



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 01:56 am GMT

PDB ID : 5CTO
Title : Crystal structure of Arabidopsis thaliana HPPD complexed with NTBC
Authors : Yang, W.C.; Yang, G.F.
Deposited on : 2015-07-24
Resolution : 2.62 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 : trunk28620
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) : recalc28949

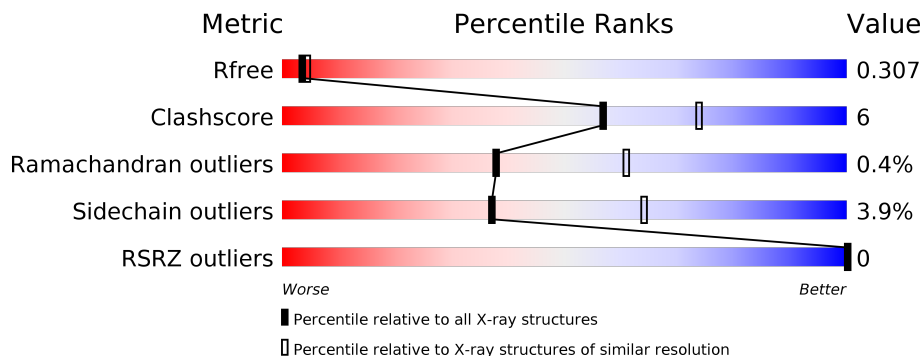
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.62 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	2983 (2.64-2.60)
Clashscore	112137	3351 (2.64-2.60)
Ramachandran outliers	110173	3298 (2.64-2.60)
Sidechain outliers	110143	3298 (2.64-2.60)
RSRZ outliers	101464	2992 (2.64-2.60)

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	445	
1	B	445	
1	C	445	
1	D	445	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11167 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 4-hydroxyphenylpyruvate dioxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	367	Total	C	N	O	S	1	0	0
			2770	1772	476	510	12			
1	B	370	Total	C	N	O	S	1	0	0
			2782	1780	478	511	13			
1	C	365	Total	C	N	O	S	1	0	0
			2750	1761	465	513	11			
1	D	370	Total	C	N	O	S	1	0	0
			2751	1758	472	509	12			

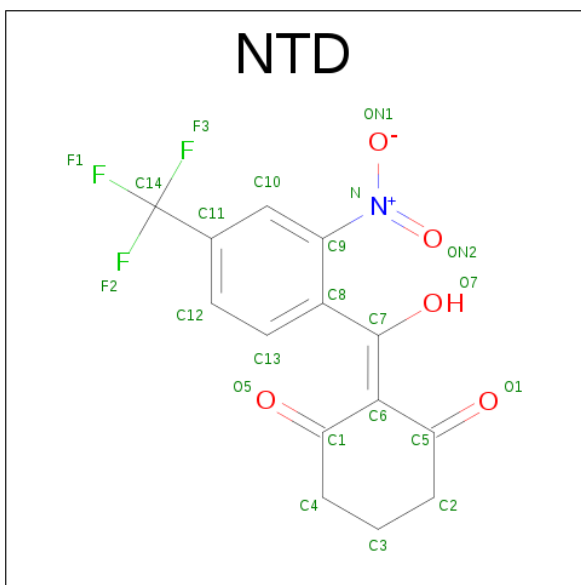
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	434	PHE	TYR	conflict	UNP P93836
B	434	PHE	TYR	conflict	UNP P93836
C	434	PHE	TYR	conflict	UNP P93836
D	434	PHE	TYR	conflict	UNP P93836

- Molecule 2 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Fe	0	0
			1	1		
2	A	1	Total	Fe	0	0
			1	1		
2	D	1	Total	Fe	0	0
			1	1		
2	C	1	Total	Fe	0	0
			1	1		

