



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:11 AM GMT

PDB ID : 3DOI
Title : Crystal Structure of a Thermostable Esterase complex with paraoxon
Authors : Levisson, M.; Sun, L.; Hendriks, S.; Dijkstra, B.W.; Van der Oost, J.; Kengen, S.W.M.
Deposited on : 2008-07-04
Resolution : 3.00 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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

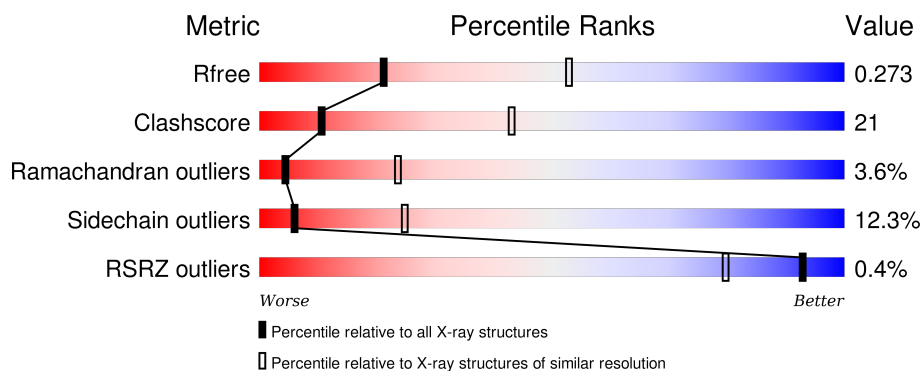
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 3.00 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}	91344	1578 (3.00-3.00)
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (3.00-3.00)

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	380	 54% 34% 8% ..
1	B	380	 61% 30% 7% .

2 Entry composition [i](#)

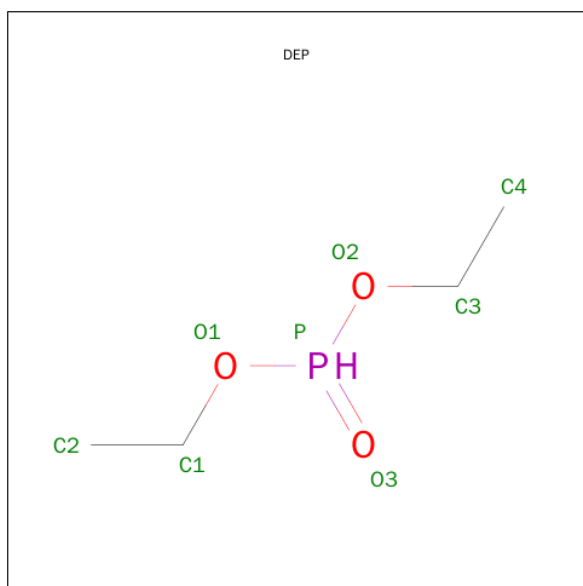
There are 2 unique types of molecules in this entry. The entry contains 5998 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 esterase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	375	Total	C	N	O	S	0	0	0
			3001	1940	493	560	8			
1	B	373	Total	C	N	O	S	0	0	0
			2981	1929	488	556	8			

- Molecule 2 is DIETHYL PHOSPHONATE (three-letter code: DEP) (formula: $C_4H_{11}O_3P$).

