



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:59 PM GMT

PDB ID : 1W52
Title : CRYSTAL STRUCTURE OF A PROTEOLYZED FORM OF PANCREATIC LIPASE RELATED PROTEIN 2 FROM HORSE
Authors : Mancheno, J.M.; Jayne, S.; Kerfelec, B.; Chapus, C.; Crenon, I.; Hermoso, J.A.
Deposited on : 2004-08-04
Resolution : 2.99 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

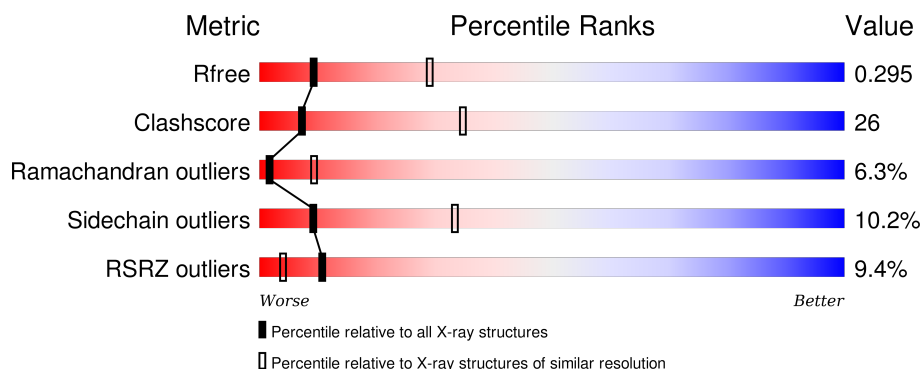
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.99 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	X	452	<div> <div>9%</div> <div>53%</div> <div>37%</div> <div>8%</div> <div>..</div> </div>

2 Entry composition [i](#)

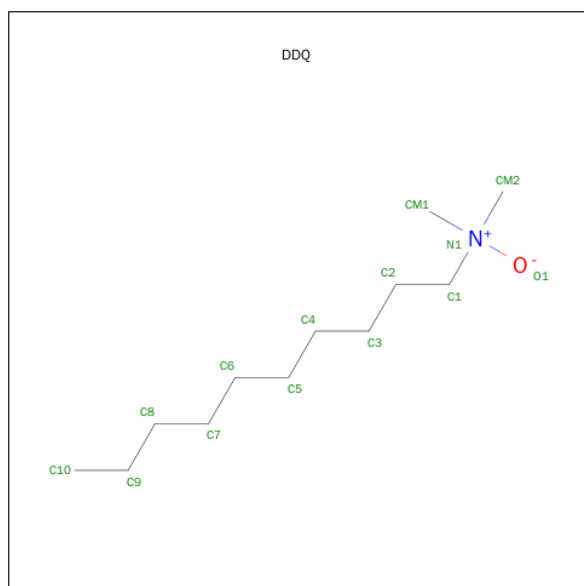
There are 3 unique types of molecules in this entry. The entry contains 3524 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PANCREATIC LIPASE RELATED PROTEIN 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	X	448	Total	C	N	O	S	0	0	0
			3494	2199	595	681	19			

- Molecule 2 is DECYLAMINE-N,N-DIMETHYL-N-OXIDE (three-letter code: DDQ) (formula: C₁₂H₂₇NO).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	X	1	Total	C	N	O	0	0
			14	12	1	1		
2	X	1	Total	C	N	O	0	0
			14	12	1	1		