- Molecule 3 is 2-{HYDROXY[2-NITRO-4-(TRIFLUOROMETHYL)PHENYL]METHYLEN E}CYCLOHEXANE-1,3-DIONE (three-letter code: NTD) (formula: C₁₄H₁₀F₃NO₅).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	0	0
			23	14	3	1	5		
3	B	1	Total	C	F	N	O	0	0
			23	14	3	1	5		
3	C	1	Total	C	F	N	O	0	0
			23	14	3	1	5		
3	D	1	Total	C	F	N	O	0	0
			23	14	3	1	5		

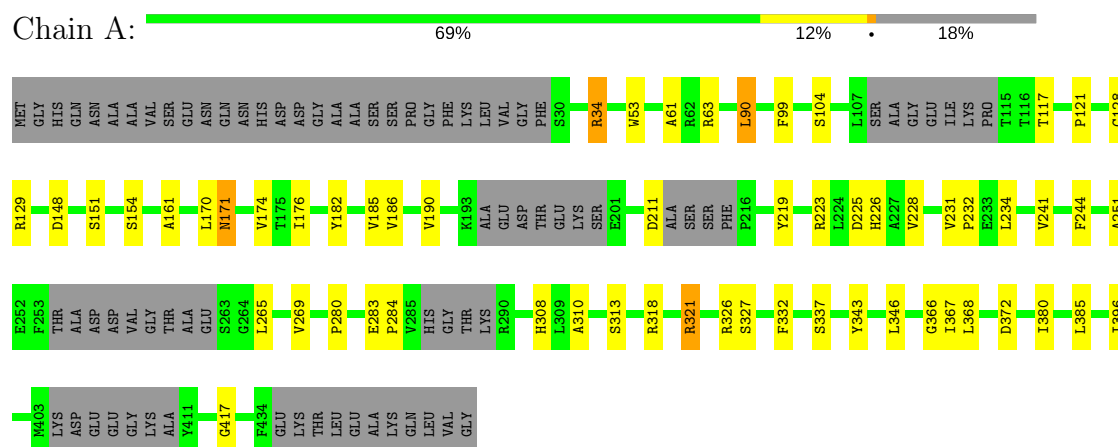
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	4	Total	O	0	0
			4	4		
4	B	5	Total	O	0	0
			5	5		
4	C	5	Total	O	0	0
			5	5		
4	D	4	Total	O	0	0
			4	4		

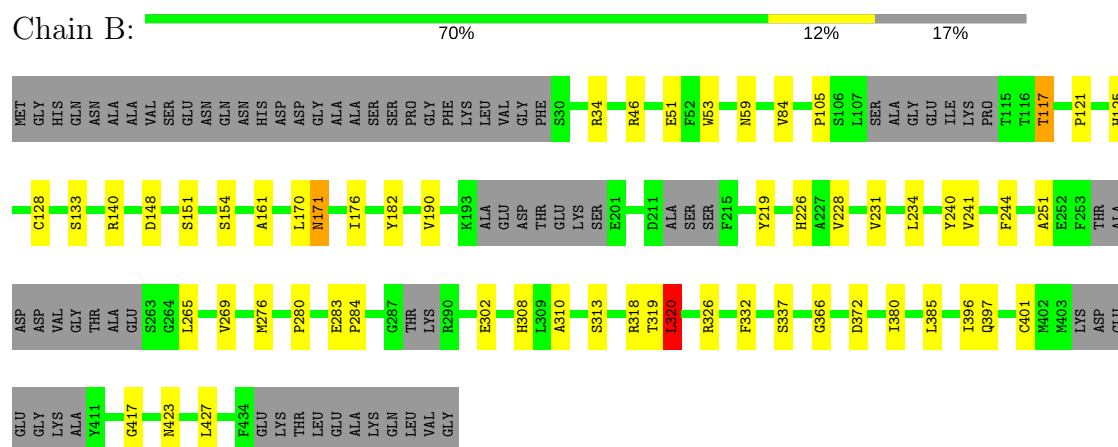
3 Residue-property plots

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.

