



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:19 PM GMT

PDB ID : 4LA0
Title : X-ray study of human serum albumin complexed with bicalutamide
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Deposited on : 2013-06-18
Resolution : 2.40 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

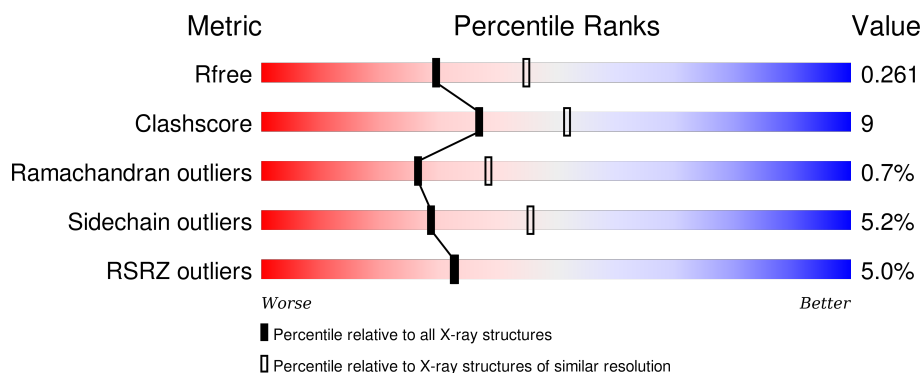
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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	585	<div> <div>3%</div> <div>75%</div> <div>22%</div> <div>..</div> </div>
1	B	585	<div> <div>7%</div> <div>78%</div> <div>19%</div> <div>..</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	198	A	601	-	-	-	X

2 Entry composition [i](#)

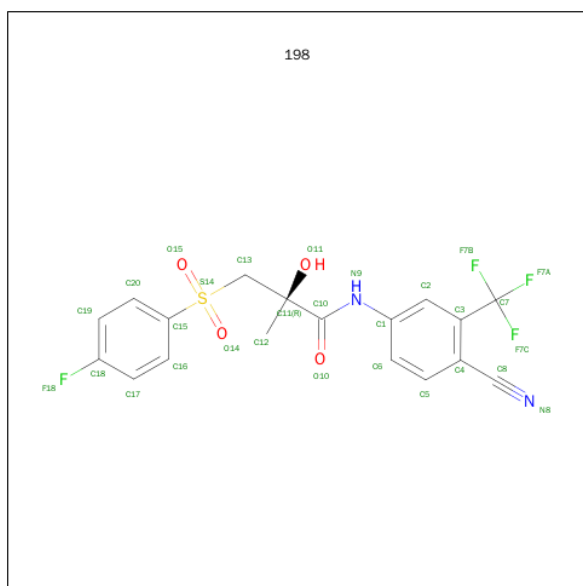
There are 3 unique types of molecules in this entry. The entry contains 9482 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 SERUM ALBUMIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	581	Total	C	N	O	S	0	0	0
			4623	2918	782	882	41			
1	B	580	Total	C	N	O	S	0	0	0
			4618	2915	781	881	41			

- Molecule 2 is R-BICALUTAMIDE (three-letter code: 198) (formula: C₁₈H₁₄F₄N₂O₄S).

