



Full wwPDB X-ray Structure Validation Report i

Feb 27, 2014 – 02:39 PM GMT

PDB ID : 1RDT
Title : Crystal Structure of a new rexinoid bound to the RXRalpha ligand binding domain in the RXRalpha/PPARgamma heterodimer
Authors : Haffner, C.D.; Lenhard, J.M.; Miller, A.B.; McDougald, D.L.; Dwornik, K.; Ittoop, O.R.; Gampe Jr., R.T.; Xu, H.E.; Blanchard, S.; Montana, V.G.
Deposited on : 2003-11-06
Resolution : 2.40 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>

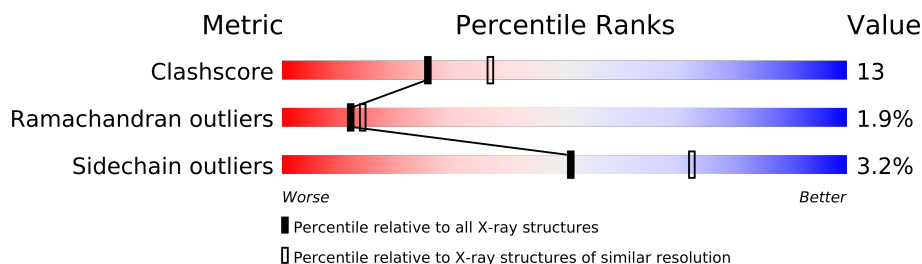
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 21963
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 2.40 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))
Clashscore	79885	2789 (2.40-2.40)
Ramachandran outliers	78287	2736 (2.40-2.40)
Sidechain outliers	78261	2737 (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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	242	
2	B	25	
3	D	284	
4	E	23	

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 4027 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 Retinoic acid receptor RXR-alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	211	Total	C	N	O	S	0	0	0
			1644	1054	285	295	10			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	221	MET	-	CLONING ARTIFACT	UNP P19793
A	222	LYS	-	CLONING ARTIFACT	UNP P19793
A	223	LYS	-	CLONING ARTIFACT	UNP P19793
A	224	GLY	-	CLONING ARTIFACT	UNP P19793

- Molecule 2 is a protein called LxxLL motif coactivator.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	9	Total	C	N	O	0	0	0
			77	50	17	10			

- Molecule 3 is a protein called Peroxisome proliferator activated receptor gamma.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	253	Total	C	N	O	S	0	0	0
			1998	1288	327	376	7			

There are 13 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	194	MET	-	CLONING ARTIFACT	UNP P37231
D	195	LYS	-	CLONING ARTIFACT	UNP P37231
D	196	LYS	-	CLONING ARTIFACT	UNP P37231
D	197	GLY	-	CLONING ARTIFACT	UNP P37231
D	198	HIS	-	CLONING ARTIFACT	UNP P37231
D	199	HIS	-	CLONING ARTIFACT	UNP P37231

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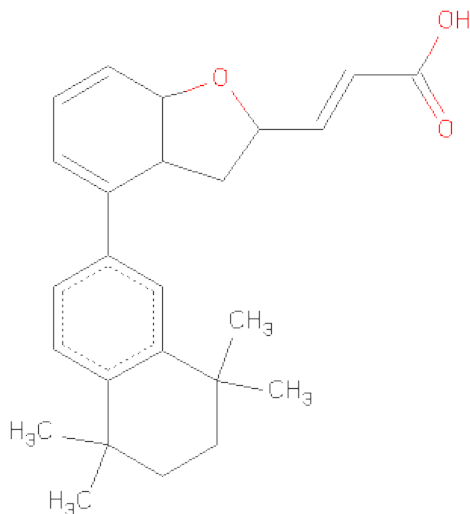
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Chain	Residue	Modelled	Actual	Comment	Reference
D	200	HIS	-	CLONING ARTIFACT	UNP P37231
D	201	HIS	-	CLONING ARTIFACT	UNP P37231
D	202	HIS	-	CLONING ARTIFACT	UNP P37231
D	203	HIS	-	CLONING ARTIFACT	UNP P37231
D	204	GLY	-	CLONING ARTIFACT	UNP P37231
D	205	ARG	-	CLONING ARTIFACT	UNP P37231
D	206	ALA	-	CLONING ARTIFACT	UNP P37231

- Molecule 4 is a protein called LxxLL motif coactivator.

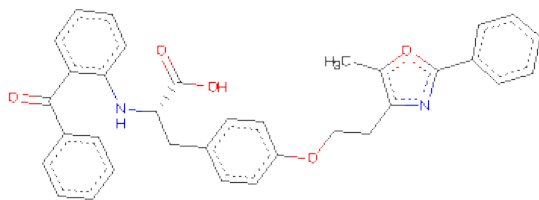
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	E	7	Total	C	N	O	0	0	0
			34	20	7	7			

- Molecule 5 is (S)-(2E)-3[4-(5,5,8,8-TETRAMETHYL-5,6,7,8-TETRAHYDRO-2-NAPHTHALENYL)TETRAHYDRO-1-BENZOFURAN-2-YL]-2-PROPENOICACID (three-letter code: L79) (formula: C₂₅H₃₀O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	O		0	0
			28	25	3			

