



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 11:59 AM GMT

PDB ID : 4JK4
Title : Crystal Structure of Bovine Serum Albumin in complex with 3,5-diiodosalicylic acid
Authors : Zielinski, K.; Bujacz, A.; Sekula, B.; Bujacz, G.
Deposited on : 2013-03-09
Resolution : 2.65 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

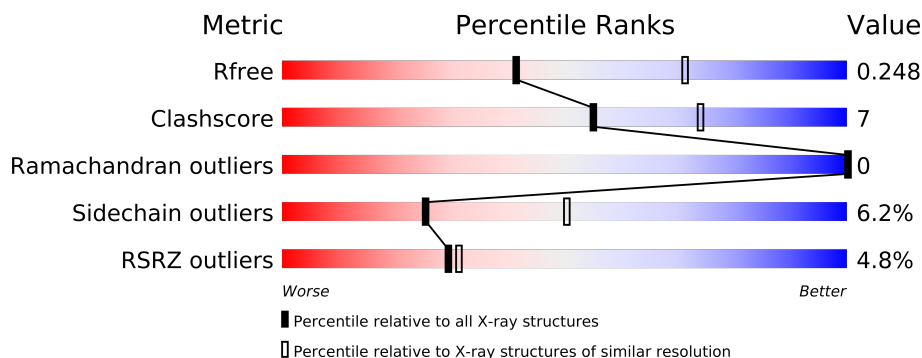
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.65 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}	66092	2232 (2.70-2.62)
Clashscore	79885	2700 (2.70-2.62)
Ramachandran outliers	78287	2657 (2.70-2.62)
Sidechain outliers	78261	2657 (2.70-2.62)
RSRZ outliers	66119	2234 (2.70-2.62)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	583	
1	B	583	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	DIU	B	602	-	X
4	1PE	B	606	-	X
5	PGE	B	607	-	X

2 Entry composition i

There are 7 unique types of molecules in this entry. The entry contains 9643 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

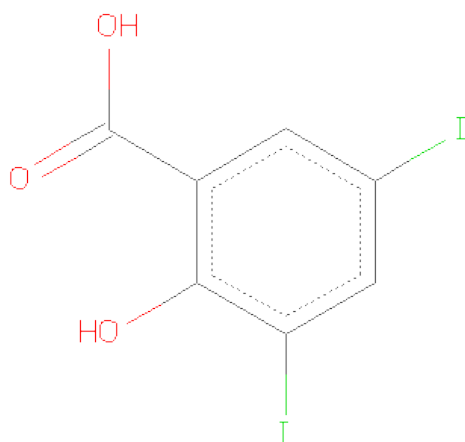
- Molecule 1 is a protein called Serum albumin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	582	Total	C	N	O	S	0	3	0
			4668	2944	782	903	39			
1	B	582	Total	C	N	O	S	0	0	0
			4645	2931	780	895	39			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	190	THR	ALA	VARAINT	UNP P02769
B	190	THR	ALA	VARAINT	UNP P02769

- Molecule 2 is 2-HYDROXY-3,5-DIIDO-BENZOICACID (three-letter code: DIU) (formula: $C_7H_4I_2O_3$).



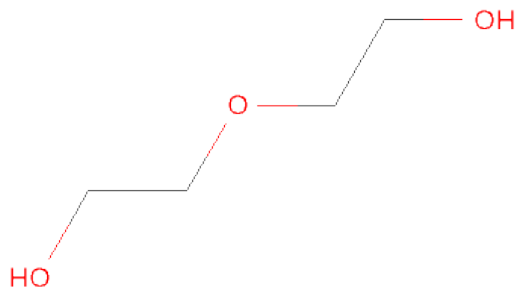
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	I	O	0	0
			12	7	2	3		

