



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2018 – 08:34 pm GMT

PDB ID : 4DSA
Title : Crystal Structure of DPP-IV with Compound C1
Authors : Xiong, B.; Zhu, L.R.; Chen, D.Q.; Zhao, Y.L.; Jiang, F.; Shen, J.K.
Deposited on : 2012-02-18
Resolution : 3.25 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.3 (157068), CSD as539be (2018)
Xtriage (Phenix) : 1.13
EDS : trunk30967
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac : 5.8.0158
CCP4 : 7.0 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk30967

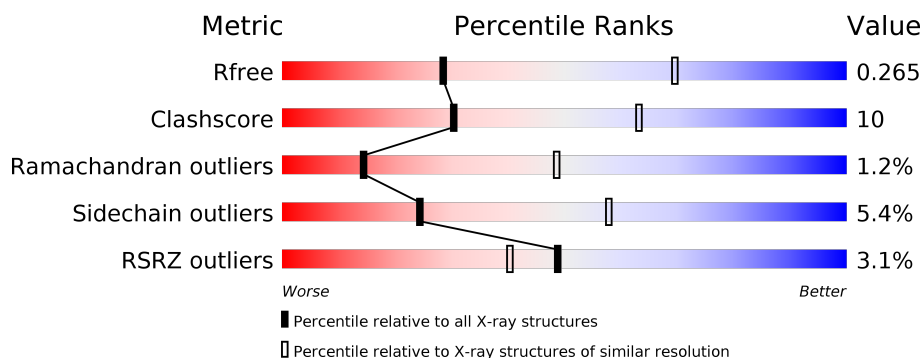
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.25 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}	111664	1008 (3.30-3.22)
Clashscore	122126	1066 (3.30-3.22)
Ramachandran outliers	120053	1046 (3.30-3.22)
Sidechain outliers	120020	1045 (3.30-3.22)
RSRZ outliers	108989	1993 (3.32-3.20)

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 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	728	<div> <div>3%</div> <div> <div></div> <div>73%</div> <div>23%</div> <div>••</div> </div> </div>
1	B	728	<div> <div>3%</div> <div> <div></div> <div>72%</div> <div>25%</div> <div>••</div> </div> </div>

2 Entry composition [i](#)

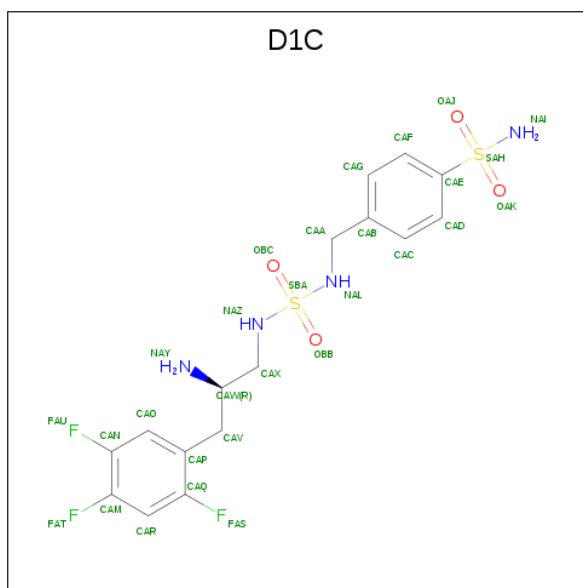
There are 2 unique types of molecules in this entry. The entry contains 11891 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

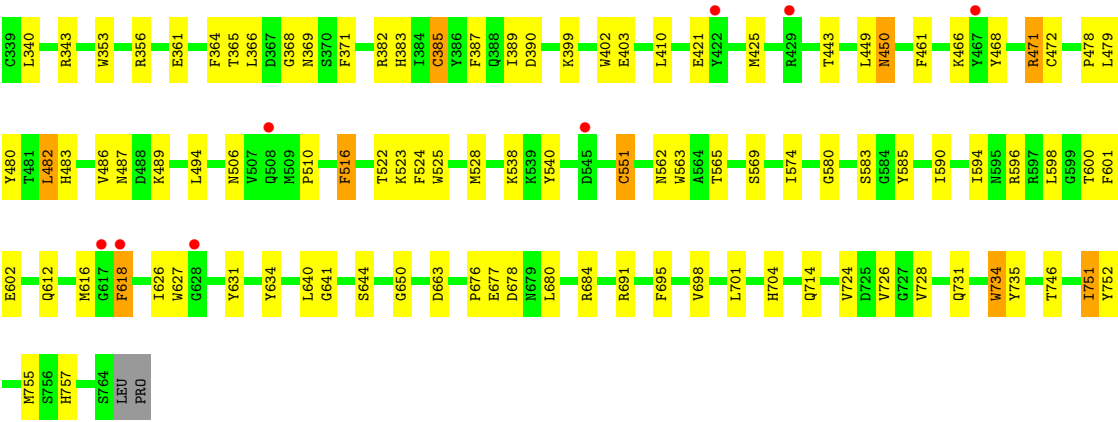
- Molecule 1 is a protein called Dipeptidyl peptidase 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	724	Total	C	N	O	S	0	0	0
			5931	3807	975	1123	26			
1	B	724	Total	C	N	O	S	0	0	0
			5931	3807	975	1123	26			