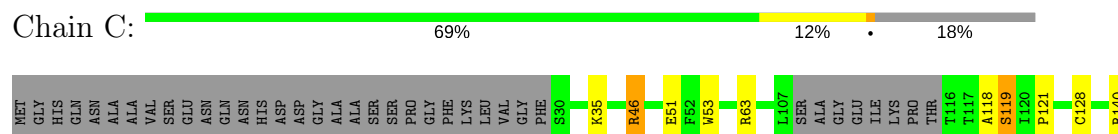
• Molecule 1: 4-hydroxyphenylpyruvate dioxygenase

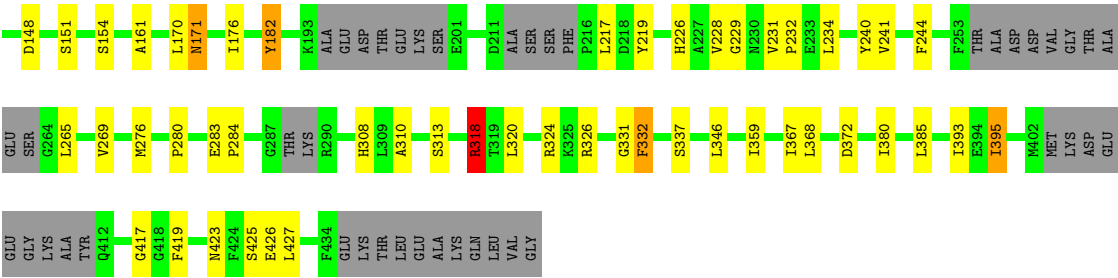


• Molecule 1: 4-hydroxyphenylpyruvate dioxygenase

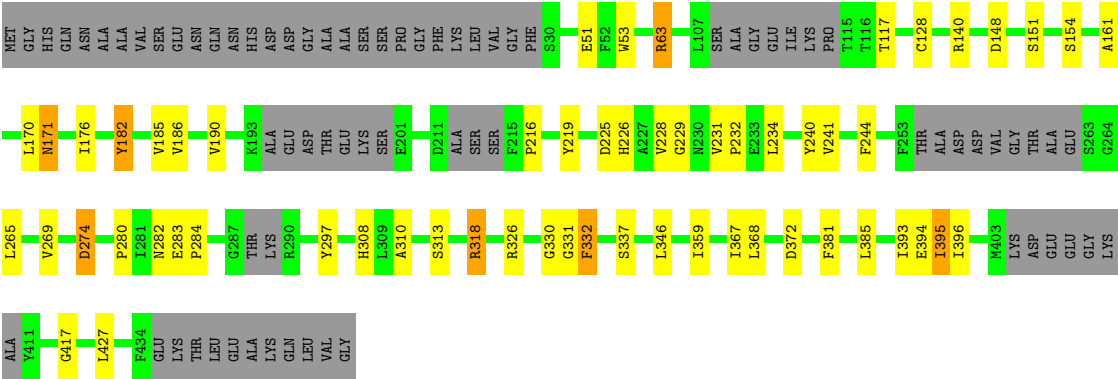


• Molecule 1: 4-hydroxyphenylpyruvate dioxygenase





● Molecule 1: 4-hydroxyphenylpyruvate dioxygenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	95.66Å 95.96Å 97.93Å 90.00° 92.15° 90.00°	Depositor
Resolution (Å)	30.26 – 2.62 29.19 – 2.62	Depositor EDS
% Data completeness (in resolution range)	98.9 (30.26-2.62) 99.0 (29.19-2.62)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.57	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.46 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.269 , 0.306 0.274 , 0.307	Depositor DCC
R_{free} test set	2672 reflections (5.36%)	DCC
Wilson B-factor (Å ²)	24.9	Xtriage
Anisotropy	0.361	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 0.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.22$	Xtriage
Estimated twinning fraction	0.034 for -h,l,k 0.047 for -h,-l,-k 0.046 for -l,k,h 0.167 for -k,-h,-l 0.165 for k,h,-l 0.044 for k,l,h 0.043 for l,h,k 0.042 for l,-h,-k 0.040 for -k,-l,h 0.145 for h,-k,-l 0.064 for l,-k,h	Xtriage
F_o, F_c correlation	0.82	EDS
Total number of atoms	11167	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.18% 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: NTD, FE