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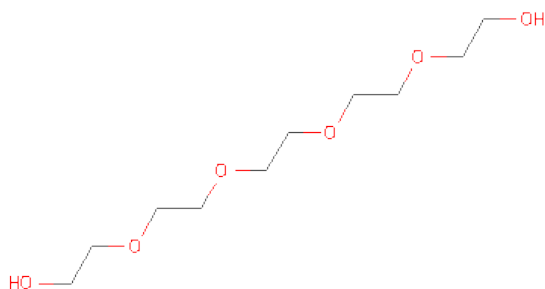
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	I	O	0	0
			12	7	2	3		
2	A	1	Total	C	I	O	0	0
			12	7	2	3		
2	A	1	Total	C	I	O	0	0
			12	7	2	3		
2	B	1	Total	C	I	O	0	0
			12	7	2	3		
2	B	1	Total	C	I	O	0	0
			12	7	2	3		
2	B	1	Total	C	I	O	0	0
			12	7	2	3		

- Molecule 3 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



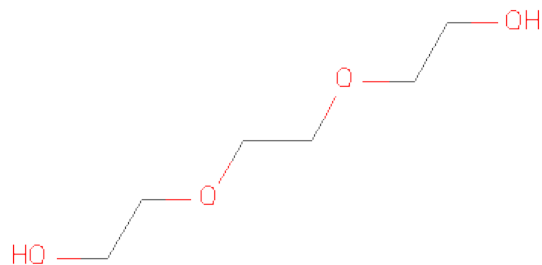
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			7	4	3		
3	B	1	Total	C	O	0	0
			7	4	3		

- Molecule 4 is PENTAETHYLENE GLYCOL (three-letter code: 1PE) (formula: C₁₀H₂₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			16	10	6		
4	A	1	Total	C	O	0	0
			16	10	6		
4	B	1	Total	C	O	0	0
			16	10	6		

- Molecule 5 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			10	6	4		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			10	6	4		

- Molecule 6 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	3	Total	Ca	0	0
			3	3		
6	A	4	Total	Ca	0	0
			4	4		

- Molecule 7 is water.

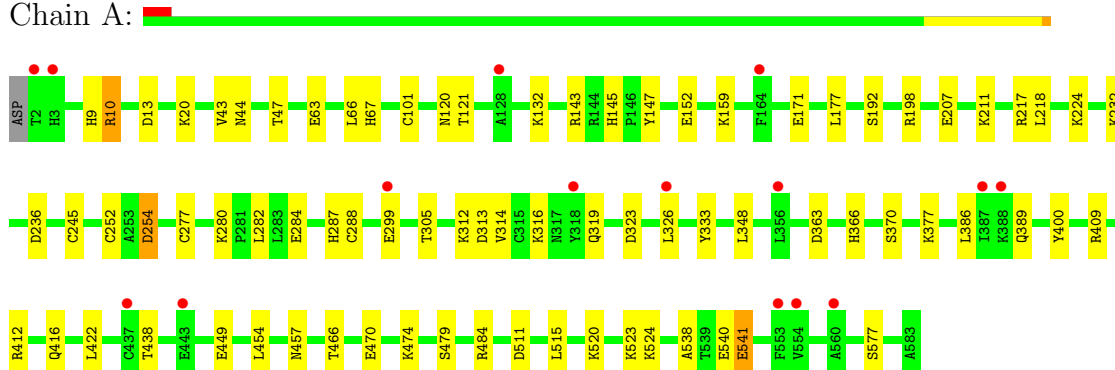
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	82	Total	O	0	0
			82	82		
7	B	63	Total	O	0	0
			63	63		

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.