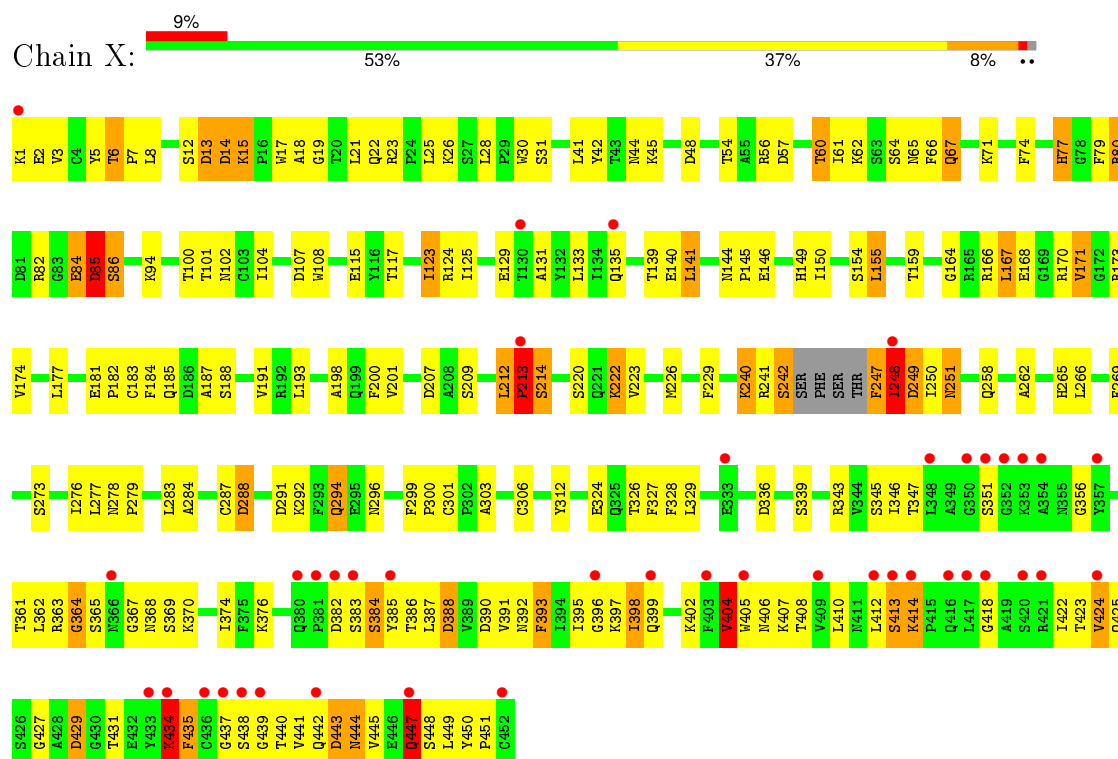
- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	X	2	Total	Ca	0	0
			2	2		

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 ($\text{RSRZ} > 2$). 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: PANCREATIC LIPASE RELATED PROTEIN 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	128.43 Å 128.43 Å 85.82 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	26.17 – 2.99 26.13 – 2.99	Depositor EDS
% Data completeness (in resolution range)	100.0 (26.17-2.99) 100.0 (26.13-2.99)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.54 (at 3.00 Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.231 , 0.294 0.240 , 0.295	Depositor DCC
R_{free} test set	847 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	45.2	Xtriage
Anisotropy	0.003	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 54.0	EDS
Estimated twinning fraction	0.035 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 16808 reflections	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	3524	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.59% 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, 0.375 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: DDQ, CA

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	X	0.93	2/3579 (0.1%)	1.06	14/4848 (0.3%)

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	X	0	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	X	404	VAL	C-N	-5.84	1.20	1.34
1	X	213	PRO	CB-CG	5.03	1.75	1.50

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	429	ASP	CB-CG-OD2	9.48	126.83	118.30
1	X	388	ASP	CB-CG-OD2	9.14	126.53	118.30
1	X	107	ASP	CB-CG-OD2	8.09	125.58	118.30
1	X	80	ARG	NE-CZ-NH2	-7.59	116.50	120.30
1	X	390	ASP	CB-CG-OD2	7.48	125.04	118.30
1	X	85	ASP	O-C-N	-7.34	110.95	122.70
1	X	336	ASP	CB-CG-OD2	6.78	124.41	118.30
1	X	404	VAL	O-C-N	-6.06	113.01	122.70
1	X	288	ASP	CB-CG-OD2	6.03	123.72	118.30
1	X	14	ASP	CB-CG-OD2	5.61	123.35	118.30
1	X	48	ASP	CB-CG-OD2	5.41	123.17	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	291	ASP	CB-CG-OD2	5.38	123.14	118.30
1	X	207	ASP	CB-CG-OD2	5.21	122.99	118.30
1	X	443	ASP	CB-CG-OD2	5.11	122.89	118.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	X	212	LEU	Peptide
1	X	85	ASP	Mainchain

5.2 Too-close contacts [i](#)

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	X	3494	0	3342	179	0
2	X	28	0	54	1	0
3	X	2	0	0	0	0
All	All	3524	0	3396	180	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 26.