- Molecule 2 is 4-([[(2R)-2-amino-3-(2,4,5-trifluorophenyl)propyl]sulfamoyl]amino)methyl]benzenesulfonamide (three-letter code: D1C) (formula: C₁₆H₁₉F₃N₄O₄S₂).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	S	0	0
			29	16	3	4	4	2		



4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	78.68Å 78.68Å 289.85Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.37 – 3.25 29.37 – 3.25	Depositor EDS
% Data completeness (in resolution range)	97.2 (29.37-3.25) 97.3 (29.37-3.25)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.09 (at 3.24Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.198 , 0.266 0.193 , 0.265	Depositor DCC
R_{free} test set	1544 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	122.7	Xtriage
Anisotropy	0.118	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 93.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.000 for -h,-k,l 0.056 for h,-h-k,-l 0.024 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11891	wwPDB-VP
Average B, all atoms (Å ²)	144.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.15% 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: D1C

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	0.61	13/6102 (0.2%)	0.67	0/8299
1	B	0.62	10/6102 (0.2%)	0.67	0/8299
All	All	0.61	23/12204 (0.2%)	0.67	0/16598

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	337	TRP	CD2-CE2	6.13	1.48	1.41
1	A	157	TRP	CD2-CE2	5.96	1.48	1.41
1	B	337	TRP	CD2-CE2	5.92	1.48	1.41
1	B	563	TRP	CD2-CE2	5.69	1.48	1.41
1	A	187	TRP	CD2-CE2	5.65	1.48	1.41
1	A	154	TRP	CD2-CE2	5.53	1.48	1.41
1	B	157	TRP	CD2-CE2	5.52	1.48	1.41
1	B	525	TRP	CD2-CE2	5.51	1.48	1.41
1	A	315	TRP	CD2-CE2	5.50	1.48	1.41
1	A	627	TRP	CD2-CE2	5.47	1.48	1.41
1	A	62	TRP	CD2-CE2	5.44	1.47	1.41
1	B	154	TRP	CD2-CE2	5.42	1.47	1.41
1	B	187	TRP	CD2-CE2	5.38	1.47	1.41
1	B	734	TRP	CD2-CE2	5.30	1.47	1.41
1	B	305	TRP	CD2-CE2	5.27	1.47	1.41
1	B	315	TRP	CD2-CE2	5.25	1.47	1.41
1	A	734	TRP	CD2-CE2	5.21	1.47	1.41
1	B	353	TRP	CD2-CE2	5.12	1.47	1.41
1	A	353	TRP	CD2-CE2	5.10	1.47	1.41
1	A	305	TRP	CD2-CE2	5.10	1.47	1.41
1	A	525	TRP	CD2-CE2	5.06	1.47	1.41
1	A	168	TRP	CD2-CE2	5.03	1.47	1.41
1	A	216	TRP	CD2-CE2	5.00	1.47	1.41

There are no bond angle outliers.

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	5931	0	5649	118	0
1	B	5931	0	5649	121	0
2	A	29	0	19	0	0
All	All	11891	0	11317	231	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 10.