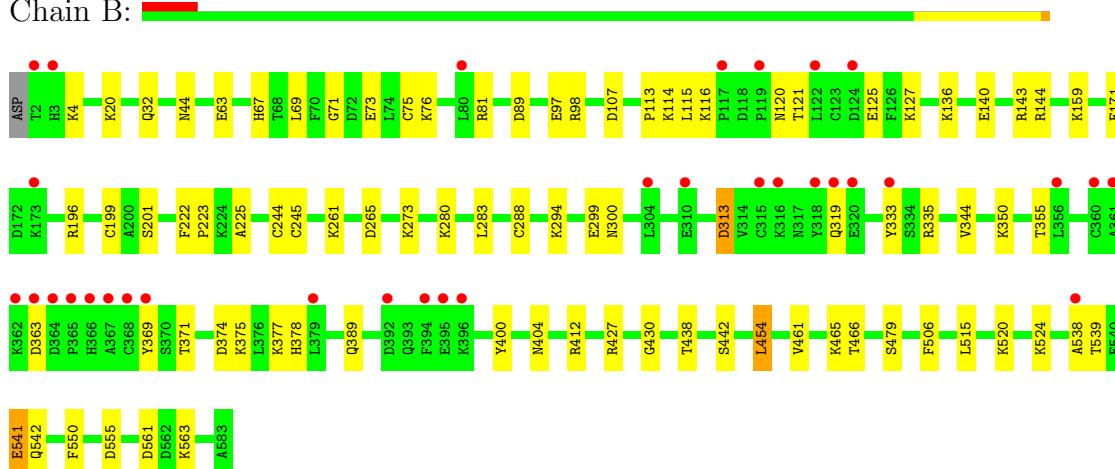
• Molecule 1: Serum albumin

Chain A:



• Molecule 1: Serum albumin

Chain B:



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	215.41 Å 44.80 Å 146.90 Å 90.00° 115.77° 90.00°	Depositor
Resolution (Å)	36.53 – 2.65 44.10 – 2.65	Depositor EDS
% Data completeness (in resolution range)	98.6 (36.53-2.65) 98.7 (44.10-2.65)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.41 (at 2.65 Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.188 , 0.246 0.191 , 0.248	Depositor DCC
R_{free} test set	1840 reflections (5.00%)	DCC
Wilson B-factor (Å ²)	60.8	Xtriage
Anisotropy	0.433	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 30.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 36861 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9643	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.45% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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: PEG, DIU, CA, PGE, 1PE

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	0.99	3/4766 (0.1%)	0.97	4/6434 (0.1%)
1	B	0.91	3/4740 (0.1%)	0.91	2/6399 (0.0%)
All	All	0.95	6/9506 (0.1%)	0.94	6/12833 (0.0%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	245	CYS	CB-SG	5.83	1.92	1.82
1	A	277	CYS	CB-SG	-5.66	1.72	1.81
1	B	199	CYS	CB-SG	5.40	1.91	1.82
1	A	245	CYS	CB-SG	5.29	1.91	1.82
1	A	147	TYR	CB-CG	-5.25	1.43	1.51
1	B	75	CYS	CB-SG	5.05	1.90	1.82

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	10	ARG	NE-CZ-NH2	-5.50	117.55	120.30
1	B	69	LEU	CA-CB-CG	5.37	127.65	115.30
1	A	348	LEU	CA-CB-CG	-5.28	103.15	115.30
1	A	254	ASP	CB-CG-OD1	5.24	123.02	118.30
1	A	454	LEU	CB-CG-CD2	-5.23	102.12	111.00
1	B	196	ARG	NE-CZ-NH2	-5.15	117.73	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4668	0	0	29	0
1	B	4645	0	0	34	0
2	A	48	0	8	2	0
2	B	48	0	9	3	0
3	A	7	0	10	0	0
3	B	7	0	10	1	0
4	A	32	0	44	0	0
4	B	16	0	22	0	0
5	A	10	0	14	0	0
5	B	10	0	14	2	0
6	A	4	0	0	0	0
6	B	3	0	0	0	0
7	A	82	0	0	3	0
7	B	63	0	0	7	0
All	All	9643	0	131	64	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 7.