All (180) 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:X:213:PRO:CB	1:X:213:PRO:CG	1.75	1.49
1:X:84:GLU:O	1:X:85:ASP:HB2	1.34	1.14
1:X:423:THR:HG23	1:X:434:LYS:HD3	1.45	0.96
1:X:115:GLU:OE1	1:X:117:THR:HB	1.69	0.91
1:X:17:TRP:CE3	1:X:124:ARG:HD3	2.14	0.83
1:X:84:GLU:O	1:X:85:ASP:CB	2.17	0.79
1:X:345:SER:CB	1:X:386:THR:HG22	2.15	0.77
1:X:167:LEU:HD23	1:X:167:LEU:N	2.00	0.77
1:X:399:GLN:HA	1:X:435:PHE:HE1	1.48	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:361:THR:O	1:X:362:LEU:HD23	1.86	0.76
1:X:346:ILE:HG12	1:X:422:ILE:HG12	1.66	0.76
1:X:94:LYS:HE2	1:X:273:SER:OG	1.85	0.76
1:X:222:LYS:NZ	1:X:326:THR:HG21	2.02	0.74
1:X:6:THR:HB	1:X:7:PRO:HD3	1.69	0.74
1:X:345:SER:OG	1:X:386:THR:HG22	1.89	0.73
1:X:144:ASN:ND2	1:X:146:GLU:HG2	2.04	0.73
1:X:276:ILE:HG22	1:X:277:LEU:HD23	1.70	0.71
1:X:54:THR:HG22	1:X:56:ARG:H	1.56	0.71
1:X:212:LEU:HB3	1:X:213:PRO:HD2	1.71	0.70
1:X:398:ILE:HD12	1:X:424:VAL:HG11	1.73	0.70
1:X:447:GLN:HG2	1:X:447:GLN:O	1.91	0.69
1:X:294:GLN:NE2	1:X:294:GLN:HA	2.07	0.69
1:X:347:THR:HA	1:X:384:SER:HA	1.73	0.68
1:X:423:THR:CG2	1:X:434:LYS:HD3	2.22	0.68
1:X:57:ASP:O	1:X:60:THR:OG1	2.11	0.68
1:X:345:SER:CA	1:X:386:THR:HG22	2.23	0.68
1:X:181:GLU:HB3	1:X:182:PRO:HD3	1.78	0.66
1:X:17:TRP:CZ3	1:X:124:ARG:HD3	2.31	0.66
1:X:398:ILE:HD12	1:X:424:VAL:CG1	2.26	0.66
1:X:247:PHE:N	1:X:247:PHE:CD2	2.63	0.65
1:X:144:ASN:HD21	1:X:146:GLU:HG2	1.62	0.65
1:X:279:PRO:HB3	1:X:312:TYR:CD1	2.33	0.63
1:X:423:THR:HA	1:X:434:LYS:HB3	1.80	0.63
1:X:247:PHE:N	1:X:247:PHE:HD2	1.96	0.63
1:X:345:SER:HA	1:X:386:THR:HG22	1.80	0.62
1:X:149:HIS:CD2	1:X:173:ARG:HD3	2.35	0.61
1:X:79:PHE:HB2	1:X:155:LEU:HD22	1.83	0.61
1:X:30:TRP:O	1:X:125:ILE:HD11	2.01	0.61
1:X:362:LEU:O	1:X:368:ASN:HA	2.00	0.61
1:X:393:PHE:CZ	1:X:395:ILE:CG2	2.84	0.61
1:X:131:ALA:HB1	1:X:167:LEU:HD21	1.83	0.60
1:X:363:ARG:O	1:X:367:GLY:O	2.19	0.60
1:X:6:THR:CB	1:X:7:PRO:HD3	2.32	0.60
1:X:6:THR:CB	1:X:7:PRO:CD	2.79	0.60
