/0/0

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	163	OCS	CB-SG	-2.66	1.66	1.77
1	A	163	OCS	CB-SG	-2.26	1.68	1.77
1	B	163	OCS	CB-SG	-2.07	1.69	1.77
1	A	163	OCS	CA-C	2.10	1.53	1.50
1	B	163	OCS	OD3-SG	2.29	1.51	1.45
1	A	163	OCS	OD3-SG	2.44	1.52	1.45
1	C	163	OCS	OD3-SG	2.84	1.53	1.45

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	163	OCS	OD2-SG-CB	-5.74	99.00	106.01
1	A	163	OCS	OD2-SG-OD1	-3.37	103.64	111.37
1	C	163	OCS	OD3-SG-OD1	-2.89	103.83	113.86
1	A	163	OCS	OD3-SG-OD1	-2.72	104.44	113.86
1	C	163	OCS	O-C-CA	-2.25	118.82	125.02
1	C	163	OCS	OD2-SG-OD1	-2.21	106.30	111.37
1	B	163	OCS	OD2-SG-OD1	-2.20	106.33	111.37
1	D	163	OCS	OD2-SG-CB	2.09	108.56	106.01
1	A	163	OCS	OD2-SG-OD3	2.39	116.84	111.37
1	C	163	OCS	OD2-SG-OD3	3.85	120.19	111.37
1	D	163	OCS	OD1-SG-CB	4.10	110.33	106.83
1	D	163	OCS	OD3-SG-CB	4.71	110.85	106.83
1	B	163	OCS	OD2-SG-CB	4.82	111.89	106.01
1	B	163	OCS	OD3-SG-CB	8.33	113.94	106.83
1	A	163	OCS	OD3-SG-CB	12.98	117.92	106.83
1	C	163	OCS	OD3-SG-CB	13.47	118.34	106.83

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	D	163	OCS	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates 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	RXB	C	300	-	29,29,29	3.56	5 (17%)	37,41,41	3.16	15 (40%)

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	RXB	C	300	-	-	1/16/32/32	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	300	RXB	CAV-CAX	-16.19	1.38	1.54
2	C	300	RXB	CAY-CAW	-4.24	1.40	1.47
2	C	300	RXB	OAF-CAX	2.16	1.28	1.23
2	C	300	RXB	OAE-CAW	3.05	1.29	1.23
2	C	300	RXB	OAQ-CAR	6.84	1.50	1.35

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	300	RXB	OAC-CAS-CBA	-5.58	113.21	120.33
2	C	300	RXB	OAQ-CAR-OAB	-3.23	116.39	122.94
2	C	300	RXB	CAN-CAY-CAW	-2.74	116.33	120.33
2	C	300	RXB	CAT-CBA-CAS	-2.73	104.54	110.31
2	C	300	RXB	CAU-NAO-CAS	-2.71	119.74	126.92
2	C	300	RXB	OAD-CAV-CAX	-2.34	117.24	119.61
2	C	300	RXB	CAN-CAY-CAZ	2.02	123.15	120.17
2	C	300	RXB	CAX-CAV-NAP	2.03	120.78	118.00
2	C	300	RXB	CAJ-CAM-CAU	2.10	123.03	118.61
2	C	300	RXB	CAY-CAW-NAP	2.18	117.04	114.84
2	C	300	RXB	OAF-CAX-CAZ	2.40	126.58	121.94
2	C	300	RXB	CAZ-CAU-NAO	3.36	123.78	118.80
2	C	300	RXB	OAQ-CAR-CAA	7.84	125.82	111.10
2	C	300	RXB	CBA-OAQ-CAR	8.28	126.39	115.48
2	C	300	RXB	OAQ-CBA-CAT	10.69	123.88	107.74

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	300	RXB	CAR-OAQ-CBA-CAT

There are no ring outliers.

1 monomer is involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	300	RXB	9	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	231/249 (92%)	-0.16	3 (1%) 77 71	36, 59, 89, 125	0
1	B	232/249 (93%)	-0.20	2 (0%) 84 79	33, 58, 89, 119	0
1	C	233/249 (93%)	0.08	4 (1%) 70 63	39, 72, 106, 161	0
1	D	233/249 (93%)	0.24	12 (5%) 28 19	54, 90, 125, 171	0
All	All	929/996 (93%)	-0.01	21 (2%) 61 51	33, 68, 113, 171	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	33	LEU	4.6
1	A	277	HIS	3.5
1	A	110	LYS	3.4
1	C	277	HIS	3.3
1	C	276	TYR	2.8
1	D	57	LYS	2.7
1	D	149	ARG	2.7
1	D	230	LEU	2.7
1	D	32	SER	2.6
1	D	130	THR	2.5
1	A	41	TYR	2.5
1	D	58	SER	2.4
1	D	132	GLY	2.4
1	D	59	THR	2.4
1	D	56	HIS	2.2
1	B	277	HIS	2.2
1	B	276	TYR	2.2
1	D	134	VAL	2.2
1	C	229	LYS	2.2
1	D	229	LYS	2.1
1	D	52	ASN	2.1

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

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
1	OCS	C	163	9/10	0.92	0.22	-	49,51,58,61	0
1	OCS	D	163	9/10	0.95	0.16	-	95,97,98,99	0
1	OCS	A	163	9/10	0.95	0.19	-	44,46,57,61	0
1	OCS	B	163	9/10	0.95	0.13	-	55,57,68,68	0

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
2	RXB	C	300	27/27	0.79	0.38	4.29	55,61,66,67	0

6.5 Other polymers ⓘ

There are no such residues in this entry.