All (231) 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:175:LYS:HG3	1:A:182:SER:HB3	1.52	0.90
1:B:175:LYS:HG3	1:B:182:SER:HB3	1.56	0.86
1:B:383:HIS:HD2	1:B:399:LYS:HA	1.41	0.85
1:A:383:HIS:HD2	1:A:399:LYS:HA	1.41	0.84
1:A:383:HIS:CD2	1:A:399:LYS:HA	2.15	0.81
1:B:383:HIS:CD2	1:B:399:LYS:HA	2.14	0.80
1:B:237:GLU:HG2	1:B:253:ARG:HG2	1.61	0.80
1:A:177:GLU:HB2	1:A:180:LEU:HD22	1.63	0.79
1:A:253:ARG:HH12	1:B:253:ARG:HH12	1.35	0.74
1:A:43:TYR:CD2	1:A:565:THR:HG22	2.21	0.74
1:B:177:GLU:HB2	1:B:180:LEU:HD22	1.70	0.72
1:B:385:CYS:HB3	1:B:387:PHE:CE1	2.26	0.71
1:B:640:LEU:HD11	1:B:650:GLY:HA3	1.73	0.70
1:A:369:ASN:HA	1:A:389:ILE:HD12	1.72	0.70
1:A:385:CYS:HB3	1:A:387:PHE:CE1	2.26	0.70
1:A:237:GLU:HG2	1:A:253:ARG:HG2	1.75	0.69
1:B:369:ASN:HA	1:B:389:ILE:HD12	1.73	0.69
1:B:516:PHE:HD1	1:B:523:LYS:HB2	1.57	0.68
1:A:149:PRO:HG2	1:A:168:TRP:CD1	2.27	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:340:LEU:HB2	1:A:343:ARG:HD2	1.76	0.68
1:B:522:THR:HG21	1:B:590:ILE:HD11	1.77	0.67
1:B:596:ARG:NH2	1:B:678:ASP:OD1	2.28	0.66
1:B:115:LEU:HD21	1:B:155:VAL:HG11	1.76	0.66
1:B:43:TYR:CD2	1:B:565:THR:HG22	2.29	0.66
1:A:516:PHE:HD1	1:A:523:LYS:HB2	1.61	0.66
1:B:149:PRO:HG2	1:B:168:TRP:CD1	2.29	0.66
1:A:382:ARG:H	1:A:403:GLU:HG2	1.61	0.65
1:B:232:GLU:HB2	1:B:262:VAL:HG11	1.78	0.65
1:B:382:ARG:H	1:B:403:GLU:HG2	1.60	0.65
1:B:195:TYR:HB2	1:B:228:PHE:HB2	1.79	0.64
1:A:290:PRO:HD3	1:A:326:ASP:OD2	1.98	0.63
1:B:487:ASN:HD21	1:B:489:LYS:HE2	1.63	0.63
1:A:403:GLU:OE1	1:A:585:TYR:HA	1.99	0.63
1:B:340:LEU:HB2	1:B:343:ARG:HD2	1.81	0.63
1:A:232:GLU:HB2	1:A:262:VAL:HG11	1.82	0.62
1:A:726:VAL:HG23	1:A:728:VAL:HG23	1.82	0.62
1:B:231:THR:HG22	1:B:232:GLU:HG3	1.82	0.61
1:A:115:LEU:HD21	1:A:155:VAL:HG11	1.83	0.61
1:A:640:LEU:HD11	1:A:650:GLY:HA3	1.83	0.60
1:A:150:ASN:O	1:A:151:ASN:HB2	2.01	0.60
1:A:55:LEU:HD22	1:A:478:PRO:HG2	1.84	0.60
1:A:487:ASN:HD21	1:A:489:LYS:HE2	1.67	0.60
1:A:195:TYR:HB2	1:A:228:PHE:HB2	1.85	0.59
1:A:364:PHE:CD1	1:A:371:PHE:HB3	2.37	0.59
1:B:146:GLU:O	1:B:175:LYS:NZ	2.34	0.59
1:B:726:VAL:HG23	1:B:728:VAL:HG23	1.85	0.59
1:A:113:PHE:CZ	1:A:178:PRO:HG2	2.38	0.59
1:B:115:LEU:HD21	1:B:155:VAL:CG1	2.33	0.59
1:A:231:THR:HG22	1:A:232:GLU:HG3	1.84	0.59
1:B:598:LEU:HD22	1:B:631:TYR:OH	2.03	0.59
1:B:170:ASN:O	1:B:196:ASN:HB2	2.04	0.58
1:B:163:LYS:NZ	1:B:220:GLY:O	2.34	0.58
1:A:482:LEU:HD23	1:A:483:HIS:H	1.69	0.58
1:A:107:ILE:HG13	1:A:114:ILE:HG13	1.86	0.58
1:B:383:HIS:HE1	1:B:402:TRP:O	1.86	0.58
1:A:383:HIS:HE1	1:A:402:TRP:O	1.87	0.57
1:B:107:ILE:HG13	1:B:114:ILE:HG13	1.86	0.57