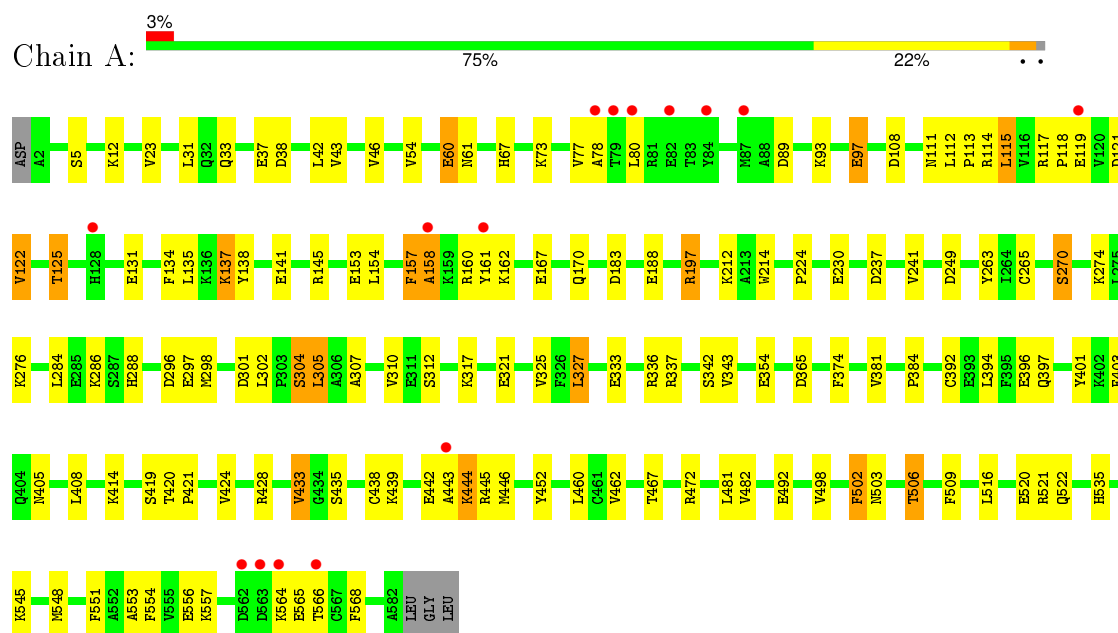


Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	123	Total 123	O 123	0	0
3	B	60	Total 60	O 60	0	0

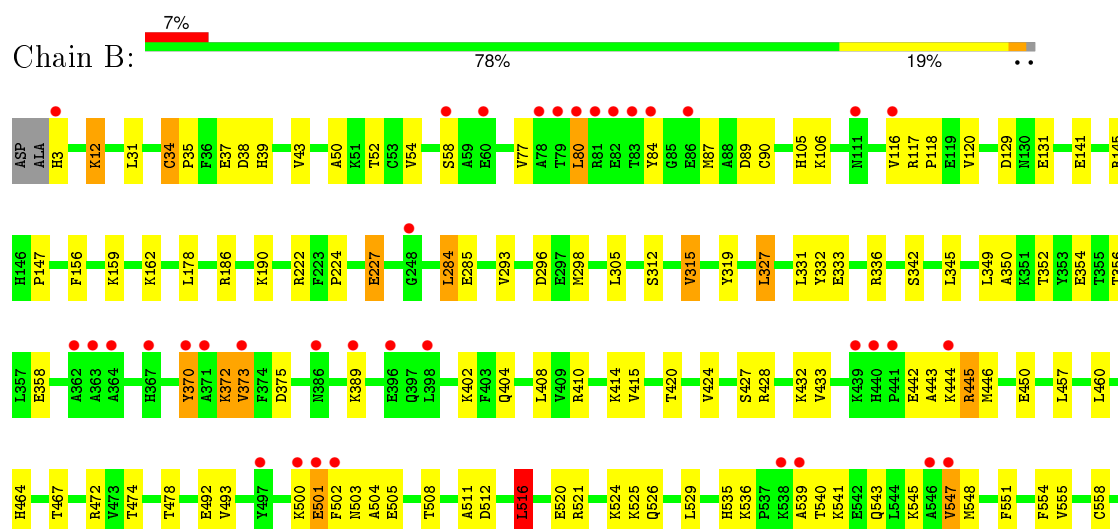
3 Residue-property plots

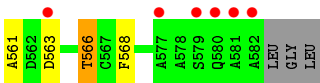
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: SERUM ALBUMIN



• Molecule 1: SERUM ALBUMIN





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	56.44Å 59.57Å 95.96Å 73.22° 83.93° 74.17°	Depositor
Resolution (Å)	29.86 – 2.40 29.86 – 2.40	Depositor EDS
% Data completeness (in resolution range)	93.3 (29.86-2.40) 80.0 (29.86-2.40)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.74 (at 2.39Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.199 , 0.254 0.214 , 0.261	Depositor DCC
R_{free} test set	1877 reflections (5.01%)	DCC
Wilson B-factor (Å ²)	35.8	Xtriage
Anisotropy	0.294	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 47.3	EDS
Estimated twinning fraction	0.004 for -h,-k,-l+1	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 41969 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9482	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.37% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section:
198

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	0.52	0/4713	0.70	4/6357 (0.1%)
1	B	0.46	0/4708	0.63	1/6350 (0.0%)
All	All	0.49	0/9421	0.67	5/12707 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	158	ALA	N-CA-CB	10.47	124.76	110.10
1	A	157	PHE	CA-C-O	-5.89	107.73	120.10
1	B	516	LEU	CA-CB-CG	5.64	128.28	115.30
1	A	305	LEU	CA-CB-CG	5.37	127.65	115.30
1	A	160	ARG	N-CA-C	5.04	124.62	111.00

There are no chirality outliers.