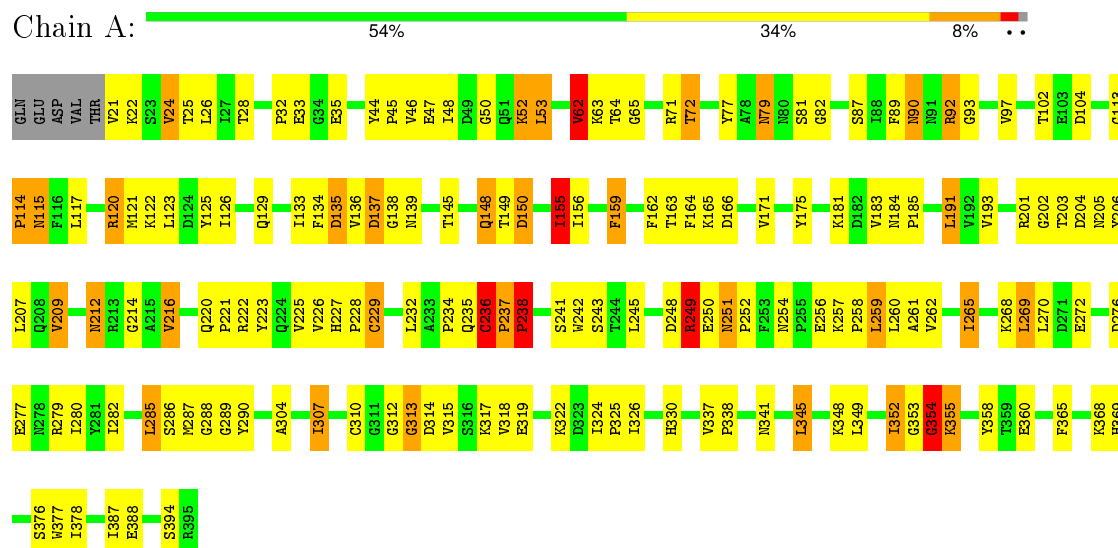


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	P	0	0
			8	4	3	1		
2	B	1	Total	C	O	P	0	0
			8	4	3	1		

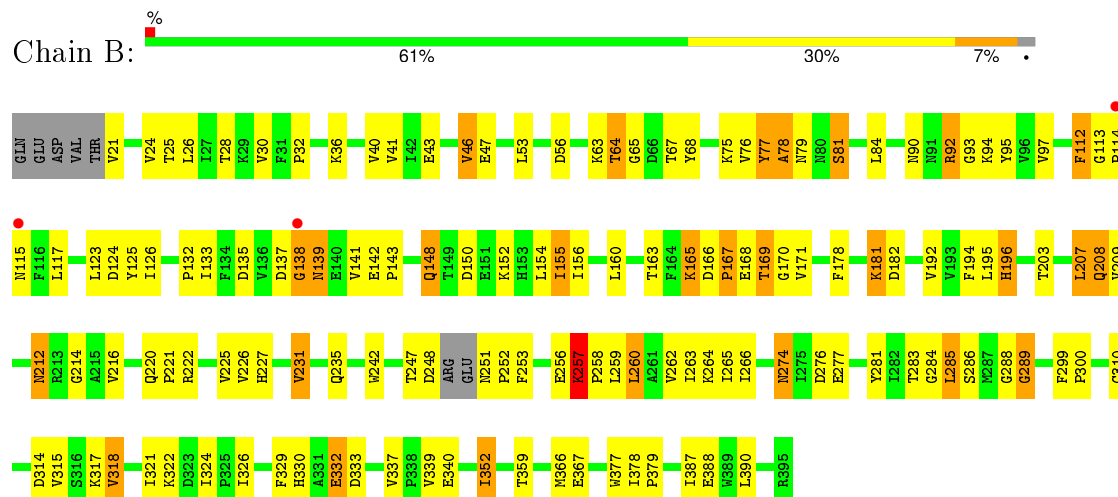
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: esterase



- Molecule 1: esterase



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	130.54Å 130.54Å 304.55Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	37.69 – 3.00 37.68 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.8 (37.69-3.00) 99.8 (37.68-3.00)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.37 (at 3.01Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.220 , 0.266 0.227 , 0.273	Depositor DCC
R_{free} test set	1041 reflections (5.39%)	DCC
Wilson B-factor (Å ²)	52.2	Xtriage
Anisotropy	0.083	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 29.8	EDS
Estimated twinning fraction	0.000 for -h,1/3*h-1/3*k-1/3*l,-4/3*h-8/3*k+1/3*l 0.000 for -1/3*h+1/3*k+1/3*l,-k,8/3*h+4/3*k+1/3*l 0.000 for -2/3*h-1/3*k-1/3*l,-1/3*h-2/3*k+1/3*l,-4/3*h+4/3*k+1/3*l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	1 of 20352 reflections (0.005%)	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5998	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 50.99 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 5.9818e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: DEP

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	0.74	1/3083 (0.0%)	0.76	1/4198 (0.0%)
1	B	0.69	1/3062 (0.0%)	0.75	0/4169
All	All	0.71	2/6145 (0.0%)	0.76	1/8367 (0.0%)

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	A	0	7
1	B	1	9
All	All	1	16

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	307	ILE	CG1-CD1	11.74	2.31	1.50
1	B	367	GLU	CG-CD	5.59	1.60	1.51

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	62	VAL	CB-CA-C	-5.91	100.17	111.40

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	B	168	GLU	CA

All (16) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	114	PRO	Peptide
1	A	120	ARG	Peptide
1	A	137	ASP	Peptide
1	A	228	PRO	Peptide
1	A	236	CYS	Peptide
1	A	354	GLY	Peptide
1	A	64	THR	Peptide
1	B	113	GLY	Peptide
1	B	114	PRO	Peptide
1	B	167	PRO	Peptide
1	B	181	LYS	Peptide
1	B	196	HIS	Peptide
1	B	284	GLY	Peptide
1	B	289	GLY	Peptide
1	B	64	THR	Peptide
1	B	77	TYR	Peptide

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	3001	0	2955	121	0
1	B	2981	0	2935	133	0
2	A	8	0	10	0	0
2	B	8	0	10	2	0
All	All	5998	0	5910	254	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 21.