1:A:224:ALA:HB1	1:A:268:PHE:CZ	2.39	0.57
1:A:163:LYS:NZ	1:A:220:GLY:O	2.33	0.57
1:B:196:ASN:OD1	1:B:227:GLN:HG3	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:600:THR:OG1	1:A:601:PHE:N	2.38	0.56
1:B:482:LEU:HD23	1:B:483:HIS:H	1.70	0.56
1:B:385:CYS:HB3	1:B:387:PHE:HE1	1.70	0.56
1:A:115:LEU:HD21	1:A:155:VAL:CG1	2.35	0.56
1:B:290:PRO:HD3	1:B:326:ASP:OD2	2.05	0.56
1:A:596:ARG:NH2	1:A:678:ASP:OD1	2.38	0.56
1:B:364:PHE:CD1	1:B:371:PHE:HB3	2.41	0.56
1:A:258:LYS:HD2	1:B:248:TYR:CE1	2.41	0.55
1:A:229:ASN:HB3	1:A:265:THR:OG1	2.06	0.55
1:A:88:VAL:HG21	1:A:91:GLU:HG2	1.88	0.55
1:B:516:PHE:HD2	1:B:516:PHE:H	1.54	0.55
1:A:510:PRO:HD3	1:A:569:SER:HB2	1.88	0.55
1:A:305:TRP:CE2	1:A:311:ILE:HD12	2.42	0.54
1:B:88:VAL:HG21	1:B:91:GLU:HG2	1.90	0.54
1:B:224:ALA:HB1	1:B:268:PHE:CZ	2.43	0.54
1:A:636:THR:HG21	1:A:651:ILE:O	2.08	0.53
1:B:74:ASN:HB3	1:B:92:ASN:HB2	1.90	0.53
1:A:528:MET:HE1	1:A:530:LEU:HD21	1.90	0.53
1:B:55:LEU:HD22	1:B:478:PRO:HG2	1.91	0.53
1:A:216:TRP:HZ3	1:A:273:THR:HG21	1.73	0.53
1:B:540:TYR:N	1:B:618:PHE:O	2.33	0.53
1:B:216:TRP:HZ3	1:B:273:THR:HG21	1.74	0.53
1:B:516:PHE:CD1	1:B:523:LYS:HB2	2.42	0.53
1:A:640:LEU:HB3	1:A:698:VAL:HG21	1.90	0.52
1:A:516:PHE:H	1:A:516:PHE:HD2	1.56	0.52
1:B:580:GLY:O	1:B:583:SER:OG	2.26	0.52
1:A:74:ASN:HB3	1:A:92:ASN:HB2	1.92	0.52
1:B:403:GLU:OE1	1:B:585:TYR:HA	2.09	0.52
1:B:113:PHE:CZ	1:B:178:PRO:HG2	2.45	0.52
1:B:594:ILE:HD11	1:B:602:GLU:OE1	2.09	0.51
1:B:600:THR:OG1	1:B:601:PHE:N	2.43	0.51
1:A:196:ASN:OD1	1:A:227:GLN:HG3	2.11	0.51
1:A:385:CYS:HB3	1:A:387:PHE:HE1	1.71	0.51
1:A:750:HIS:CD2	1:B:724:VAL:HG22	2.46	0.51
1:B:538:LYS:O	1:B:618:PHE:HA	2.11	0.51
1:B:334:SER:OG	1:B:336:ARG:HG2	2.11	0.51
1:B:594:ILE:HD11	1:B:602:GLU:H	1.76	0.51
1:A:154:TRP:CE2	1:A:212:SER:HB3	2.46	0.50
1:A:146:GLU:O	1:A:175:LYS:NZ	2.42	0.50
1:B:191:GLU:O	1:B:193:ILE:HG13	2.11	0.50
1:A:224:ALA:HB1	1:A:268:PHE:HZ	1.74	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:74:ASN:HB3	1:A:92:ASN:CB	2.42	0.50
1:B:154:TRP:CE2	1:B:212:SER:HB3	2.47	0.50
1:B:224:ALA:HB1	1:B:268:PHE:HZ	1.75	0.50
1:B:230:ASP:OD1	1:B:264:PRO:HB3	2.12	0.50
1:A:402:TRP:NE1	1:A:421:GLU:HG3	2.27	0.50
1:A:461:PHE:CD1	1:A:468:TYR:HB3	2.47	0.50
1:A:594:ILE:HD11	1:A:602:GLU:H	1.77	0.50
1:B:330:TYR:CE1	1:B:335:GLY:HA2	2.47	0.50
1:A:594:ILE:HD11	1:A:602:GLU:OE1	2.12	0.49
1:B:74:ASN:HB3	1:B:92:ASN:CB	2.43	0.49
1:A:538:LYS:O	1:A:618:PHE:HA	2.13	0.49
1:B:229:ASN:HB3	1:B:265:THR:OG1	2.12	0.49
1:A:258:LYS:HD2	1:B:248:TYR:CZ	2.48	0.49
1:B:402:TRP:NE1	1:B:421:GLU:HG3	2.28	0.49
1:B:68:TYR:HE2	1:B:482:LEU:HD11	1.77	0.49