There are no planarity outliers.

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	A	4623	0	4543	89	0
1	B	4618	0	4538	82	1
2	A	29	0	14	4	0
2	B	29	0	14	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	123	0	0	10	0
3	B	60	0	0	6	0
All	All	9482	0	9109	173	1

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 9.

All (173) 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:197:ARG:NH2	3:A:725:HOH:O	1.82	1.08
1:A:135:LEU:HG	1:A:161:TYR:CE1	2.05	0.90
1:B:89:ASP:OD2	3:B:754:HOH:O	1.88	0.90
1:A:333:GLU:OE1	1:A:337:ARG:NH2	2.08	0.85
1:B:501:GLU:O	1:B:535:HIS:NE2	2.10	0.85
1:A:158:ALA:HA	1:A:161:TYR:HE2	1.40	0.84
2:B:601:198:F7A	3:B:745:HOH:O	1.86	0.82
1:B:34:CYS:HG	1:B:84:TYR:HH	0.84	0.82
1:B:80:LEU:HD11	1:B:84:TYR:CE2	2.19	0.76
1:B:414:LYS:O	1:B:472:ARG:NH1	2.16	0.76
1:B:500:LYS:O	1:B:502:PHE:N	2.18	0.76
1:A:31:LEU:HD11	1:A:78:ALA:HB2	1.67	0.75
1:A:158:ALA:HA	1:A:161:TYR:CE2	2.25	0.71
1:A:405:ASN:HA	1:A:408:LEU:HD12	1.72	0.71
1:A:522:GLN:OE1	3:A:742:HOH:O	2.09	0.69
1:A:135:LEU:HG	1:A:161:TYR:HE1	1.57	0.68
1:A:428:ARG:NH2	3:A:742:HOH:O	2.27	0.66
1:B:80:LEU:HD11	1:B:84:TYR:CD2	2.31	0.66
1:A:270:SER:O	3:A:733:HOH:O	2.14	0.66
1:A:119:GLU:HB2	1:A:122:VAL:HG22	1.76	0.66
1:B:159:LYS:NZ	1:B:285:GLU:OE1	2.28	0.66
1:A:230:GLU:OE1	3:A:766:HOH:O	2.15	0.65
1:B:186:ARG:NH2	3:B:729:HOH:O	1.88	0.65
1:B:87:MET:HG2	1:B:105:HIS:CD2	2.31	0.65
1:A:435:SER:O	1:A:439:LYS:HE3	1.96	0.65
1:A:157:PHE:HZ	1:A:188:GLU:HG2	1.62	0.63
1:B:12:LYS:HE2	1:B:54:VAL:HG13	1.81	0.63
1:B:116:VAL:N	3:B:720:HOH:O	2.30	0.63
1:A:12:LYS:HD3	1:A:54:VAL:HG13	1.81	0.62
1:B:370:TYR:O	1:B:373:VAL:HG22	2.00	0.61
1:A:298:MET:HE3	1:A:302:LEU:HD12	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:114:ARG:NH2	1:A:520:GLU:OE2	2.28	0.60
1:A:214:TRP:CD1	1:A:343:VAL:HG11	2.36	0.60
1:A:135:LEU:HD21	1:A:162:LYS:HB2	1.84	0.60
1:A:419:SER:OG	1:A:421:PRO:HD2	2.02	0.59
1:A:424:VAL:O	1:A:428:ARG:HG3	2.02	0.59
1:A:167:GLU:O	1:A:170:GLN:HG2	2.04	0.58
1:A:502:PHE:HB2	1:A:535:HIS:CE1	2.39	0.58
1:A:46:VAL:HG23	1:A:73:LYS:HG2	1.85	0.58
1:A:154:LEU:O	1:A:158:ALA:N	2.36	0.57
1:B:558:CYS:HA	1:B:561:ALA:HB2	1.87	0.57
1:B:467:THR:OG1	1:B:467:THR:O	2.19	0.57
1:B:80:LEU:O	1:B:80:LEU:HD12	2.05	0.56
1:B:327:LEU:HD21	1:B:354:GLU:HB2	1.87	0.56
1:A:153:GLU:OE2	1:A:288:HIS:ND1	2.35	0.56
1:B:332:TYR:O	1:B:336:ARG:HG3	2.06	0.56
1:B:525:LYS:HG2	1:B:551:PHE:CE1	2.41	0.56
1:B:402:LYS:HG2	1:B:545:LYS:HZ1	1.71	0.55
1:B:224:PRO:HD2	1:B:296:ASP:HB3	1.88	0.55
1:A:516:LEU:O	1:A:521:ARG:NH2	2.39	0.55
1:B:516:LEU:HB3	1:B:520:GLU:HB2	1.89	0.55
1:B:543:GLN:O	1:B:547:VAL:HG12	2.07	0.54
1:A:138:TYR:CD2	1:A:161:TYR:CE2	2.96	0.54