All (254) 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:30:VAL:CG2	1:B:156:ILE:HD11	1.65	1.26
1:A:285:LEU:HD22	1:A:285:LEU:H	1.03	1.15

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:307:ILE:CD1	1:A:307:ILE:CG1	2.31	1.07
1:B:30:VAL:CG2	1:B:156:ILE:CD1	2.36	1.03
1:B:274:ASN:H	1:B:274:ASN:HD22	1.06	1.00
1:B:169:THR:HG23	1:B:171:VAL:HG12	1.42	1.00
1:B:112:PHE:HD1	1:B:112:PHE:N	1.58	0.98
1:B:155:ILE:HG21	1:B:216:VAL:HG13	1.43	0.96
1:B:285:LEU:HD23	1:B:285:LEU:O	1.67	0.94
1:A:285:LEU:HD22	1:A:285:LEU:N	1.83	0.93
1:B:285:LEU:HD23	1:B:285:LEU:C	1.88	0.93
1:B:166:ASP:HB3	1:B:169:THR:HG23	1.51	0.93
1:A:285:LEU:CD2	1:A:285:LEU:H	1.83	0.92
1:A:220:GLN:NE2	1:A:222:ARG:HH11	1.68	0.91
1:B:30:VAL:HG21	1:B:156:ILE:HD11	1.52	0.91
1:B:30:VAL:HG22	1:B:156:ILE:HD11	1.50	0.91
1:A:72:THR:HG23	1:A:104:ASP:OD1	1.72	0.87
1:B:169:THR:CG2	1:B:171:VAL:HG12	2.04	0.87
1:A:313:GLY:HA2	1:A:345:LEU:HD21	1.55	0.86
1:B:112:PHE:H	1:B:112:PHE:HD1	0.88	0.86
1:A:249:ARG:HG3	1:A:249:ARG:O	1.76	0.85
1:B:112:PHE:CD1	1:B:112:PHE:N	2.30	0.84
1:A:135:ASP:OD1	1:A:138:GLY:HA2	1.78	0.84
1:B:274:ASN:N	1:B:274:ASN:HD22	1.76	0.83
1:B:166:ASP:HB3	1:B:169:THR:CG2	2.09	0.83
1:B:178:PHE:HB3	1:B:231:VAL:HG13	1.62	0.82
1:B:166:ASP:CG	1:B:169:THR:HG22	2.01	0.80
1:A:220:GLN:HE21	1:A:222:ARG:HH11	1.25	0.79
1:B:285:LEU:O	1:B:285:LEU:CD2	2.30	0.79
1:B:30:VAL:HG23	1:B:156:ILE:CD1	2.14	0.77
1:B:330:HIS:CE1	1:B:339:VAL:HA	2.19	0.77
1:A:286:SER:O	1:A:289:GLY:N	2.18	0.77
1:A:24:VAL:CG2	1:A:148:GLN:HA	2.15	0.76
1:B:43:GLU:HG3	1:B:95:TYR:CE1	2.20	0.76
1:A:268:LYS:O	1:A:272:GLU:HG3	1.85	0.75
1:A:114:PRO:O	1:A:115:ASN:HB2	1.86	0.74
1:A:21:VAL:HA	1:A:45:PRO:HD3	1.69	0.74
1:B:24:VAL:HG22	1:B:148:GLN:HA	1.70	0.73
1:A:148:GLN:H	1:A:148:GLN:HE21	1.35	0.72
1:B:310:CYS:HA	1:B:330:HIS:CD2	2.23	0.72
1:B:212:ASN:HD22	1:B:214:GLY:H	1.38	0.71
1:A:155:ILE:O	1:A:216:VAL:HG22	1.89	0.71
1:B:166:ASP:CB	1:B:169:THR:CG2	2.69	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:274:ASN:ND2	1:B:274:ASN:H	1.83	0.70
1:A:125:TYR:H	1:A:148:GLN:HE22	1.38	0.70
1:A:47:GLU:HG2	1:A:93:GLY:HA2	1.72	0.70
1:A:212:ASN:C	1:A:212:ASN:HD22	1.96	0.69
1:A:25:THR:HG23	1:A:150:ASP:HB3	1.73	0.69
1:B:166:ASP:CB	1:B:169:THR:HG23	2.23	0.69