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	0.64	0/2832	0.84	3/3825 (0.1%)
1	B	0.64	1/2846 (0.0%)	0.77	2/3845 (0.1%)
1	C	0.67	1/2813 (0.0%)	0.82	6/3802 (0.2%)
1	D	0.67	1/2813 (0.0%)	0.80	4/3806 (0.1%)
All	All	0.66	3/11304 (0.0%)	0.81	15/15278 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1
1	D	0	1
All	All	0	2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	318	ARG	CB-CG	-8.94	1.28	1.52
1	D	318	ARG	CZ-NH1	7.69	1.43	1.33
1	B	318	ARG	CZ-NH1	5.81	1.40	1.33

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	318	ARG	NE-CZ-NH1	17.56	129.08	120.30
1	A	318	ARG	NH1-CZ-NH2	-8.40	110.17	119.40
1	C	318	ARG	CG-CD-NE	7.92	128.43	111.80
1	C	318	ARG	CA-CB-CG	7.42	129.72	113.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	119	SER	N-CA-CB	-7.36	99.46	110.50
1	B	318	ARG	NE-CZ-NH1	6.78	123.69	120.30
1	B	320	LEU	CA-CB-CG	-6.67	99.95	115.30
1	D	427	LEU	CA-CB-CG	6.52	130.30	115.30
1	D	274	ASP	CB-CG-OD2	-6.49	112.46	118.30
1	C	395	ILE	CG1-CB-CG2	-6.42	97.26	111.40
1	C	331	GLY	N-CA-C	6.26	128.75	113.10
1	C	318	ARG	CB-CG-CD	6.24	127.83	111.60
1	D	395	ILE	CG1-CB-CG2	-6.00	98.21	111.40
1	D	331	GLY	N-CA-C	5.45	126.73	113.10
1	A	321	ARG	NE-CZ-NH1	5.35	122.98	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	118	ALA	Peptide
1	D	318	ARG	Sidechain

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	2770	0	2640	29	0
1	B	2782	0	2626	32	0
1	C	2750	0	2600	33	0
1	D	2751	0	2584	31	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	23	0	9	1	0
3	B	23	0	9	1	0
3	C	23	0	9	2	0
3	D	23	0	9	2	0
4	A	4	0	0	0	0
4	B	5	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	5	0	0	0	0
4	D	4	0	0	0	0
All	All	11167	0	10486	124	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 6.