1:X:213:PRO:O	1:X:214:SER:HB2	2.02	0.59
1:X:141:LEU:N	1:X:141:LEU:HD23	2.16	0.59
1:X:251:ASN:OD1	1:X:258:GLN:HB2	2.03	0.59
1:X:345:SER:HB2	1:X:423:THR:HB	1.86	0.58
1:X:184:PHE:O	1:X:185:GLN:C	2.41	0.58
1:X:45:LYS:HB2	1:X:65:ASN:HD22	1.69	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:384:SER:O	1:X:385:TYR:CD2	2.57	0.57
1:X:212:LEU:HB3	1:X:213:PRO:CD	2.34	0.57
1:X:80:ARG:HD3	1:X:247:PHE:HB3	1.87	0.56
1:X:45:LYS:CB	1:X:65:ASN:HD22	2.16	0.56
1:X:188:SER:O	1:X:191:VAL:HG22	2.05	0.56
1:X:437:GLY:CA	1:X:449:LEU:HG	2.35	0.56
1:X:284:ALA:HB1	1:X:327:PHE:HB3	1.86	0.56
1:X:294:GLN:HE21	1:X:294:GLN:CA	2.17	0.56
1:X:133:LEU:HD12	1:X:133:LEU:O	2.06	0.56
1:X:77:HIS:CE1	1:X:82:ARG:HA	2.41	0.55
1:X:402:LYS:NZ	1:X:448:SER:OG	2.26	0.55
1:X:345:SER:OG	1:X:386:THR:CG2	2.54	0.55
1:X:284:ALA:HB1	1:X:327:PHE:CB	2.36	0.55
1:X:100:THR:C	1:X:101:THR:OG1	2.44	0.55
1:X:61:ILE:O	1:X:64:SER:OG	2.12	0.55
1:X:166:ARG:C	1:X:167:LEU:HD23	2.27	0.54
1:X:14:ASP:O	1:X:15:LYS:O	2.25	0.54
1:X:300:PRO:HB2	1:X:427:GLY:O	2.08	0.54
1:X:405:TRP:NE1	1:X:443:ASP:HA	2.22	0.54
1:X:404:VAL:O	1:X:405:TRP:HB3	2.08	0.54
1:X:296:ASN:HA	1:X:388:ASP:OD2	2.07	0.54
1:X:77:HIS:CD2	1:X:77:HIS:C	2.81	0.53
1:X:123:ILE:HD12	1:X:155:LEU:HG	1.91	0.53
1:X:242:SER:O	1:X:242:SER:OG	2.20	0.53
1:X:435:PHE:CD2	1:X:435:PHE:N	2.76	0.53
1:X:391:VAL:O	1:X:393:PHE:N	2.42	0.53
1:X:361:THR:HG22	1:X:362:LEU:H	1.74	0.53
1:X:424:VAL:CG1	1:X:425:GLN:N	2.72	0.52
1:X:17:TRP:CD2	1:X:124:ARG:NH1	2.78	0.52
1:X:167:LEU:CD2	1:X:167:LEU:N	2.72	0.52
1:X:418:GLY:HA2	1:X:447:GLN:HE22	1.75	0.52
1:X:213:PRO:O	1:X:214:SER:CB	2.58	0.51
1:X:443:ASP:O	1:X:445:VAL:HG23	2.10	0.51
1:X:283:LEU:HD23	1:X:306:CYS:HB2	1.92	0.51
1:X:222:LYS:HZ1	1:X:326:THR:HG21	1.74	0.51
1:X:398:ILE:HG22	1:X:399:GLN:N	2.26	0.51
1:X:177:LEU:HD13	1:X:269:PHE:HA	1.92	0.51
1:X:150:ILE:O	1:X:174:VAL:HA	2.11	0.51
1:X:327:PHE:N	1:X:327:PHE:CD1	2.78	0.50
1:X:364:GLY:O	1:X:397:LYS:O	2.30	0.50