1:A:414:LYS:O	1:A:472:ARG:NH2	2.41	0.54
1:B:408:LEU:HD13	1:B:427:SER:HB2	1.91	0.54
1:A:114:ARG:HH22	1:A:520:GLU:CD	2.08	0.53
1:B:141:GLU:OE1	1:B:145:ARG:NH1	2.37	0.53
1:A:442:GLU:O	1:A:444:LYS:N	2.41	0.53
1:A:302:LEU:HB3	1:A:337:ARG:NH1	2.23	0.53
1:A:224:PRO:HD2	1:A:296:ASP:HB3	1.90	0.53
1:B:561:ALA:O	1:B:563:ASP:N	2.35	0.52
1:A:108:ASP:OD1	1:A:197:ARG:HD2	2.08	0.52
1:B:227:GLU:HB2	3:B:724:HOH:O	2.10	0.52
1:B:516:LEU:O	1:B:521:ARG:NH2	2.42	0.52
1:A:230:GLU:OE2	1:A:263:TYR:OH	2.18	0.51
1:B:50:ALA:O	1:B:54:VAL:HG23	2.10	0.51
1:A:77:VAL:O	1:A:80:LEU:HD12	2.09	0.51
1:B:39:HIS:O	1:B:43:VAL:HG23	2.10	0.51
1:A:118:PRO:HG2	2:A:601:198:H121	1.92	0.51
1:A:274:LYS:NZ	1:A:297:GLU:OE2	2.28	0.51
1:B:428:ARG:NH1	1:B:526:GLN:OE1	2.42	0.51
1:B:410:ARG:HG2	1:B:414:LYS:HE2	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:135:LEU:HG	1:A:161:TYR:CZ	2.46	0.50
1:A:212:LYS:NZ	3:A:789:HOH:O	2.39	0.50
1:A:237:ASP:O	1:A:241:VAL:HG23	2.11	0.50
1:A:553:ALA:O	1:A:556:GLU:HB3	2.11	0.50
1:B:331:LEU:HD13	1:B:350:ALA:HB2	1.93	0.50
1:A:381:VAL:O	1:A:384:PRO:HD2	2.12	0.50
1:A:317:LYS:HE2	1:A:321:GLU:OE2	2.12	0.49
1:A:60:GLU:HG2	1:A:61:ASN:ND2	2.26	0.49
1:A:503:ASN:HB3	1:A:506:THR:HG23	1.93	0.49
1:B:89:ASP:OD1	1:B:90:CYS:N	2.45	0.49
1:B:3:HIS:O	1:B:3:HIS:CD2	2.66	0.49
1:B:37:GLU:CD	1:B:37:GLU:H	2.15	0.48
1:B:501:GLU:O	1:B:535:HIS:CE1	2.66	0.48
1:A:117:ARG:NH2	1:A:183:ASP:OD1	2.46	0.48
1:A:438:CYS:O	1:A:445:ARG:NH2	2.46	0.48
1:B:305:LEU:HD21	1:B:333:GLU:HB3	1.96	0.47
1:A:433:VAL:HG23	1:A:452:TYR:CD2	2.49	0.47
1:B:536:LYS:O	1:B:540:THR:OG1	2.30	0.47
1:A:89:ASP:OD2	3:A:778:HOH:O	2.20	0.47
1:A:115:LEU:HD21	1:A:141:GLU:HB3	1.95	0.47
1:B:319:TYR:OH	1:B:358:GLU:OE2	2.26	0.47
1:A:119:GLU:OE2	3:A:757:HOH:O	2.20	0.47
1:A:310:VAL:HG13	1:A:374:PHE:HE2	1.79	0.47
1:B:349:LEU:HA	1:B:352:THR:HG22	1.97	0.47
1:A:420:THR:O	1:A:424:VAL:HG23	2.15	0.47
1:A:134:PHE:CE1	2:A:601:198:H131	2.50	0.47
1:B:420:THR:O	1:B:424:VAL:HG23	2.14	0.47
1:A:342:SER:HB3	1:A:446:MET:HG2	1.96	0.47
1:A:60:GLU:OE1	1:A:61:ASN:N	2.44	0.47
1:B:561:ALA:HB1	1:B:563:ASP:OD1	2.16	0.46
1:B:156:PHE:CD1	1:B:284:LEU:HB3	2.51	0.46
1:A:516:LEU:HB3	1:A:520:GLU:HB2	1.96	0.46
1:A:33:GLN:NE2	3:A:822:HOH:O	1.98	0.46
1:B:511:ALA:HA	1:B:568:PHE:CE2	2.51	0.46
1:B:186:ARG:O	1:B:190:LYS:HG2	2.16	0.46
1:B:80:LEU:CD1	1:B:84:TYR:CD2	2.97	0.46
1:A:327:LEU:HD21	1:A:354:GLU:HG3	1.98	0.45
1:A:154:LEU:HA	1:A:157:PHE:HB2	1.98	0.45
1:B:131:GLU:OE2	1:B:162:LYS:HE3	2.17	0.45
1:B:356:THR:HG21	1:B:373:VAL:HA	1.99	0.45
1:A:564:LYS:C	1:A:566:THR:H	2.20	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:464:HIS:HE1	1:B:474:THR:OG1	2.00	0.45