All (64) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:223:PRO:O	1:B:335:ARG:NH1	2.23	0.72
1:B:273:LYS:NZ	7:B:757:HOH:O	2.25	0.69
1:B:319:GLN:NE2	7:B:721:HOH:O	2.26	0.69
1:B:541:GLU:OE1	1:B:542:GLN:N	2.26	0.69
1:A:63:GLU:OE1	1:A:63:GLU:N	2.27	0.67
1:A:412:ARG:NH2	1:A:538:ALA:O	2.28	0.66
1:A:466:THR:OG1	1:A:466:THR:O	2.13	0.65
1:A:470:GLU:N	1:A:470:GLU:OE1	2.31	0.64
1:A:515:LEU:O	1:A:520:LYS:NZ	2.31	0.63
1:B:125:GLU:OE2	1:B:136:LYS:NZ	2.31	0.63
1:A:20:LYS:NZ	1:A:44:ASN:OD1	2.32	0.63
1:A:280:LYS:NZ	1:A:288:CYS:SG	2.74	0.60
1:B:524:LYS:NZ	7:B:758:HOH:O	2.34	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:43:VAL:O	1:A:47:THR:OG1	2.20	0.59
1:A:400:TYR:OH	1:A:524:LYS:NZ	2.36	0.58
1:B:283:LEU:N	7:B:727:HOH:O	2.37	0.57
1:A:449:GLU:OE2	1:A:484:ARG:NH2	2.37	0.57
1:A:217:ARG:NH1	7:A:703:HOH:O	2.36	0.57
1:B:313:ASP:N	1:B:313:ASP:OD1	2.38	0.57
1:B:63:GLU:OE1	1:B:63:GLU:N	2.37	0.56
1:A:323:ASP:OD1	1:A:323:ASP:N	2.40	0.54
1:A:120:ASN:OD1	1:A:121:THR:N	2.40	0.54
1:B:115:LEU:CB	2:B:603:DIU:H6	2.38	0.53
1:A:9:HIS:NE2	1:A:13:ASP:OD2	2.41	0.53
1:B:171:GLU:OE1	1:B:171:GLU:N	2.42	0.53
1:B:120:ASN:OD1	1:B:121:THR:N	2.42	0.53
1:A:198:ARG:NH2	2:A:604:DIU:O2	2.43	0.52
1:A:152:GLU:OE2	1:A:287:HIS:ND1	2.43	0.51
1:B:261:LYS:NZ	1:B:265:ASP:OD1	2.43	0.51
1:A:207:GLU:OE2	1:A:211:LYS:NZ	2.45	0.50
1:B:140:GLU:OE1	1:B:144:ARG:NH1	2.45	0.50
1:B:466:THR:O	1:B:466:THR:OG1	2.29	0.49
1:B:412:ARG:NH2	1:B:538:ALA:O	2.46	0.49
1:B:97:GLU:OE1	1:B:97:GLU:N	2.46	0.49
1:B:32:GLN:OE1	1:B:107:ASP:N	2.45	0.49
1:B:73:GLU:OE2	1:B:76:LYS:NZ	2.45	0.49
1:B:20:LYS:NZ	1:B:44:ASN:OD1	2.46	0.48
1:B:454:LEU:CD1	2:B:604:DIU:I1	3.31	0.48
1:B:400:TYR:CE2	1:B:404:ASN:ND2	2.82	0.48
1:A:232:LYS:NZ	1:A:236:ASP:OD2	2.47	0.48
1:A:145:HIS:ND1	1:A:192:SER:OG	2.47	0.47
1:B:427:ARG:O	1:B:430:GLY:N	2.47	0.47
1:B:222:PHE:O	1:B:225:ALA:N	2.47	0.47
1:B:67:HIS:CD2	7:B:759:HOH:O	2.66	0.47
1:B:71:GLY:CA	7:B:704:HOH:O	2.62	0.47
1:B:115:LEU:O	1:B:116:LYS:NZ	2.48	0.46
1:B:506:PHE:CD1	5:B:607:PGE:H12	2.50	0.46
1:B:375:LYS:O	1:B:378:HIS:ND1	2.48	0.46
1:B:98:ARG:NH1	7:B:705:HOH:O	2.49	0.46
1:A:479:SER:O	1:A:479:SER:OG	2.32	0.45
1:A:313:ASP:O	1:A:314:VAL:C	2.55	0.45
1:B:81:ARG:NH2	1:B:89:ASP:OD1	2.50	0.44
1:A:67:HIS:CE1	7:A:763:HOH:O	2.71	0.44
1:B:550:PHE:CD1	5:B:607:PGE:H42	2.53	0.44
1:A:541:GLU:CD	1:A:541:GLU:N	2.71	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:9:HIS:CD2	1:A:13:ASP:OD2	2.70	0.43
1:A:10:ARG:NH1	1:A:254:ASP:OD2	2.51	0.43
2:B:603:DIU:I2	3:B:605:PEG:H32	2.89	0.43
1:A:366:HIS:O	1:A:370:SER:OG	2.37	0.43
1:B:515:LEU:O	1:B:520:LYS:NZ	2.52	0.42
1:A:313:ASP:O	1:A:316:LYS:N	2.53	0.42
1:B:113:PRO:O	1:B:144:ARG:NH2	2.53	0.41
1:A:409:ARG:NH2	2:A:602:DIU:O1	2.53	0.41
1:A:457:ASN:ND2	7:A:750:HOH:O	2.53	0.41