1:B:256:GLU:HG3	1:B:258:PRO:HD2	1.75	0.69
1:A:220:GLN:NE2	1:A:222:ARG:NH1	2.40	0.68
1:A:62:VAL:HG13	1:A:125:TYR:CE1	2.28	0.68
1:A:318:VAL:HG21	1:A:348:LYS:HG3	1.73	0.68
1:B:285:LEU:CD2	1:B:285:LEU:C	2.61	0.68
1:A:290:TYR:HA	1:A:312:GLY:O	1.93	0.68
1:A:262:VAL:HA	1:A:265:ILE:HG23	1.76	0.67
1:B:78:ALA:O	1:B:92:ARG:N	2.27	0.67
1:B:285:LEU:CD1	1:B:377:TRP:CG	2.79	0.66
1:B:285:LEU:HD11	1:B:377:TRP:CG	2.30	0.66
1:A:314:ASP:OD2	1:A:317:LYS:HE3	1.96	0.66
1:A:330:HIS:HE1	1:A:337:VAL:O	1.80	0.65
1:A:212:ASN:HD22	1:A:214:GLY:H	1.45	0.65
1:A:285:LEU:HD23	1:A:285:LEU:O	1.97	0.65
1:A:258:PRO:O	1:A:262:VAL:HG23	1.97	0.64
1:B:154:LEU:O	1:B:156:ILE:N	2.30	0.64
1:A:276:ASP:C	1:A:276:ASP:OD1	2.36	0.64
1:B:314:ASP:OD2	1:B:317:LYS:HE3	1.97	0.64
1:A:79:ASN:HD22	1:A:79:ASN:C	2.02	0.63
1:A:47:GLU:HG2	1:A:93:GLY:CA	2.28	0.63
1:B:148:GLN:H	1:B:148:GLN:HE21	1.48	0.62
1:B:352:ILE:HG23	1:B:352:ILE:O	2.00	0.62
1:B:154:LEU:C	1:B:156:ILE:H	2.04	0.61
1:A:32:PRO:HB3	1:A:378:ILE:HG23	1.82	0.61
1:B:285:LEU:HD11	1:B:377:TRP:CD1	2.36	0.60
1:A:251:ASN:HD21	1:A:254:ASN:ND2	1.99	0.60
1:B:220:GLN:NE2	1:B:222:ARG:HH11	2.00	0.60
1:B:30:VAL:HG23	1:B:156:ILE:HD12	1.83	0.60
1:A:47:GLU:OE1	1:A:92:ARG:HG3	2.01	0.60
1:A:330:HIS:CE1	1:A:337:VAL:O	2.55	0.60
1:A:265:ILE:O	1:A:269:LEU:HD22	2.02	0.59
1:A:279:ARG:HD3	1:A:394:SER:HB2	1.84	0.59
1:A:87:SER:C	1:A:89:PHE:H	2.04	0.59
1:B:169:THR:C	1:B:171:VAL:H	2.06	0.58
1:B:79:ASN:HA	1:B:92:ARG:O	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:247:THR:O	1:B:248:ASP:CB	2.50	0.58
1:B:226:VAL:HG13	1:B:227:HIS:CD2	2.39	0.58
1:B:337:VAL:CG2	2:B:401:DEP:H23	2.34	0.58
1:A:148:GLN:H	1:A:148:GLN:NE2	1.99	0.57
1:A:354:GLY:O	1:A:355:LYS:HB2	2.01	0.57
1:A:338:PRO:O	1:A:341:ASN:HB2	2.03	0.57
1:A:353:GLY:O	1:A:354:GLY:C	2.41	0.57
1:B:257:LYS:HB2	1:B:258:PRO:HD3	1.86	0.57
1:B:281:TYR:HB3	1:B:390:LEU:HD11	1.85	0.57
1:A:202:GLY:HA2	1:A:238:PRO:HG3	1.87	0.56
1:B:166:ASP:CB	1:B:169:THR:HG22	2.36	0.56
1:A:63:LYS:CD	1:A:65:GLY:O	2.54	0.56
1:A:79:ASN:HD21	1:A:82:GLY:H	1.52	0.56
1:B:196:HIS:CD2	1:B:208:GLN:HB3	2.41	0.56
1:A:79:ASN:ND2	1:A:82:GLY:H	2.03	0.55
1:A:226:VAL:HG23	1:A:227:HIS:CE1	2.41	0.55
1:A:72:THR:CG2	1:A:104:ASP:OD1	2.50	0.55
1:A:212:ASN:ND2	1:A:214:GLY:H	2.04	0.55