1:X:266:LEU:HD22	1:X:269:PHE:HE2	1.76	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:8:LEU:HD11	1:X:129:GLU:OE2	2.10	0.50
1:X:391:VAL:C	1:X:393:PHE:H	2.16	0.50
1:X:74:PHE:CE2	1:X:104:ILE:HD12	2.46	0.49
1:X:140:GLU:CB	1:X:141:LEU:HD23	2.43	0.49
1:X:424:VAL:HG13	1:X:425:GLN:N	2.26	0.49
1:X:383:SER:O	1:X:384:SER:HB3	2.11	0.49
1:X:249:ASP:OD1	1:X:251:ASN:HB2	2.13	0.49
1:X:441:VAL:HG21	1:X:445:VAL:HG11	1.94	0.49
1:X:67:GLN:HG2	1:X:102:ASN:OD1	2.13	0.49
1:X:294:GLN:HE21	1:X:294:GLN:HA	1.72	0.48
1:X:41:LEU:HD21	1:X:66:PHE:HB2	1.95	0.48
1:X:435:PHE:HB2	1:X:449:LEU:HB3	1.94	0.48
1:X:422:ILE:HG13	1:X:449:LEU:HD22	1.95	0.48
1:X:262:ALA:O	1:X:266:LEU:HG	2.13	0.48
1:X:1:LYS:O	1:X:28:LEU:HD11	2.14	0.48
1:X:406:ASN:ND2	1:X:444:ASN:HD21	2.12	0.48
1:X:324:GLU:O	1:X:326:THR:HG23	2.14	0.48
1:X:77:HIS:CD2	1:X:108:TRP:NE1	2.82	0.48
1:X:13:ASP:HB3	1:X:28:LEU:HD21	1.96	0.47
1:X:443:ASP:OD1	1:X:444:ASN:HB2	2.14	0.47
1:X:413:SER:O	1:X:414:LYS:HB2	2.14	0.47
1:X:17:TRP:CG	1:X:124:ARG:NH1	2.83	0.47
1:X:277:LEU:N	1:X:277:LEU:HD23	2.30	0.47
1:X:30:TRP:O	1:X:125:ILE:CD1	2.63	0.47
1:X:140:GLU:C	1:X:141:LEU:HD23	2.35	0.47
1:X:71:LYS:O	1:X:101:THR:CG2	2.62	0.47
1:X:2:GLU:HB3	1:X:12:SER:HA	1.96	0.47
1:X:164:GLY:HA2	1:X:171:VAL:HG11	1.97	0.47
1:X:329:LEU:N	1:X:329:LEU:HD23	2.29	0.47
1:X:22:GLN:O	1:X:183:CYS:HA	2.14	0.47
1:X:402:LYS:HD3	1:X:402:LYS:HA	1.72	0.46
1:X:229:PHE:HB2	1:X:328:PHE:CD2	2.50	0.46
1:X:41:LEU:HD12	1:X:42:TYR:N	2.30	0.46
1:X:345:SER:CB	1:X:386:THR:CG2	2.91	0.46
1:X:182:PRO:O	1:X:183:CYS:HB2	2.15	0.46
1:X:149:HIS:HA	1:X:173:ARG:O	2.15	0.46
1:X:247:PHE:N	1:X:250:ILE:HG13	2.30	0.46
1:X:395:ILE:HG13	1:X:396:GLY:H	1.82	0.45
1:X:6:THR:OG1	1:X:7:PRO:HD2	2.16	0.45
1:X:5:TYR:CZ	1:X:31:SER:HA	2.52	0.45
1:X:14:ASP:O	1:X:15:LYS:C	2.55	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:374:ILE:HG23	1:X:387:LEU:HB3	1.97	0.45
1:X:240:LYS:HE2	1:X:240:LYS:HB3	1.81	0.45
1:X:404:VAL:HG12	1:X:405:TRP:N	2.31	0.45
1:X:441:VAL:CG2	1:X:445:VAL:HG21	2.47	0.45