All (124) 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:C:320:LEU:HD21	1:C:380:ILE:HD13	1.75	0.68
3:A:502:NTD:O5	3:A:502:NTD:C8	2.40	0.67
1:C:53:TRP:CZ3	1:C:128:CYS:HA	2.30	0.67
1:B:53:TRP:CZ3	1:B:128:CYS:HA	2.30	0.67
1:A:90:LEU:HD22	1:A:99:PHE:CE1	2.31	0.66
1:D:53:TRP:CZ3	1:D:128:CYS:HA	2.31	0.65
1:A:53:TRP:CZ3	1:A:128:CYS:HA	2.32	0.64
3:D:502:NTD:O5	3:D:502:NTD:C8	2.48	0.60
1:A:225:ASP:CB	1:A:396:ILE:HD11	2.31	0.60
1:C:121:PRO:HG3	1:C:170:LEU:HD21	1.85	0.58
1:A:121:PRO:HG3	1:A:170:LEU:HD21	1.87	0.57
1:B:121:PRO:HG3	1:B:170:LEU:HD21	1.87	0.56
1:C:228:VAL:HG21	1:C:308:HIS:CE1	2.41	0.56
1:A:228:VAL:HG21	1:A:308:HIS:CE1	2.41	0.56
1:B:228:VAL:HG21	1:B:308:HIS:CE1	2.42	0.55
1:C:423:ASN:O	1:C:427:LEU:HB2	2.06	0.55
1:C:346:LEU:HD23	1:C:359:ILE:HG23	1.88	0.55
1:C:320:LEU:HD21	1:C:380:ILE:HG21	1.90	0.54
1:B:423:ASN:O	1:B:427:LEU:HB2	2.07	0.54
1:A:321:ARG:HD2	1:C:318:ARG:NH2	2.22	0.54
3:B:502:NTD:O5	3:B:502:NTD:C8	2.56	0.54
1:D:228:VAL:HG21	1:D:308:HIS:CE1	2.42	0.54
1:D:171:ASN:OD1	1:D:171:ASN:N	2.41	0.53
1:D:346:LEU:HD23	1:D:359:ILE:HG23	1.89	0.53
1:B:320:LEU:HD21	1:B:380:ILE:HG21	1.90	0.52
1:D:225:ASP:HB3	1:D:396:ILE:HD11	1.92	0.52
1:A:225:ASP:HB3	1:A:396:ILE:HD11	1.92	0.51
1:D:225:ASP:CB	1:D:396:ILE:HD11	2.39	0.51
1:B:148:ASP:O	1:B:151:SER:HB3	2.11	0.51
1:D:381:PHE:CE2	1:D:394:GLU:HB3	2.46	0.51
1:B:171:ASN:OD1	1:B:171:ASN:N	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:148:ASP:O	1:C:151:SER:HB3	2.11	0.51
1:C:171:ASN:N	1:C:171:ASN:OD1	2.44	0.50
1:A:148:ASP:O	1:A:151:SER:HB3	2.11	0.50
1:A:219:TYR:OH	1:A:326:ARG:NE	2.45	0.49
1:B:34:ARG:NH1	1:B:251:ALA:O	2.42	0.49
1:D:381:PHE:CD2	1:D:394:GLU:HB3	2.47	0.49
1:A:171:ASN:OD1	1:A:171:ASN:N	2.45	0.49
1:D:51:GLU:O	1:D:140:ARG:N	2.43	0.49
1:D:148:ASP:O	1:D:151:SER:HB3	2.13	0.48
1:C:161:ALA:HB2	1:C:244:PHE:CG	2.49	0.47
1:D:161:ALA:HB2	1:D:244:PHE:CG	2.50	0.47
1:C:46:ARG:CZ	1:C:276:MET:HG2	2.44	0.47
1:A:226:HIS:CE1	1:A:310:ALA:CB	2.97	0.47
1:D:282:ASN:HD21	3:D:502:NTD:H31	1.79	0.47
1:A:265:LEU:C	1:A:265:LEU:HD12	2.35	0.47
1:A:63:ARG:HD3	1:B:59:ASN:OD1	2.15	0.47
1:B:219:TYR:OH	1:B:326:ARG:NE	2.48	0.46
1:A:34:ARG:NH2	1:A:251:ALA:O	2.46	0.46
1:D:274:ASP:N	1:D:274:ASP:OD1	2.48	0.46
1:B:161:ALA:HB2	1:B:244:PHE:CG	2.51	0.46
1:A:61:ALA:O	1:A:90:LEU:HD21	2.16	0.46
1:A:283:GLU:HG3	1:A:284:PRO:HD2	1.98	0.46
1:D:283:GLU:HG3	1:D:284:PRO:HD2	1.99	0.46
1:D:231:VAL:HB	1:D:232:PRO:HD2	1.97	0.45
1:D:226:HIS:CE1	1:D:310:ALA:CB	2.99	0.45
1:C:419:PHE:HB3	3:C:502:NTD:O1	2.16	0.45
1:D:63:ARG:HD2	1:D:330:GLY:O	2.16	0.45
1:C:51:GLU:O	1:C:140:ARG:N	2.46	0.45
1:A:343:TYR:O	1:A:346:LEU:HB2	2.17	0.45
1:B:46:ARG:CZ	1:B:276:MET:HG3	2.47	0.45
1:C:226:HIS:CE1	1:C:310:ALA:CB	3.00	0.45
1:B:105:PRO:HG3	1:B:125:HIS:CD2	2.52	0.45
1:B:234:LEU:HB2	1:B:283:GLU:HB2	1.98	0.44
1:D:332:PHE:CE1	1:D:385:LEU:HD22	2.52	0.44