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

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	583/583 (100%)	570 (98%)	13 (2%)	0	100	100
1	B	580/583 (100%)	563 (97%)	17 (3%)	0	100	100
All	All	1163/1166 (100%)	1133 (97%)	30 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	523/521 (100%)	491 (94%)	32 (6%)	26	52
1	B	520/521 (100%)	486 (94%)	34 (6%)	24	48
All	All	1043/1042 (100%)	977 (94%)	66 (6%)	25	50

All (66) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	66	LEU
1	A	101	CYS
1	A	132	LYS
1	A	143	ARG
1	A	159	LYS
1	A	171	GLU
1	A	177	LEU
1	A	218	LEU
1	A	224	LYS
1	A	252	CYS
1	A	282	LEU
1	A	284	GLU
1	A	299[A]	GLU
1	A	299[B]	GLU
1	A	305	THR
1	A	312	LYS
1	A	319	GLN
1	A	326	LEU
1	A	333	TYR
1	A	363	ASP
1	A	377	LYS
1	A	386	LEU
1	A	389	GLN
1	A	416	GLN
1	A	422	LEU
1	A	438	THR
1	A	474	LYS
1	A	511	ASP
1	A	523	LYS
1	A	540	GLU
1	A	541	GLU
1	A	577	SER
1	B	4	LYS
1	B	114	LYS
1	B	127	LYS
1	B	143	ARG
1	B	159	LYS
1	B	201	SER
1	B	244	CYS
1	B	280	LYS
1	B	288	CYS
1	B	294	LYS

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Mol	Chain	Res	Type
1	B	299	GLU
1	B	300	ASN
1	B	313	ASP
1	B	333	TYR
1	B	344	VAL
1	B	350	LYS
1	B	355	THR
1	B	363	ASP
1	B	369	TYR
1	B	371	THR
1	B	374	ASP
1	B	377	LYS
1	B	389	GLN
1	B	438	THR
1	B	442	SER
1	B	454	LEU
1	B	461	VAL
1	B	465	LYS
1	B	479	SER
1	B	539	THR
1	B	541	GLU
1	B	555	ASP
1	B	561	ASP
1	B	563	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA chains 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 22 ligands modelled in this entry, 7 are monoatomic - leaving 15 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	DIU	A	601	-	12,12,12	1.93	4 (33%)	17,17,17	1.33	1 (5%)
2	DIU	A	602	-	12,12,12	1.82	4 (33%)	17,17,17	0.89	1 (5%)
2	DIU	A	603	-	12,12,12	1.92	4 (33%)	17,17,17	1.03	1 (5%)
2	DIU	A	604	-	12,12,12	2.71	5 (41%)	17,17,17	0.89	0
3	PEG	A	605	-	6,6,6	0.70	0	5,5,5	0.91	0
4	1PE	A	606	-	15,15,15	0.84	0	14,14,14	1.20	1 (7%)
4	1PE	A	607	-	15,15,15	0.99	0	14,14,14	0.90	0
5	PGE	A	608	-	9,9,9	0.61	0	8,8,8	0.75	0
2	DIU	B	601	-	12,12,12	2.12	5 (41%)	17,17,17	1.08	1 (5%)
2	DIU	B	602	-	12,12,12	1.41	2 (16%)	17,17,17	0.84	0
2	DIU	B	603	-	12,12,12	2.18	5 (41%)	17,17,17	1.17	2 (11%)
2	DIU	B	604	-	12,12,12	2.79	7 (58%)	17,17,17	0.99	0
3	PEG	B	605	-	6,6,6	0.63	0	5,5,5	0.59	0
4	1PE	B	606	-	15,15,15	0.82	0	14,14,14	0.65	0
5	PGE	B	607	-	9,9,9	0.68	0	8,8,8	0.27	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	DIU	A	601	-	-	0/4/4/4	0/1/1/1
2	DIU	A	602	-	-	0/4/4/4	0/1/1/1
2	DIU	A	603	-	-	0/4/4/4	0/1/1/1
2	DIU	A	604	-	-	0/4/4/4	0/1/1/1
3	PEG	A	605	-	-	0/4/4/4	0/0/0/0
4	1PE	A	606	-	-	0/13/13/13	0/0/0/0
4	1PE	A	607	-	-	0/13/13/13	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	PGE	A	608	-	-	0/7/7/7	0/0/0/0
2	DIU	B	601	-	-	0/4/4/4	0/1/1/1
2	DIU	B	602	-	-	0/4/4/4	0/1/1/1
2	DIU	B	603	-	-	0/4/4/4	0/1/1/1
2	DIU	B	604	-	-	0/4/4/4	0/1/1/1
3	PEG	B	605	-	-	0/4/4/4	0/0/0/0
4	1PE	B	606	-	-	0/13/13/13	0/0/0/0
5	PGE	B	607	-	-	0/7/7/7	0/0/0/0