1:X:181:GLU:HG3	1:X:185:GLN:HB2	2.00	0.44
1:X:299:PHE:HB3	1:X:343:ARG:NH2	2.32	0.44
1:X:356:GLY:O	1:X:376:LYS:HA	2.17	0.44
1:X:200:PHE:HZ	1:X:226:MET:HE3	1.82	0.44
1:X:399:GLN:HA	1:X:435:PHE:CE1	2.39	0.44
1:X:450:TYR:HB3	1:X:451:PRO:HD2	2.00	0.44
1:X:170:ARG:HD3	1:X:170:ARG:HA	1.81	0.44
1:X:278:ASN:N	1:X:279:PRO:CD	2.81	0.43
1:X:328:PHE:O	1:X:329:LEU:HB3	2.17	0.43
1:X:200:PHE:HZ	1:X:226:MET:CE	2.31	0.43
1:X:44:ASN:HD21	1:X:100:THR:HA	1.83	0.43
1:X:94:LYS:CE	1:X:273:SER:OG	2.60	0.43
1:X:22:GLN:HG2	1:X:187:ALA:HB2	2.00	0.43
1:X:22:GLN:C	1:X:23:ARG:HG3	2.39	0.43
1:X:173:ARG:HA	1:X:200:PHE:O	2.18	0.43
1:X:77:HIS:NE2	1:X:108:TRP:CD1	2.87	0.43
1:X:294:GLN:NE2	1:X:294:GLN:CA	2.71	0.43
1:X:71:LYS:O	1:X:101:THR:HG22	2.19	0.43
1:X:54:THR:HG22	1:X:56:ARG:HG3	2.00	0.43
1:X:397:LYS:O	1:X:397:LYS:HG3	2.19	0.43
1:X:222:LYS:HZ3	1:X:326:THR:HG21	1.81	0.42
1:X:363:ARG:NH2	1:X:402:LYS:HE3	2.34	0.42
1:X:300:PRO:O	1:X:301:CYS:C	2.57	0.42
1:X:181:GLU:HB3	1:X:182:PRO:CD	2.48	0.42
1:X:123:ILE:HG12	1:X:159:THR:OG1	2.19	0.42
1:X:393:PHE:CZ	1:X:395:ILE:HG23	2.55	0.42
1:X:3:VAL:HG23	1:X:3:VAL:O	2.20	0.42
1:X:19:GLY:HA3	1:X:26:LYS:HE2	2.02	0.42
1:X:154:SER:HB2	1:X:265:HIS:NE2	2.35	0.42
2:X:502:DDQ:H22	2:X:502:DDQ:HM22	1.81	0.42
1:X:18:ALA:HB1	1:X:26:LYS:HA	2.01	0.42
1:X:248:ILE:O	1:X:248:ILE:HG23	2.20	0.42
1:X:185:GLN:NE2	1:X:220:SER:OG	2.49	0.42
1:X:100:THR:C	1:X:101:THR:HG1	2.21	0.42
1:X:174:VAL:HG12	1:X:198:ALA:CB	2.49	0.42
1:X:288:ASP:OD2	1:X:292:LYS:NZ	2.36	0.42
1:X:174:VAL:O	1:X:201:VAL:HA	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:174:VAL:HG12	1:X:198:ALA:HB2	2.03	0.41
1:X:135:GLN:O	1:X:139:THR:HG23	2.21	0.41
1:X:402:LYS:HZ2	1:X:448:SER:CB	2.29	0.41
1:X:393:PHE:CZ	1:X:395:ILE:HG22	2.55	0.41
1:X:44:ASN:ND2	1:X:100:THR:HA	2.36	0.41
1:X:418:GLY:HA3	1:X:440:THR:HB	2.03	0.40
1:X:193:LEU:O	1:X:223:VAL:HB	2.21	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	X	444/452 (98%)	338 (76%)	78 (18%)	28 (6%)	2 9