1:B:305:TRP:CE2	1:B:311:ILE:HD12	2.48	0.49
1:B:640:LEU:HB3	1:B:698:VAL:HG21	1.95	0.49
1:A:195:TYR:HB3	1:A:198:ILE:O	2.13	0.48
1:B:169:ASN:O	1:B:170:ASN:HB2	2.13	0.48
1:A:65:ASP:O	1:A:466:LYS:HB2	2.13	0.48
1:A:68:TYR:HE2	1:A:482:LEU:HD11	1.78	0.48
1:A:701:LEU:HA	1:A:731:GLN:HB2	1.95	0.48
1:B:125:ARG:HD2	1:B:126:HIS:NE2	2.28	0.48
1:B:235:LEU:HD13	1:B:253:ARG:HB3	1.95	0.48
1:B:516:PHE:HD1	1:B:523:LYS:CB	2.25	0.48
1:A:235:LEU:HD13	1:A:253:ARG:HB3	1.94	0.48
1:A:334:SER:OG	1:A:336:ARG:HG2	2.13	0.48
1:A:237:GLU:HG2	1:A:253:ARG:CG	2.44	0.48
1:A:680:LEU:HD11	1:A:684:ARG:CZ	2.43	0.48
1:A:594:ILE:HA	1:A:594:ILE:HD12	1.78	0.48
1:A:714:GLN:HA	1:B:241:TYR:CE1	2.48	0.48
1:B:184:ARG:HH21	1:B:187:TRP:HA	1.79	0.48
1:B:510:PRO:HD3	1:B:569:SER:HB2	1.95	0.48
1:B:150:ASN:O	1:B:151:ASN:HB2	2.14	0.47
1:A:248:TYR:CE1	1:B:258:LYS:HD2	2.49	0.47
1:B:735:TYR:OH	1:B:751:ILE:HA	2.13	0.47
1:A:598:LEU:HD22	1:A:631:TYR:OH	2.15	0.47
1:B:365:THR:O	1:B:368:GLY:N	2.46	0.47
1:A:522:THR:HG21	1:A:590:ILE:HD11	1.96	0.47
1:A:516:PHE:CD1	1:A:523:LYS:HB2	2.46	0.47
1:B:594:ILE:HD12	1:B:594:ILE:HA	1.82	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:524:PHE:HE1	1:A:590:ILE:HG12	1.80	0.47
1:B:110:ASP:OD2	1:B:162:HIS:ND1	2.29	0.47
1:B:528:MET:HE2	1:B:528:MET:HB3	1.74	0.46
1:B:302:ASP:HB3	1:B:314:GLN:HB2	1.96	0.46
1:B:524:PHE:HE1	1:B:590:ILE:HG12	1.81	0.46
1:B:402:TRP:CD1	1:B:421:GLU:HG3	2.50	0.46
1:A:528:MET:HE2	1:A:528:MET:HB3	1.68	0.46
1:A:402:TRP:CD1	1:A:421:GLU:HG3	2.50	0.46
1:A:644:SER:O	1:A:646:VAL:N	2.49	0.46
1:A:305:TRP:CZ2	1:A:311:ILE:HD12	2.51	0.45
1:A:170:ASN:O	1:A:196:ASN:HB2	2.16	0.45
1:A:159:PRO:HD3	1:A:216:TRP:CG	2.51	0.45
1:A:184:ARG:HH21	1:A:187:TRP:HA	1.82	0.45
1:A:461:PHE:CE1	1:A:468:TYR:HB3	2.52	0.45
1:A:509:MET:O	1:A:532:PRO:HB3	2.17	0.45
1:B:680:LEU:HD11	1:B:684:ARG:CZ	2.46	0.45
1:B:383:HIS:CE1	1:B:402:TRP:O	2.69	0.45
1:A:528:MET:CE	1:A:530:LEU:HD21	2.46	0.45
1:A:735:TYR:OH	1:A:751:ILE:HA	2.17	0.45
1:B:153:GLN:OE1	1:B:198:ILE:HD13	2.17	0.45
1:B:471:ARG:HD3	1:B:480:TYR:CE2	2.52	0.44
1:A:72:GLN:O	1:A:74:ASN:N	2.50	0.44
1:B:117:GLU:OE1	1:B:128:TYR:HE1	2.00	0.44
1:A:725:ASP:HA	1:B:746:THR:HG21	1.99	0.44
1:A:383:HIS:CE1	1:A:402:TRP:O	2.69	0.44
1:A:241:TYR:CE1	1:B:714:GLN:HA	2.52	0.44
1:A:56:LYS:HB2	1:A:497:ASN:OD1	2.17	0.44
1:B:45:LEU:HG	1:B:49:LEU:HD22	1.99	0.44
1:B:48:TYR:CE1	1:B:562:ASN:HA	2.53	0.44
1:A:330:TYR:CE1	1:A:335:GLY:HA2	2.53	0.44
1:A:531:PRO:HB3	1:A:572:ASN:HA	2.00	0.44
1:A:43:TYR:CD2	1:A:565:THR:CG2	2.97	0.44
1:B:676:PRO:HD2	1:B:677:GLU:OE1	2.18	0.44
1:B:70:TYR:HB3	1:B:79:PHE:HE1	1.84	0.43