1:B:63:LYS:HB3	1:B:124:ASP:HB3	1.88	0.55
1:B:257:LYS:CB	1:B:258:PRO:HD3	2.37	0.55
1:B:227:HIS:NE2	1:B:388:GLU:OE2	2.38	0.55
1:B:26:LEU:HD23	1:B:40:VAL:HG22	1.89	0.55
1:B:169:THR:O	1:B:171:VAL:N	2.40	0.55
1:A:285:LEU:CD2	1:A:285:LEU:N	2.50	0.54
1:B:30:VAL:HG22	1:B:156:ILE:CD1	2.24	0.54
1:B:212:ASN:HD22	1:B:212:ASN:C	2.10	0.54
1:A:155:ILE:O	1:A:216:VAL:CG2	2.56	0.54
1:A:191:LEU:HD23	1:A:280:ILE:HG23	1.89	0.54
1:B:155:ILE:CG2	1:B:216:VAL:HG13	2.27	0.54
1:B:166:ASP:CG	1:B:169:THR:CG2	2.74	0.54
1:A:148:GLN:N	1:A:148:GLN:HE21	2.03	0.53
1:A:175:TYR:CD1	1:A:265:ILE:HD12	2.44	0.53
1:B:274:ASN:N	1:B:274:ASN:ND2	2.48	0.53
1:A:90:ASN:N	1:A:90:ASN:OD1	2.42	0.53
1:B:166:ASP:OD2	1:B:169:THR:HG22	2.09	0.52
1:B:262:VAL:HA	1:B:265:ILE:HD12	1.90	0.52
1:B:46:VAL:HG13	1:B:47:GLU:N	2.25	0.52
1:B:94:LYS:HD2	1:B:95:TYR:CE2	2.45	0.52
1:A:24:VAL:HG23	1:A:148:GLN:HA	1.92	0.52
1:B:332:GLU:HG3	1:B:333:ASP:N	2.24	0.51
1:A:220:GLN:HE21	1:A:222:ARG:NH1	2.02	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:208:GLN:HG2	1:B:235:GLN:HB2	1.92	0.51
1:B:337:VAL:HG22	2:B:401:DEP:H23	1.93	0.51
1:A:354:GLY:O	1:A:355:LYS:CB	2.58	0.51
1:B:171:VAL:O	1:B:171:VAL:HG13	2.10	0.51
1:B:256:GLU:HG3	1:B:258:PRO:CD	2.40	0.51
1:B:340:GLU:HA	1:B:340:GLU:OE1	2.11	0.50
1:A:162:PHE:HB2	1:A:175:TYR:CZ	2.46	0.50
1:A:77:TYR:CZ	1:A:97:VAL:HG11	2.47	0.50
1:B:285:LEU:HG	1:B:377:TRP:CE3	2.47	0.50
1:A:242:TRP:O	1:A:259:LEU:HB2	2.11	0.50
1:B:148:GLN:H	1:B:148:GLN:NE2	2.09	0.49
1:B:46:VAL:O	1:B:94:LYS:HA	2.13	0.49
1:A:324:ILE:O	1:A:326:ILE:HG13	2.12	0.49
1:A:87:SER:C	1:A:89:PHE:N	2.66	0.49
1:A:353:GLY:O	1:A:354:GLY:O	2.30	0.49
1:A:48:ILE:HD13	1:A:133:ILE:HD12	1.94	0.49
1:B:315:VAL:HG22	1:B:315:VAL:O	2.11	0.49
1:B:252:PRO:HG2	1:B:253:PHE:CE1	2.48	0.49
1:A:63:LYS:HD2	1:A:65:GLY:O	2.12	0.49
1:A:319:GLU:OE2	1:A:322:LYS:HE2	2.12	0.49
1:B:78:ALA:O	1:B:92:ARG:O	2.31	0.48
1:B:125:TYR:H	1:B:148:GLN:HE22	1.62	0.48
1:A:63:LYS:HD3	1:A:65:GLY:O	2.14	0.48
1:A:33:GLU:OE1	1:A:33:GLU:N	2.45	0.48
1:A:156:ILE:HG12	1:A:209:VAL:HG13	1.96	0.48
1:A:193:VAL:HG22	1:A:282:ILE:HG13	1.95	0.47
1:A:201:ARG:HD3	1:A:287:MET:SD	2.54	0.47
1:B:324:ILE:O	1:B:326:ILE:HG13	2.14	0.47
1:B:53:LEU:HD23	1:B:76:VAL:HG12	1.97	0.47