1:B:31:LEU:HB2	1:B:39:HIS:HE1	1.82	0.45
1:B:372:LYS:HB2	1:B:375:ASP:HB2	1.99	0.45
1:A:115:LEU:HD22	1:A:145:ARG:CZ	2.47	0.45
1:B:521:ARG:HD3	1:B:555:VAL:HG11	1.98	0.44
1:A:554:PHE:HA	1:A:557:LYS:HB3	1.98	0.44
1:B:408:LEU:HD21	1:B:526:GLN:HB3	1.98	0.44
1:B:389:LYS:HA	1:B:445:ARG:HH12	1.83	0.44
1:A:298:MET:CE	1:A:302:LEU:HD12	2.46	0.44
1:A:121:ASP:O	1:A:125:THR:OG1	2.35	0.44
1:A:23:VAL:HG12	1:A:43:VAL:HG22	2.00	0.44
1:B:120:VAL:HG13	1:B:178:LEU:HD23	1.99	0.44
2:A:601:198:H2	2:A:601:198:O10	2.18	0.44
1:B:80:LEU:HD11	1:B:84:TYR:HE2	1.74	0.44
1:B:35:PRO:HG2	1:B:38:ASP:OD2	2.18	0.44
1:B:404:GLN:O	1:B:408:LEU:HB2	2.18	0.43
1:A:93:LYS:HD2	1:A:97:GLU:HB2	1.99	0.43
1:B:117:ARG:HA	1:B:118:PRO:HD3	1.85	0.43
1:A:131:GLU:OE2	1:A:162:LYS:HE3	2.18	0.43
1:A:38:ASP:O	1:A:42:LEU:HG	2.18	0.43
1:A:433:VAL:HG23	1:A:452:TYR:HD2	1.83	0.43
1:B:504:ALA:O	1:B:508:THR:HG23	2.19	0.43
1:A:111:ASN:OD1	3:A:808:HOH:O	2.21	0.43
1:A:401:TYR:CE1	1:A:522:GLN:HG2	2.54	0.43
1:B:408:LEU:HD23	1:B:529:LEU:HD23	1.99	0.43
1:A:117:ARG:HA	1:A:118:PRO:HD3	1.75	0.43
1:A:67:HIS:NE2	1:A:249:ASP:OD1	2.52	0.43
1:B:415:VAL:HA	1:B:472:ARG:HH22	1.83	0.43
1:B:331:LEU:HD12	1:B:331:LEU:HA	1.83	0.43
1:A:37:GLU:CD	1:A:37:GLU:H	2.22	0.43
1:B:512:ASP:O	1:B:516:LEU:HD13	2.19	0.43
1:A:60:GLU:N	1:A:60:GLU:OE1	2.52	0.42
1:B:457:LEU:HA	1:B:457:LEU:HD23	1.86	0.42
1:B:284:LEU:HA	1:B:284:LEU:HD12	1.90	0.42
1:B:222:ARG:HD2	1:B:293:VAL:CG1	2.50	0.42
1:A:265:CYS:SG	1:A:286:LYS:HD2	2.58	0.42
1:B:222:ARG:HD2	1:B:293:VAL:HG13	2.01	0.42
1:A:137:LYS:HD3	2:A:601:198:O15	2.20	0.42
1:A:197:ARG:HD3	1:A:462:VAL:HG21	2.02	0.42
1:A:394:LEU:HD23	1:A:403:PHE:HE1	1.83	0.42
1:B:545:LYS:HD2	1:B:545:LYS:HA	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:392:CYS:O	1:A:396:GLU:HG2	2.19	0.41
1:A:112:LEU:HA	1:A:113:PRO:HD2	1.84	0.41
1:B:345:LEU:O	1:B:349:LEU:HG	2.20	0.41
1:B:442:GLU:O	1:B:444:LYS:N	2.53	0.41
1:B:410:ARG:O	1:B:414:LYS:HG3	2.21	0.41
1:A:307:ALA:O	1:A:312:SER:HB2	2.20	0.41
1:B:561:ALA:C	1:B:563:ASP:H	2.22	0.41
1:B:342:SER:HB3	1:B:446:MET:HG2	2.03	0.41
1:A:304:SER:O	1:A:305:LEU:CB	2.69	0.41
1:B:503:ASN:OD1	1:B:505:GLU:N	2.53	0.41
1:B:106:LYS:HD3	1:B:147:PRO:HB2	2.03	0.41
1:A:394:LEU:HD12	1:A:397:GLN:HE21	1.86	0.40
1:A:509:PHE:O	1:A:568:PHE:HB3	2.21	0.40
1:B:116:VAL:HG12	3:B:720:HOH:O	2.20	0.40
1:B:428:ARG:O	1:B:432:LYS:HG3	2.22	0.40
1:A:481:LEU:HA	1:A:481:LEU:HD12	1.74	0.40
1:B:563:ASP:O	1:B:566:THR:HG23	2.21	0.40
1:B:31:LEU:HB2	1:B:39:HIS:CE1	2.56	0.40
1:B:312:SER:O	1:B:315:VAL:HG12	2.22	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:129:ASP:OD1	1:B:541:LYS:NZ[1_455]	2.08	0.12