1:B:482:LEU:HD12	1:B:494:LEU:HD21	2.00	0.43
1:A:369:ASN:HA	1:A:389:ILE:CD1	2.44	0.43
1:B:594:ILE:CD1	1:B:601:PHE:HB2	2.48	0.43
1:A:113:PHE:CE2	1:A:178:PRO:HG2	2.53	0.43
1:A:516:PHE:HD1	1:A:523:LYS:CB	2.30	0.43
1:B:95:PHE:HB3	1:B:98:PHE:HB2	2.00	0.43
1:B:162:HIS:NE2	1:B:177:GLU:OE1	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:84:GLY:HA3	1:A:492:ARG:HH22	1.83	0.43
1:A:658:ARG:HB2	1:A:689:MET:CE	2.49	0.43
1:B:237:GLU:HG2	1:B:253:ARG:CG	2.41	0.43
1:B:734:TRP:C	1:B:734:TRP:CD1	2.90	0.43
1:A:169:ASN:O	1:A:170:ASN:HB2	2.19	0.42
1:A:48:TYR:CE1	1:A:562:ASN:HA	2.54	0.42
1:B:701:LEU:HA	1:B:731:GLN:HB2	2.00	0.42
1:B:65:ASP:O	1:B:466:LYS:HB2	2.19	0.42
1:A:446:SER:HA	1:A:449:LEU:HG	2.01	0.42
1:B:468:TYR:CE2	1:B:483:HIS:HB2	2.54	0.42
1:B:72:GLN:O	1:B:74:ASN:N	2.51	0.42
1:A:301:CYS:SG	1:A:359:PRO:HD2	2.59	0.42
1:A:45:LEU:HG	1:A:49:LEU:HD22	2.01	0.42
1:B:528:MET:CE	1:B:574:ILE:HG21	2.50	0.42
1:A:677:GLU:N	1:A:677:GLU:OE1	2.51	0.42
1:B:256:TYR:CZ	1:B:663:ASP:HB3	2.54	0.42
1:A:471:ARG:HD3	1:A:480:TYR:CE2	2.55	0.42
1:B:195:TYR:HB3	1:B:198:ILE:O	2.19	0.42
1:B:109:PRO:HG2	1:B:158:SER:O	2.18	0.42
1:B:461:PHE:CD1	1:B:468:TYR:HB3	2.55	0.42
1:B:735:TYR:CE2	1:B:751:ILE:HG13	2.55	0.41
1:A:612:GLN:O	1:A:616:MET:HG3	2.20	0.41
1:A:540:TYR:N	1:A:618:PHE:O	2.36	0.41
1:A:256:TYR:CZ	1:A:663:ASP:HB3	2.55	0.41
1:A:159:PRO:HD3	1:A:216:TRP:CB	2.50	0.41
1:A:365:THR:O	1:A:368:GLY:N	2.53	0.41
1:A:95:PHE:HB3	1:A:98:PHE:HB2	2.02	0.41
1:A:230:ASP:OD1	1:A:264:PRO:HB3	2.19	0.41
1:A:676:PRO:HD2	1:A:677:GLU:OE1	2.21	0.41
1:A:562:ASN:O	1:A:565:THR:HB	2.20	0.41
1:B:157:TRP:HA	1:B:163:LYS:O	2.21	0.41
1:B:369:ASN:HA	1:B:389:ILE:CD1	2.46	0.41
1:B:191:GLU:C	1:B:193:ILE:H	2.24	0.41
1:B:641:GLY:O	1:B:691:ARG:HB3	2.21	0.41
1:A:383:HIS:CD2	1:A:399:LYS:CA	2.98	0.41
1:B:612:GLN:O	1:B:616:MET:HG3	2.21	0.41
1:A:530:LEU:HA	1:A:531:PRO:HD3	1.99	0.41
1:A:302:ASP:HB3	1:A:314:GLN:HB2	2.03	0.40
1:A:509:MET:HB3	1:A:509:MET:HE2	1.98	0.40
1:B:626:ILE:O	1:B:650:GLY:HA2	2.20	0.40
1:A:117:GLU:OE1	1:A:128:TYR:HE1	2.05	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:237:GLU:HA	1:B:252:VAL:O	2.21	0.40
1:B:461:PHE:CE1	1:B:468:TYR:HB3	2.57	0.40
1:B:695:PHE:HB3	1:B:728:VAL:HG11	2.04	0.40
1:B:752:TYR:HA	1:B:755:MET:HE2	2.02	0.40
1:B:113:PHE:CE2	1:B:178:PRO:HG2	2.56	0.40
1:B:631:TYR:O	1:B:634:TYR:HB3	2.20	0.40
1:B:356:ARG:HD3	1:B:551:CYS:SG	2.61	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	722/728 (99%)	643 (89%)	69 (10%)	10 (1%)	12	45
1	B	722/728 (99%)	640 (89%)	75 (10%)	7 (1%)	17	52
All	All	1444/1456 (99%)	1283 (89%)	144 (10%)	17 (1%)	14	48