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	604	DIU	C1-C2	4.80	1.49	1.41
2	B	604	DIU	C4-C5	4.71	1.47	1.38
2	B	603	DIU	C6-C5	4.69	1.47	1.38
2	B	604	DIU	C1-C2	4.36	1.48	1.41
2	A	603	DIU	C6-C5	4.33	1.46	1.38
2	B	604	DIU	C2-C3	4.30	1.50	1.40
2	B	601	DIU	C2-C3	4.29	1.50	1.40
2	A	604	DIU	C4-C5	4.24	1.46	1.38
2	A	604	DIU	C2-C3	3.79	1.49	1.40
2	A	604	DIU	C3-I1	3.59	2.20	2.10
2	A	601	DIU	C1-C2	3.55	1.47	1.41
2	B	601	DIU	C1-C2	3.47	1.47	1.41
2	B	604	DIU	C3-I1	3.46	2.20	2.10
2	A	601	DIU	C2-C3	3.37	1.48	1.40
2	A	601	DIU	O1-C	3.19	1.32	1.23
2	A	604	DIU	C6-C5	3.06	1.44	1.38
2	A	602	DIU	C6-C5	3.04	1.44	1.38
2	A	602	DIU	C1-C2	3.03	1.46	1.41
2	B	604	DIU	C5-I2	2.84	2.18	2.10
2	B	602	DIU	C2-C3	2.78	1.46	1.40
2	B	603	DIU	C2-C3	2.67	1.46	1.40
2	B	601	DIU	C5-I2	-2.64	2.02	2.10
2	A	602	DIU	C2-C3	2.62	1.46	1.40
2	B	603	DIU	C3-I1	-2.58	2.02	2.10
2	A	603	DIU	C1-C2	2.55	1.45	1.41
2	A	603	DIU	C2-C3	2.54	1.46	1.40
2	B	601	DIU	C6-C5	2.51	1.43	1.38
2	B	602	DIU	C6-C5	2.43	1.43	1.38
2	B	603	DIU	C1-C2	2.32	1.45	1.41
2	B	603	DIU	C1-C	2.28	1.54	1.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	604	DIU	C6-C5	2.28	1.42	1.38
2	A	602	DIU	C4-C5	2.24	1.42	1.38
2	A	601	DIU	C6-C1	2.09	1.43	1.39
2	B	601	DIU	O3-C2	-2.08	1.32	1.37
2	B	604	DIU	C4-C3	2.04	1.44	1.39
2	A	603	DIU	C1-C	2.02	1.54	1.49

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	DIU	C3-C4-C5	4.29	122.90	119.30
4	A	606	1PE	OH5-C14-C24	-3.22	95.85	110.47
2	B	603	DIU	C3-C4-C5	3.06	121.87	119.30
2	B	601	DIU	C3-C4-C5	2.75	121.61	119.30
2	A	603	DIU	C3-C4-C5	2.55	121.44	119.30
2	A	602	DIU	C3-C4-C5	2.23	121.18	119.30
2	B	603	DIU	C6-C5-I2	2.05	122.26	119.38