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	579/585 (99%)	553 (96%)	21 (4%)	5 (1%)	21 30

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	578/585 (99%)	549 (95%)	26 (4%)	3 (0%)	34	48
All	All	1157/1170 (99%)	1102 (95%)	47 (4%)	8 (1%)	26	38

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	444	LYS
1	B	501	GLU
1	A	502	PHE
1	B	443	ALA
1	B	539	ALA
1	A	443	ALA
1	A	5	SER
1	A	565	GLU

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	508/511 (99%)	482 (95%)	26 (5%)	29	46
1	B	508/511 (99%)	481 (95%)	27 (5%)	28	44
All	All	1016/1022 (99%)	963 (95%)	53 (5%)	29	45

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	60	GLU
1	A	97	GLU
1	A	115	LEU
1	A	122	VAL
1	A	125	THR
1	A	137	LYS
1	A	197	ARG
1	A	270	SER

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Mol	Chain	Res	Type
1	A	276	LYS
1	A	284	LEU
1	A	301	ASP
1	A	304	SER
1	A	325	VAL
1	A	327	LEU
1	A	336	ARG
1	A	365	ASP
1	A	433	VAL
1	A	460	LEU
1	A	467	THR
1	A	482	VAL
1	A	492	GLU
1	A	498	VAL
1	A	506	THR
1	A	545	LYS
1	A	548	MET
1	A	551	PHE
1	B	12	LYS
1	B	34	CYS
1	B	52	THR
1	B	58	SER
1	B	77	VAL
1	B	80	LEU
1	B	227	GLU
1	B	284	LEU
1	B	298	MET
1	B	315	VAL
1	B	327	LEU
1	B	370	TYR
1	B	372	LYS
1	B	373	VAL
1	B	433	VAL
1	B	445	ARG
1	B	450	GLU
1	B	460	LEU
1	B	478	THR
1	B	492	GLU
1	B	493	VAL
1	B	516	LEU
1	B	524	LYS
1	B	547	VAL