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	178	PRO
1	A	644	SER
1	A	645	GLY
1	B	169	ASN
1	B	178	PRO
1	B	644	SER
1	A	73	GLU
1	A	169	ASN
1	A	341	VAL
1	A	551	CYS
1	A	760	LYS

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Mol	Chain	Res	Type
1	B	73	GLU
1	A	565	THR
1	B	551	CYS
1	A	279	VAL
1	B	279	VAL
1	B	450	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	649/653 (99%)	615 (95%)	34 (5%)	25	58
1	B	649/653 (99%)	613 (94%)	36 (6%)	24	57
All	All	1298/1306 (99%)	1228 (95%)	70 (5%)	24	58

All (70) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	LEU
1	A	51	ASN
1	A	87	SER
1	A	156	THR
1	A	180	LEU
1	A	182	SER
1	A	231	THR
1	A	242	SER
1	A	244	GLU
1	A	288	THR
1	A	301	CYS
1	A	312	SER
1	A	329	ASP
1	A	361	GLU
1	A	366	LEU
1	A	382	ARG
1	A	385	CYS

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Mol	Chain	Res	Type
1	A	390	ASP
1	A	410	LEU
1	A	412	SER
1	A	425	MET
1	A	443	THR
1	A	450	ASN
1	A	471	ARG
1	A	479	LEU
1	A	482	LEU
1	A	486	VAL
1	A	506	ASN
1	A	516	PHE
1	A	522	THR
1	A	618	PHE
1	A	627	TRP
1	A	704	HIS
1	A	751	ILE
1	B	49	LEU
1	B	51	ASN
1	B	87	SER
1	B	156	THR
1	B	169	ASN
1	B	180	LEU
1	B	182	SER
1	B	211	TYR
1	B	231	THR
1	B	242	SER
1	B	244	GLU
1	B	284	SER
1	B	288	THR
1	B	299	TYR
1	B	329	ASP
1	B	361	GLU
1	B	366	LEU
1	B	385	CYS
1	B	390	ASP
1	B	410	LEU
1	B	425	MET
1	B	443	THR
1	B	449	LEU
1	B	450	ASN
1	B	471	ARG