1:A:241:SER:OG	1:A:243:SER:O	2.29	0.47
1:A:212:ASN:C	1:A:212:ASN:ND2	2.66	0.47
1:B:256:GLU:O	1:B:257:LYS:C	2.52	0.47
1:A:226:VAL:HG23	1:A:227:HIS:ND1	2.30	0.47
1:A:205:ASN:N	1:A:235:GLN:OE1	2.45	0.47
1:B:169:THR:C	1:B:171:VAL:N	2.68	0.47
1:A:44:TYR:CD2	1:A:48:ILE:HD12	2.50	0.47
1:A:28:THR:HB	1:A:35:GLU:HG3	1.97	0.47
1:B:25:THR:HG23	1:B:150:ASP:HB3	1.97	0.47
1:B:195:LEU:O	1:B:288:GLY:HA3	2.15	0.47
1:A:358:TYR:CZ	1:A:360:GLU:HB2	2.50	0.47
1:B:192:VAL:O	1:B:231:VAL:HA	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:134:PHE:HA	1:A:139:ASN:O	2.15	0.46
1:B:132:PRO:HA	1:B:141:VAL:O	2.15	0.46
1:B:30:VAL:CG2	1:B:156:ILE:HD12	2.36	0.46
1:B:212:ASN:ND2	1:B:214:GLY:H	2.10	0.46
1:A:204:ASP:OD2	1:A:206:TYR:HB3	2.16	0.46
1:B:135:ASP:OD1	1:B:138:GLY:HA2	2.16	0.46
1:B:352:ILE:CG2	1:B:352:ILE:O	2.63	0.46
1:A:242:TRP:CD1	1:A:288:GLY:HA2	2.50	0.46
1:B:315:VAL:O	1:B:315:VAL:CG2	2.63	0.46
1:A:159:PHE:CD1	1:A:159:PHE:N	2.84	0.46
1:B:194:PHE:HA	1:B:283:THR:O	2.16	0.46
1:B:68:TYR:CE1	1:B:126:ILE:HG22	2.50	0.45
1:B:166:ASP:HB3	1:B:169:THR:HG22	1.94	0.45
1:A:277:GLU:H	1:A:277:GLU:CD	2.20	0.45
1:A:92:ARG:HB3	1:A:92:ARG:NH1	2.30	0.45
1:B:276:ASP:OD1	1:B:276:ASP:C	2.54	0.45
1:B:154:LEU:C	1:B:156:ILE:N	2.70	0.45
1:B:166:ASP:HA	1:B:167:PRO:HD3	1.41	0.45
1:B:329:PHE:CD1	1:B:359:THR:HB	2.52	0.45
1:A:135:ASP:OD1	1:A:138:GLY:CA	2.59	0.45
1:B:32:PRO:HB3	1:B:378:ILE:HG23	1.97	0.45
1:B:78:ALA:HB1	1:B:92:ARG:HA	1.99	0.44
1:A:352:ILE:HA	1:A:352:ILE:HD13	1.83	0.44
1:B:366:MET:CE	1:B:379:PRO:HG3	2.47	0.44
1:A:248:ASP:C	1:A:250:GLU:H	2.21	0.44
1:A:285:LEU:HD12	1:A:377:TRP:CG	2.52	0.44
1:B:259:LEU:O	1:B:259:LEU:HD12	2.17	0.44
1:B:166:ASP:OD2	1:B:169:THR:CG2	2.66	0.44
1:A:249:ARG:O	1:A:249:ARG:CG	2.55	0.44
1:B:220:GLN:NE2	1:B:222:ARG:NH1	2.65	0.44
1:B:166:ASP:OD1	1:B:264:LYS:NZ	2.46	0.44
1:B:387:ILE:HD12	1:B:387:ILE:H	1.82	0.44
1:A:171:VAL:HG21	1:A:257:LYS:HB3	2.00	0.44
1:A:50:GLY:HA2	1:A:53:LEU:HB2	1.99	0.44
1:B:165:LYS:O	1:B:167:PRO:HD3	2.18	0.43
1:B:169:THR:CG2	1:B:171:VAL:CG1	2.88	0.43
1:A:62:VAL:HG13	1:A:125:TYR:CZ	2.52	0.43
1:A:129:GLN:NE2	1:A:133:ILE:HG12	2.33	0.43
1:A:270:LEU:HD21	1:A:280:ILE:HD11	2.00	0.43
1:B:242:TRP:CD1	1:B:288:GLY:HA2	2.53	0.43