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Mol	Chain	Res	Type
1	B	548	MET
1	B	554	PHE
1	B	566	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	61	ASN
1	B	3	HIS
1	B	105	HIS

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	198	A	601	-	28,30,30	1.36	2 (7%)	38,46,46	2.55	6 (15%)
2	198	B	601	-	28,30,30	1.40	3 (10%)	38,46,46	2.27	4 (10%)

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	198	A	601	-	-	0/30/30/30	0/2/2/2
2	198	B	601	-	-	0/30/30/30	0/2/2/2

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	198	C7-C3	2.06	1.55	1.50
2	A	601	198	C10-N9	2.72	1.42	1.35
2	B	601	198	C10-N9	3.14	1.43	1.35
2	B	601	198	C4-C8	5.05	1.52	1.44
2	A	601	198	C4-C8	5.23	1.52	1.44

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	198	O15-S14-O14	-13.00	101.22	118.40
2	B	601	198	O15-S14-O14	-12.14	102.36	118.40
2	A	601	198	F7C-C7-C3	-3.03	107.36	112.68
2	A	601	198	F7A-C7-C3	-2.56	108.18	112.68
2	A	601	198	F7A-C7-F7B	2.04	113.06	105.71
2	B	601	198	C2-C3-C7	2.17	121.34	116.51
2	B	601	198	O14-S14-C15	2.20	110.79	108.38
2	A	601	198	O15-S14-C15	2.32	110.92	108.38
2	B	601	198	O15-S14-C15	3.12	111.80	108.38
2	A	601	198	O14-S14-C15	4.98	113.84	108.38

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	198	4	0
2	B	601	198	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	A	581/585 (99%)	-0.10	15 (2%) 59 58	20, 38, 67, 94	0
1	B	580/585 (99%)	0.23	43 (7%) 17 17	30, 53, 87, 101	0
All	All	1161/1170 (99%)	0.07	58 (4%) 32 33	20, 45, 82, 101	0

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	582	ALA	10.7
1	A	161	TYR	6.6
1	B	79	THR	6.3
1	B	579	SER	6.1
1	A	79	THR	5.8
1	B	581	ALA	5.2
1	B	84	TYR	5.1
1	A	158	ALA	4.6
1	B	370	TYR	4.3
1	B	371	ALA	4.2
1	B	539	ALA	4.2
1	B	83	THR	4.0
1	B	367	HIS	3.9
1	B	364	ALA	3.9
1	B	81	ARG	3.9
1	B	563	ASP	3.7
1	B	502	PHE	3.6
1	B	577	ALA	3.5
1	A	564	LYS	3.4
1	B	580	GLN	3.4
1	B	538	LYS	3.3
1	B	80	LEU	3.3
1	B	3	HIS	3.2
1	A	562	ASP	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	373	VAL	3.2
1	B	78	ALA	3.1
1	B	82	GLU	3.1
1	A	443	ALA	3.1
1	B	501	GLU	2.9
1	B	439	LYS	2.8
1	A	87	MET	2.7
1	B	86	GLU	2.7
1	B	396	GLU	2.6
1	B	546	ALA	2.6
1	B	116	VAL	2.5
1	B	440	HIS	2.5
1	B	363	ALA	2.5
1	B	547	VAL	2.5
1	B	60	GLU	2.4
1	A	84	TYR	2.4
1	A	128	HIS	2.4
1	A	82	GLU	2.4
1	A	80	LEU	2.3
1	B	58	SER	2.3
1	B	497	TYR	2.3
1	B	389	LYS	2.3
1	B	441	PRO	2.3
1	B	248	GLY	2.3
1	B	500	LYS	2.2
1	A	563	ASP	2.2
1	A	119	GLU	2.2
1	B	362	ALA	2.1
1	B	444	LYS	2.1
1	A	78	ALA	2.1
1	B	398	LEU	2.0
1	A	566	THR	2.0
1	B	111	ASN	2.0
1	B	386	ASN	2.0

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

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	198	A	601	29/29	0.89	0.27	3.22	42,60,72,77	0
2	198	B	601	29/29	0.90	0.21	0.93	49,58,68,75	0

6.5 Other polymers [i](#)

There are no such residues in this entry.