1:C:219:TYR:OH	1:C:326:ARG:NE	2.50	0.44
1:B:283:GLU:HG3	1:B:284:PRO:HD2	1.99	0.44
1:C:393:ILE:HG22	1:C:395:ILE:HD12	2.00	0.44
1:A:161:ALA:HB2	1:A:244:PHE:CG	2.52	0.44
1:B:240:TYR:CE2	1:B:244:PHE:HE2	2.36	0.44
1:D:240:TYR:CE2	1:D:244:PHE:HE2	2.35	0.44
1:C:372:ASP:OD1	1:C:417:GLY:HA2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:419:PHE:HB3	3:C:502:NTD:C5	2.48	0.43
1:A:367:ILE:HG22	1:A:368:LEU:O	2.19	0.43
1:A:372:ASP:OD1	1:A:417:GLY:HA2	2.17	0.43
1:A:104:SER:OG	1:B:302:GLU:OE2	2.23	0.43
1:C:231:VAL:HB	1:C:232:PRO:HD2	1.99	0.43
1:D:265:LEU:C	1:D:265:LEU:HD12	2.38	0.43
1:A:129:ARG:CZ	1:B:133:SER:HA	2.48	0.43
1:B:51:GLU:O	1:B:140:ARG:N	2.48	0.43
1:D:367:ILE:HG22	1:D:368:LEU:O	2.19	0.43
1:C:240:TYR:CE2	1:C:244:PHE:HE2	2.36	0.43
1:D:269:VAL:HA	1:D:280:PRO:HA	2.00	0.43
1:A:332:PHE:CE1	1:A:385:LEU:HD22	2.54	0.42
1:B:332:PHE:CE1	1:B:385:LEU:HD22	2.53	0.42
1:C:234:LEU:HB2	1:C:283:GLU:HB2	2.00	0.42
1:A:231:VAL:HB	1:A:232:PRO:HD2	2.00	0.42
1:A:269:VAL:HA	1:A:280:PRO:HA	2.01	0.42
1:D:234:LEU:HB2	1:D:283:GLU:HB2	1.99	0.42
1:C:332:PHE:CE1	1:C:385:LEU:HD22	2.54	0.42
1:C:283:GLU:HG3	1:C:284:PRO:HD2	2.01	0.42
1:B:269:VAL:HA	1:B:280:PRO:HA	2.00	0.42
1:D:219:TYR:OH	1:D:326:ARG:NE	2.53	0.42
1:D:297:TYR:CD2	1:D:297:TYR:C	2.93	0.42
1:C:367:ILE:HG22	1:C:368:LEU:O	2.20	0.42
1:B:320:LEU:HD21	1:B:380:ILE:HD13	2.01	0.42
1:A:241:VAL:HA	1:A:244:PHE:CE2	2.54	0.41
1:C:425:SER:O	1:C:426:GLU:C	2.58	0.41
1:D:182:TYR:CZ	1:D:229:GLY:HA3	2.55	0.41
1:A:327:SER:OG	1:A:332:PHE:O	2.34	0.41
1:C:53:TRP:CE3	1:C:128:CYS:HA	2.56	0.41
1:C:265:LEU:C	1:C:265:LEU:HD12	2.41	0.41
1:B:396:ILE:CG1	1:B:397:GLN:N	2.80	0.41
1:B:84:VAL:HG23	1:B:117:THR:HG23	2.02	0.41
1:C:241:VAL:HA	1:C:244:PHE:CE2	2.55	0.41
1:D:228:VAL:HG22	1:D:280:PRO:HB2	2.01	0.41
1:D:241:VAL:HA	1:D:244:PHE:CE2	2.55	0.41
1:A:185:VAL:HG12	1:A:186:VAL:N	2.36	0.41
1:B:401:CYS:N	4:B:603:HOH:O	2.49	0.41
1:B:226:HIS:CE1	1:B:310:ALA:CB	3.04	0.41
1:B:241:VAL:HA	1:B:244:PHE:CE2	2.56	0.41
1:B:372:ASP:OD1	1:B:417:GLY:HA2	2.21	0.41
1:D:185:VAL:HG12	1:D:186:VAL:N	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:228:VAL:HG22	1:C:280:PRO:HB2	2.03	0.40
1:C:320:LEU:O	1:C:324:ARG:HG3	2.21	0.40
1:B:228:VAL:HG22	1:B:280:PRO:HB2	2.03	0.40
1:B:265:LEU:HD12	1:B:265:LEU:C	2.42	0.40
1:C:182:TYR:CZ	1:C:229:GLY:HA3	2.56	0.40
1:C:269:VAL:HA	1:C:280:PRO:HA	2.03	0.40
1:A:366:GLY:O	1:A:380:ILE:HD12	2.20	0.40
1:B:219:TYR:O	1:B:319:THR:OG1	2.31	0.40
1:D:393:ILE:HG22	1:D:395:ILE:HD12	2.03	0.40
1:B:366:GLY:O	1:B:380:ILE:HD12	2.22	0.40
1:D:372:ASP:OD1	1:D:417:GLY:HA2	2.20	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	353/445 (79%)	328 (93%)	25 (7%)	0	100	100
1	B	356/445 (80%)	332 (93%)	24 (7%)	0	100	100
1	C	351/445 (79%)	322 (92%)	26 (7%)	3 (1%)	20	38
1	D	356/445 (80%)	331 (93%)	23 (6%)	2 (1%)	28	51
All	All	1416/1780 (80%)	1313 (93%)	98 (7%)	5 (0%)	38	61