1:B:263:ILE:HA	1:B:266:ILE:HD12	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:164:PHE:CE2	1:A:166:ASP:HA	2.54	0.43
1:B:133:ILE:O	1:B:141:VAL:N	2.47	0.43
1:B:285:LEU:HG	1:B:377:TRP:CD2	2.54	0.43
1:B:286:SER:O	1:B:289:GLY:N	2.44	0.43
1:A:260:LEU:O	1:A:261:ALA:C	2.54	0.43
1:B:41:VAL:HG22	1:B:97:VAL:HG22	2.00	0.43
1:A:304:ALA:O	1:A:325:PRO:HD2	2.18	0.42
1:A:232:LEU:HG	1:A:234:PRO:HD3	2.00	0.42
1:A:365:PHE:O	1:A:369:HIS:HD2	2.01	0.42
1:A:48:ILE:O	1:A:92:ARG:HA	2.19	0.42
1:B:148:GLN:N	1:B:148:GLN:HE21	2.13	0.42
1:A:50:GLY:HA2	1:A:53:LEU:HD22	2.01	0.42
1:B:138:GLY:HA2	1:B:139:ASN:O	2.20	0.42
1:B:64:THR:HA	1:B:65:GLY:HA2	1.77	0.42
1:A:26:LEU:HD23	1:A:123:LEU:HD13	2.00	0.42
1:A:245:LEU:HA	1:A:252:PRO:CB	2.50	0.42
1:B:28:THR:HA	1:B:36:LYS:O	2.19	0.42
1:A:62:VAL:HG22	1:A:71:ARG:HG2	2.00	0.42
1:A:352:ILE:O	1:A:352:ILE:HG23	2.19	0.42
1:A:126:ILE:HD11	1:A:145:THR:CG2	2.49	0.42
1:A:184:ASN:HA	1:A:185:PRO:HD3	1.92	0.42
1:A:236:CYS:HA	1:A:237:PRO:HD3	1.78	0.42
1:A:256:GLU:HG3	1:A:257:LYS:N	2.35	0.42
1:B:212:ASN:HD22	1:B:214:GLY:N	2.13	0.42
1:A:71:ARG:NE	1:A:104:ASP:OD2	2.39	0.41
1:A:155:ILE:O	1:A:155:ILE:HG22	2.20	0.41
1:B:227:HIS:HE2	1:B:388:GLU:CD	2.24	0.41
1:B:135:ASP:OD1	1:B:138:GLY:N	2.50	0.41
1:B:366:MET:HE1	1:B:379:PRO:HG3	2.02	0.41
1:B:221:PRO:O	1:B:225:VAL:HG22	2.20	0.41
1:B:299:PHE:N	1:B:300:PRO:HD3	2.35	0.41
1:A:223:TYR:CE2	1:A:387:ILE:HD13	2.55	0.41
1:B:329:PHE:CE1	1:B:359:THR:HB	2.55	0.41
1:B:318:VAL:O	1:B:321:ILE:HG13	2.21	0.41
1:A:90:ASN:O	1:A:92:ARG:N	2.46	0.41
1:B:260:LEU:HD12	1:B:260:LEU:HA	1.77	0.41
1:B:208:GLN:CG	1:B:235:GLN:HB2	2.51	0.41
1:B:207:LEU:HD12	1:B:207:LEU:HA	1.98	0.41
1:B:142:GLU:HA	1:B:143:PRO:HD3	1.99	0.41
1:B:212:ASN:ND2	1:B:212:ASN:C	2.74	0.40
1:B:359:THR:HG22	1:B:359:THR:O	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:117:LEU:HD12	1:A:117:LEU:HA	1.81	0.40
1:B:257:LYS:CB	1:B:258:PRO:CD	2.99	0.40
1:A:50:GLY:C	1:A:52:LYS:N	2.75	0.40
1:A:24:VAL:HG22	1:A:148:GLN:HA	2.00	0.40
1:B:330:HIS:HE1	1:B:337:VAL:O	2.04	0.40
1:A:227:HIS:HE1	1:A:388:GLU:OE1	2.04	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	373/380 (98%)	313 (84%)	47 (13%)	13 (4%)	4	24
1	B	369/380 (97%)	311 (84%)	44 (12%)	14 (4%)	4	22
All	All	742/760 (98%)	624 (84%)	91 (12%)	27 (4%)	4	24