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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 ⓘ

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	582/583 (99%)	0.10	15 (2%) 53 56	14, 40, 72, 103	0
1	B	582/583 (99%)	0.27	33 (5%) 23 24	17, 47, 92, 134	0
All	All	1164/1166 (99%)	0.18	48 (4%) 29 38	14, 43, 85, 134	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	361	ALA	7.3
1	B	363	ASP	5.3
1	B	316	LYS	4.9
1	A	388	LYS	4.5
1	B	304	LEU	4.1
1	A	3	HIS	4.1
1	B	360	CYS	4.0
1	B	367	ALA	3.9
1	B	117	PRO	3.6
1	B	362	LYS	3.6
1	B	379	LEU	3.4
1	B	2	THR	3.3
1	A	437	CYS	3.1
1	B	366	HIS	3.1
1	B	395	GLU	3.1
1	A	2	THR	3.0
1	B	538	ALA	3.0
1	B	310	GLU	2.8
1	B	369	TYR	2.8
1	B	122	LEU	2.8
1	A	299[A]	GLU	2.7
1	A	356	LEU	2.6
1	B	396	LYS	2.6
1	B	319	GLN	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	443	GLU	2.5
1	B	368	CYS	2.5
1	A	387	ILE	2.4
1	A	326	LEU	2.4
1	A	560	ALA	2.4
1	A	554	VAL	2.4
1	B	3	HIS	2.4
1	B	392	ASP	2.4
1	B	365	PRO	2.3
1	B	364	ASP	2.3
1	B	315	CYS	2.3
1	B	124	ASP	2.3
1	B	119	PRO	2.2
1	A	553	PHE	2.2
1	A	164	PHE	2.2
1	A	128	ALA	2.2
1	B	356	LEU	2.2
1	B	80	LEU	2.1
1	B	173	LYS	2.1
1	B	320	GLU	2.1
1	B	318	TYR	2.1
1	B	394	PHE	2.1
1	A	318	TYR	2.1
1	B	333	TYR	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 ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	PGE	B	607	10/10	0.46	4.80	62,65,68,69	0
4	1PE	B	606	16/16	0.36	4.01	52,69,81,81	0
2	DIU	B	602	12/12	0.36	2.90	48,50,91,123	12
4	1PE	A	607	16/16	0.30	1.23	18,59,66,67	0
5	PGE	A	608	10/10	0.26	1.10	40,44,47,49	0
6	CA	A	609	1/1	0.15	1.03	62,62,62,62	0
3	PEG	A	605	7/7	0.21	0.23	13,26,31,32	0
6	CA	B	608	1/1	0.20	0.11	70,70,70,70	0
6	CA	A	611	1/1	0.16	-0.37	74,74,74,74	0
4	1PE	A	606	16/16	0.16	-0.41	12,31,40,40	0
2	DIU	A	603	12/12	0.18	-0.62	61,64,76,120	0
3	PEG	B	605	7/7	0.18	-0.64	22,36,58,59	0
2	DIU	A	602	12/12	0.15	-0.85	48,53,96,98	0
2	DIU	B	604	12/12	0.16	-0.87	60,94,96,101	0
2	DIU	B	603	12/12	0.12	-1.14	57,61,78,126	0
2	DIU	B	601	12/12	0.15	-1.44	31,38,48,112	0
2	DIU	A	601	12/12	0.14	-1.49	25,36,44,116	0
6	CA	B	609	1/1	0.07	-1.77	60,60,60,60	0
2	DIU	A	604	12/12	0.12	-1.96	49,52,62,76	0
6	CA	A	610	1/1	0.08	-3.14	47,47,47,47	0
6	CA	B	610	1/1	0.08	-4.37	65,65,65,65	0
6	CA	A	612	1/1	0.10	-9.66	48,48,48,48	0

6.5 Other polymers ⓘ

There are no such residues in this entry.