All (28) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	X	6	THR
1	X	15	LYS
1	X	213	PRO
1	X	214	SER
1	X	248	ILE
1	X	398	ILE
1	X	404	VAL
1	X	414	LYS
1	X	431	THR
1	X	434	LYS
1	X	442	GLN
1	X	447	GLN
1	X	85	ASP
1	X	86	SER

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Mol	Chain	Res	Type
1	X	241	ARG
1	X	249	ASP
1	X	365	SER
1	X	369	SER
1	X	392	ASN
1	X	413	SER
1	X	303	ALA
1	X	364	GLY
1	X	384	SER
1	X	168	GLU
1	X	382	ASP
1	X	412	LEU
1	X	251	ASN
1	X	439	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	X	383/387 (99%)	344 (90%)	39 (10%)	9 33

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	X	13	ASP
1	X	21	LEU
1	X	25	LEU
1	X	60	THR
1	X	62	LYS
1	X	67	GLN
1	X	77	HIS
1	X	84	GLU
1	X	85	ASP
1	X	86	SER
1	X	123	ILE
1	X	141	LEU

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Mol	Chain	Res	Type
1	X	145	PRO
1	X	155	LEU
1	X	167	LEU
1	X	171	VAL
1	X	209	SER
1	X	213	PRO
1	X	222	LYS
1	X	240	LYS
1	X	242	SER
1	X	247	PHE
1	X	248	ILE
1	X	287	CYS
1	X	294	GLN
1	X	339	SER
1	X	351	SER
1	X	370	LYS
1	X	393	PHE
1	X	407	LYS
1	X	408	THR
1	X	410	LEU
1	X	424	VAL
1	X	429	ASP
1	X	434	LYS
1	X	435	PHE
1	X	438	SER
1	X	444	ASN
1	X	447	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	X	36	ASN
1	X	65	ASN
1	X	144	ASN
1	X	294	GLN
1	X	315	GLN
1	X	355	ASN
1	X	380	GLN
1	X	406	ASN
1	X	425	GLN

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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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$
2	DDQ	X	501	-	13,13,13	3.32	2 (15%)	14,15,15	3.00	4 (28%)
2	DDQ	X	502	-	13,13,13	3.98	1 (7%)	14,15,15	2.00	1 (7%)

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	DDQ	X	501	-	-	0/11/11/11	0/0/0/0
2	DDQ	X	502	-	-	0/11/11/11	0/0/0/0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	X	502	DDQ	O1-N1	-14.28	1.26	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	X	501	DDQ	O1-N1	-11.33	1.28	1.39
2	X	501	DDQ	CM2-N1	3.24	1.54	1.49

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	X	501	DDQ	O1-N1-C1	-10.21	98.79	110.27
2	X	502	DDQ	O1-N1-C1	-6.89	102.52	110.27
2	X	501	DDQ	CM2-N1-C1	2.09	116.51	109.77
2	X	501	DDQ	CM2-N1-CM1	2.28	111.41	108.83
2	X	501	DDQ	O1-N1-CM1	2.74	112.72	109.05

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
2	X	502	DDQ	1	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	X	448/452 (99%)	0.36	42 (9%) 11 4	12, 47, 100, 100	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	X	438	SER	6.7
1	X	351	SER	6.5
1	X	434	LYS	5.9
1	X	436	CYS	5.6
1	X	350	GLY	5.5
1	X	442	GLN	5.3
1	X	352	GLY	5.3
1	X	399	GLN	5.1
1	X	366	ASN	4.9
1	X	353	LYS	4.9
1	X	412	LEU	4.8
1	X	383	SER	4.2
1	X	416	GLN	4.2
1	X	396	GLY	3.9
1	X	452	CYS	3.8
1	X	417	LEU	3.7
1	X	380	GLN	3.7
1	X	437	GLY	3.7
1	X	354	ALA	3.6
1	X	403	PHE	3.0
1	X	420	SER	2.9
1	X	421	ARG	2.9
1	X	382	ASP	2.8
1	X	381	PRO	2.8
1	X	1	LYS	2.8
1	X	348	LEU	2.7
1	X	424	VAL	2.7

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Mol	Chain	Res	Type	RSRZ
1	X	439	GLY	2.7
1	X	433	TYR	2.6
1	X	413	SER	2.6
1	X	357	TYR	2.5
1	X	130	THR	2.5
1	X	414	LYS	2.5
1	X	409	VAL	2.5
1	X	418	GLY	2.5
1	X	333	GLU	2.4
1	X	213	PRO	2.3
1	X	447	GLN	2.3
1	X	405	TRP	2.2
1	X	385	TYR	2.2
1	X	248	ILE	2.1
1	X	135	GLN	2.1

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
2	DDQ	X	502	14/14	0.93	0.20	0.86	21,26,36,37	0
2	DDQ	X	501	14/14	0.92	0.21	0.56	20,28,35,41	0
3	CA	X	602	1/1	0.97	0.11	-1.63	28,28,28,28	0
3	CA	X	601	1/1	0.95	0.10	-	50,50,50,50	0

6.5 Other polymers ⓘ

There are no such residues in this entry.