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	119	SER
1	C	217	LEU
1	C	332	PHE
1	D	216	PRO

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Mol	Chain	Res	Type
1	D	332	PHE

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	285/371 (77%)	271 (95%)	14 (5%)	29	53
1	B	283/371 (76%)	273 (96%)	10 (4%)	41	67
1	C	282/371 (76%)	272 (96%)	10 (4%)	41	67
1	D	278/371 (75%)	268 (96%)	10 (4%)	40	67
All	All	1128/1484 (76%)	1084 (96%)	44 (4%)	37	64

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	34	ARG
1	A	90	LEU
1	A	117	THR
1	A	154	SER
1	A	171	ASN
1	A	174	VAL
1	A	176	ILE
1	A	182	TYR
1	A	190	VAL
1	A	211	ASP
1	A	223	ARG
1	A	234	LEU
1	A	313	SER
1	A	337	SER
1	B	117	THR
1	B	154	SER
1	B	171	ASN
1	B	176	ILE
1	B	182	TYR
1	B	190	VAL

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Mol	Chain	Res	Type
1	B	231	VAL
1	B	313	SER
1	B	320	LEU
1	B	337	SER
1	C	35	LYS
1	C	46	ARG
1	C	63	ARG
1	C	154	SER
1	C	171	ASN
1	C	176	ILE
1	C	182	TYR
1	C	313	SER
1	C	318	ARG
1	C	337	SER
1	D	63	ARG
1	D	117	THR
1	D	154	SER
1	D	170	LEU
1	D	171	ASN
1	D	176	ILE
1	D	182	TYR
1	D	190	VAL
1	D	313	SER
1	D	337	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	159	ASN
1	D	159	ASN