All (27) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	121	MET
1	A	229	CYS
1	A	354	GLY
1	A	355	LYS
1	B	81	SER
1	B	168	GLU
1	A	113	GLY
1	B	115	ASN
1	B	138	GLY
1	B	155	ILE
1	A	238	PRO
1	A	313	GLY

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Mol	Chain	Res	Type
1	B	77	TYR
1	B	78	ALA
1	B	170	GLY
1	A	115	ASN
1	A	310	CYS
1	A	368	LYS
1	B	56	ASP
1	B	139	ASN
1	A	249	ARG
1	B	84	LEU
1	B	152	LYS
1	B	257	LYS
1	A	237	PRO
1	B	93	GLY
1	A	155	ILE

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	333/338 (98%)	285 (86%)	48 (14%)	4	18
1	B	331/338 (98%)	297 (90%)	34 (10%)	9	33
All	All	664/676 (98%)	582 (88%)	82 (12%)	6	25

All (82) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	LYS
1	A	24	VAL
1	A	46	VAL
1	A	52	LYS
1	A	53	LEU
1	A	62	VAL
1	A	72	THR
1	A	79	ASN

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Mol	Chain	Res	Type
1	A	81	SER
1	A	90	ASN
1	A	92	ARG
1	A	102	THR
1	A	120	ARG
1	A	122	LYS
1	A	135	ASP
1	A	136	VAL
1	A	137	ASP
1	A	148	GLN
1	A	149	THR
1	A	150	ASP
1	A	155	ILE
1	A	159	PHE
1	A	163	THR
1	A	165	LYS
1	A	181	LYS
1	A	183	VAL
1	A	191	LEU
1	A	203	THR
1	A	207	LEU
1	A	209	VAL
1	A	212	ASN
1	A	216	VAL
1	A	221	PRO
1	A	225	VAL
1	A	229	CYS
1	A	236	CYS
1	A	238	PRO
1	A	249	ARG
1	A	251	ASN
1	A	259	LEU
1	A	265	ILE
1	A	269	LEU
1	A	285	LEU
1	A	315	VAL
1	A	345	LEU
1	A	349	LEU
1	A	352	ILE
1	A	376	SER
1	B	21	VAL
1	B	46	VAL

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Mol	Chain	Res	Type
1	B	67	THR
1	B	75	LYS
1	B	81	SER
1	B	90	ASN
1	B	92	ARG
1	B	112	PHE
1	B	117	LEU
1	B	123	LEU
1	B	137	ASP
1	B	148	GLN
1	B	160	LEU
1	B	163	THR
1	B	165	LYS
1	B	169	THR
1	B	181	LYS
1	B	182	ASP
1	B	203	THR
1	B	207	LEU
1	B	208	GLN
1	B	209	VAL
1	B	212	ASN
1	B	231	VAL
1	B	251	ASN
1	B	257	LYS
1	B	260	LEU
1	B	274	ASN
1	B	277	GLU
1	B	285	LEU
1	B	318	VAL
1	B	322	LYS
1	B	332	GLU
1	B	352	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	79	ASN
1	A	148	GLN
1	A	212	ASN
1	A	220	GLN
1	A	227	HIS
1	A	254	ASN

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Mol	Chain	Res	Type
1	A	278	ASN
1	A	330	HIS
1	A	369	HIS
1	B	106	HIS
1	B	148	GLN
1	B	208	GLN
1	B	212	ASN
1	B	220	GLN
1	B	254	ASN
1	B	274	ASN
1	B	330	HIS
1	B	341	ASN

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](#)

2 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	DEP	A	401	-	2,7,7	0.88	0	2,7,7	0.80	0
2	DEP	B	401	-	2,7,7	0.71	0	2,7,7	0.23	0

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	DEP	A	401	-	-	0/2/6/6	0/0/0/0
2	DEP	B	401	-	-	0/2/6/6	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	401	DEP	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 [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	375/380 (98%)	-0.36	0	100 100	32, 44, 64, 74	0
1	B	373/380 (98%)	-0.37	3 (0%)	87 67	30, 44, 63, 74	0
All	All	748/760 (98%)	-0.37	3 (0%)	93 80	30, 44, 64, 74	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	114	PRO	3.0
1	B	115	ASN	2.6
1	B	138	GLY	2.2

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(\AA^2)	Q<0.9
2	DEP	A	401	8/8	0.91	0.26	1.06	45,47,48,48	0
2	DEP	B	401	8/8	0.98	0.15	-0.56	36,37,44,45	0

6.5 Other polymers [i](#)

There are no such residues in this entry.