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Mol	Chain	Res	Type
1	B	472	CYS
1	B	479	LEU
1	B	482	LEU
1	B	486	VAL
1	B	506	ASN
1	B	516	PHE
1	B	618	PHE
1	B	627	TRP
1	B	704	HIS
1	B	751	ILE
1	B	757	HIS

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

Mol	Chain	Res	Type
1	A	138	ASN
1	A	247	GLN
1	A	430	ASN
1	A	748	HIS
1	B	66	HIS
1	B	138	ASN
1	B	247	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

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	D1C	A	800	-	30,30,30	5.25	11 (36%)	36,44,44	2.30	8 (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	D1C	A	800	-	-	0/21/22/22	0/2/2/2

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	800	D1C	SBA-NAZ	-17.48	1.40	1.61
2	A	800	D1C	CAV-CAP	-10.03	1.37	1.51
2	A	800	D1C	CAA-CAB	-4.57	1.41	1.51
2	A	800	D1C	CAE-SAH	-4.25	1.70	1.77
2	A	800	D1C	FAU-CAN	-2.85	1.28	1.35
2	A	800	D1C	FAS-CAQ	2.28	1.41	1.35
2	A	800	D1C	OAK-SAH	5.08	1.53	1.43
2	A	800	D1C	OAJ-SAH	5.19	1.54	1.43
2	A	800	D1C	OBB-SBA	7.06	1.53	1.43
2	A	800	D1C	OBC-SBA	7.49	1.54	1.43
2	A	800	D1C	SBA-NAL	14.08	1.78	1.61

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	800	D1C	OBC-SBA-OBB	-8.97	107.01	120.33
2	A	800	D1C	OAJ-SAH-OAK	-4.38	111.09	118.71
2	A	800	D1C	CAR-CAQ-CAP	-3.26	119.48	123.97
2	A	800	D1C	FAU-CAN-CAO	2.33	123.07	118.58
2	A	800	D1C	OBC-SBA-NAL	2.58	111.27	106.73
2	A	800	D1C	FAT-CAM-CAR	2.58	123.55	118.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	800	D1C	OBB-SBA-NAL	4.06	113.89	106.73
2	A	800	D1C	CAO-CAP-CAQ	4.55	121.91	116.70

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	724/728 (99%)	0.04	22 (3%)	50 42	87, 141, 186, 246	0
1	B	724/728 (99%)	0.06	23 (3%)	47 38	86, 144, 192, 229	0
All	All	1448/1456 (99%)	0.05	45 (3%)	49 40	86, 143, 189, 246	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	524	PHE	5.5
1	A	82	GLU	4.9
1	B	309	GLU	4.7
1	A	83	TYR	4.5
1	B	138	ASN	3.7
1	A	455	GLN	3.7
1	A	628	GLY	3.4
1	A	632	GLY	3.2
1	B	467	TYR	3.2
1	A	545	ASP	3.1
1	B	100	HIS	3.0
1	A	42	THR	3.0
1	B	618	PHE	3.0
1	A	633	GLY	2.8
1	B	628	GLY	2.8
1	B	103	ASN	2.7
1	B	97	GLU	2.7
1	B	101	SER	2.7
1	B	429	ARG	2.7
1	B	179	ASN	2.6
1	B	334	SER	2.6
1	A	630	SER	2.5
1	B	96	ASP	2.5
1	A	471	ARG	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	636	THR	2.4
1	A	138	ASN	2.4
1	B	42	THR	2.4
1	B	545	ASP	2.3
1	A	674	PRO	2.3
1	A	657	SER	2.3
1	B	617	GLY	2.2
1	A	122	LYS	2.2
1	A	309	GLU	2.2
1	A	91	GLU	2.2
1	B	122	LYS	2.1
1	A	222	PHE	2.1
1	B	125	ARG	2.1
1	B	338	ASN	2.1
1	B	41	LYS	2.1
1	A	41	LYS	2.1
1	A	726	VAL	2.1
1	B	422	TYR	2.0
1	B	89	PHE	2.0
1	B	508	GLN	2.0
1	A	539	LYS	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 carbohydrates 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(\AA^2)	Q<0.9
2	D1C	A	800	29/29	0.90	0.27	98,148,204,207	0

6.5 Other polymers [i](#)

There are no such residues in this entry.