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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
3	NTD	A	502	2	21,24,24	1.98	3 (14%)	28,36,36	1.52	4 (14%)
3	NTD	B	502	2	21,24,24	1.56	2 (9%)	28,36,36	1.59	4 (14%)
3	NTD	C	502	2	21,24,24	1.65	3 (14%)	28,36,36	1.43	4 (14%)
3	NTD	D	502	2	21,24,24	1.86	3 (14%)	28,36,36	1.50	4 (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	NTD	A	502	2	-	0/16/32/32	0/2/2/2
3	NTD	B	502	2	-	0/16/32/32	0/2/2/2
3	NTD	C	502	2	-	0/16/32/32	0/2/2/2
3	NTD	D	502	2	-	0/16/32/32	0/2/2/2

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	502	NTD	O7-C7	-3.91	1.22	1.33
3	B	502	NTD	O7-C7	-3.78	1.22	1.33
3	D	502	NTD	O7-C7	-3.73	1.22	1.33
3	C	502	NTD	O7-C7	-3.70	1.22	1.33
3	C	502	NTD	C6-C1	2.63	1.52	1.46
3	D	502	NTD	C6-C5	4.24	1.56	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	502	NTD	C6-C5	5.21	1.58	1.46
3	B	502	NTD	C6-C7	5.52	1.52	1.39
3	C	502	NTD	C6-C7	5.67	1.52	1.39
3	A	502	NTD	C6-C7	5.76	1.52	1.39
3	D	502	NTD	C6-C7	6.04	1.53	1.39

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	502	NTD	C8-C7-C6	-4.56	117.57	126.84
3	C	502	NTD	C8-C7-C6	-4.39	117.91	126.84
3	A	502	NTD	C8-C7-C6	-4.27	118.16	126.84
3	B	502	NTD	C8-C7-C6	-4.21	118.27	126.84
3	D	502	NTD	C3-C2-C5	-2.57	109.06	113.52
3	C	502	NTD	C1-C6-C5	-2.23	116.30	118.85
3	C	502	NTD	C10-C9-C8	2.36	122.91	119.42
3	D	502	NTD	C10-C9-C8	2.44	123.02	119.42
3	B	502	NTD	C10-C9-C8	2.49	123.09	119.42
3	A	502	NTD	C10-C9-C8	2.68	123.37	119.42
3	A	502	NTD	C1-C6-C5	2.78	122.03	118.85
3	B	502	NTD	C1-C6-C5	3.58	122.95	118.85
3	C	502	NTD	O7-C7-C8	3.82	121.08	113.46
3	B	502	NTD	O7-C7-C8	4.10	121.63	113.46
3	D	502	NTD	O7-C7-C8	4.11	121.66	113.46
3	A	502	NTD	O7-C7-C8	4.21	121.86	113.46

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	502	NTD	1	0
3	B	502	NTD	1	0
3	C	502	NTD	2	0
3	D	502	NTD	2	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

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

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

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	367/445 (82%)	-0.87	0 100 100	10, 22, 33, 59	1 (0%)
1	B	370/445 (83%)	-0.81	0 100 100	10, 23, 37, 52	1 (0%)
1	C	365/445 (82%)	-0.86	0 100 100	12, 22, 34, 53	1 (0%)
1	D	370/445 (83%)	-0.84	0 100 100	10, 23, 39, 48	1 (0%)
All	All	1472/1780 (82%)	-0.85	0 100 100	10, 23, 36, 59	4 (0%)

There are no RSRZ outliers to report.

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. 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
3	NTD	D	502	23/23	0.97	0.13	0.52	22,29,37,39	0
3	NTD	C	502	23/23	0.96	0.11	-0.32	26,31,41,44	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	NTD	A	502	23/23	0.98	0.10	-0.75	17,19,21,24	0
3	NTD	B	502	23/23	0.98	0.10	-0.83	23,26,37,39	0
2	FE	B	501	1/1	1.00	0.07	-2.51	11,11,11,11	0
2	FE	D	501	1/1	1.00	0.08	-2.63	14,14,14,14	0
2	FE	A	501	1/1	1.00	0.08	-2.77	20,20,20,20	0
2	FE	C	501	1/1	0.98	0.08	-3.27	19,19,19,19	0

6.5 Other polymers [i](#)

There are no such residues in this entry.