- Molecule 6 is 2-(2-BENZOYL-PHENYLAMINO)-3-{4-[2-(5-METHYL-2-PHENYL-OXAZOL-4-YL)-ETHOXY]-PHENYL}-PROPIONICACID (three-letter code: 570) (formula: C₃₄H₃₀N₂O₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	D	1	Total	C	N	O	0	0
			41	34	2	5		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	72	Total	O	0	0
			72	72		
7	B	2	Total	O	0	0
			2	2		
7	D	123	Total	O	0	0
			123	123		
7	E	8	Total	O	0	0
			8	8		

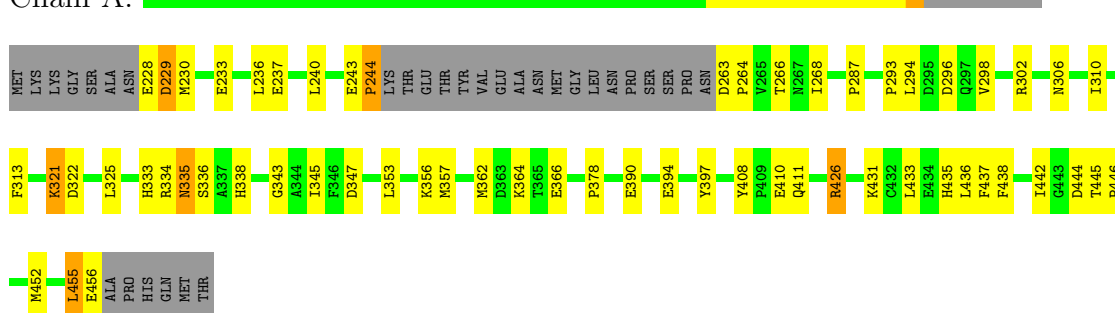
3 Residue-property plots

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.

Note EDS was not executed.

- Molecule 1: Retinoic acid receptor RXR-alpha

Chain A:



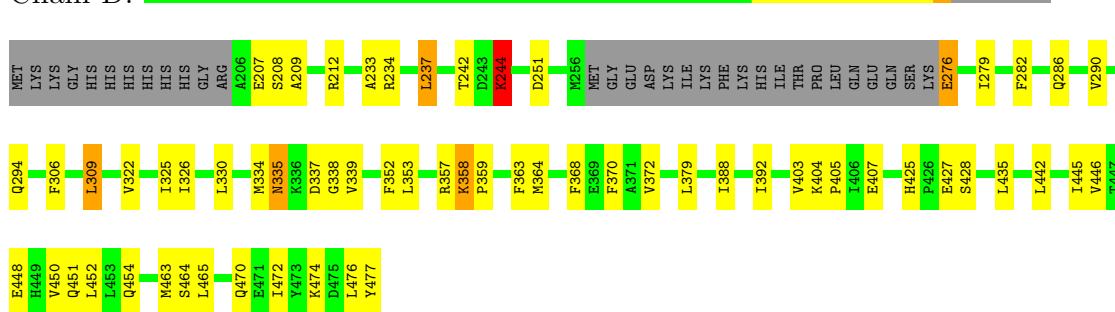
- Molecule 2: LxxLL motif coactivator

Chain B:



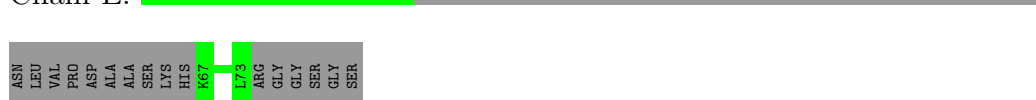
- Molecule 3: Peroxisome proliferator activated receptor gamma

Chain D:



- Molecule 4: LxxLL motif coactivator

Chain E:



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	49.69Å 54.58Å 211.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.40	Depositor
% Data completeness (in resolution range)	97.1 (20.00-2.40)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNX 2000	Depositor
R, R_{free}	0.221 , 0.259	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4027	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 570, L79

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	0.26	0/1676	0.52	0/2261
2	B	0.27	0/78	0.54	0/104
3	D	0.26	0/2031	0.50	0/2739
4	E	0.43	0/33	0.33	0/44
All	All	0.27	0/3818	0.51	0/5148

There are no bond length outliers.

There are no bond angle outliers.

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	1644	0	1662	48	2
2	B	77	0	80	7	2
3	D	1998	0	2028	53	0
4	E	34	0	12	0	0
5	A	28	0	26	2	0
6	D	41	0	29	0	0
7	A	72	0	0	2	0
7	B	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	D	123	0	0	3	0
7	E	8	0	0	0	0
All	All	4027	0	3837	103	2

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 13.

All (103) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:D:358:LYS:HB3	3:D:359:PRO:HD2	1.30	1.06
3:D:242:THR:HG22	7:D:566:HOH:O	1.55	1.04
3:D:358:LYS:HB3	3:D:359:PRO:CD	1.90	1.02
1:A:233:GLU:HG3	7:A:103:HOH:O	0.84	1.00
3:D:357:ARG:NH2	3:D:358:LYS:HD2	2.00	0.76
3:D:358:LYS:CB	3:D:359:PRO:HD2	2.15	0.75
3:D:335:ASN:ND2	3:D:338:GLY:H	1.83	0.74
1:A:335:ASN:ND2	1:A:336:SER:H	1.86	0.74
3:D:335:ASN:C	3:D:335:ASN:HD22	1.92	0.73
1:A:345:ILE:HD13	1:A:431:LYS:HD2	1.70	0.72
1:A:243:GLU:HB2	1:A:244:PRO:C	2.09	0.72
1:A:426:ARG:HH11	1:A:426:ARG:HB2	1.55	0.71
1:A:333:HIS:HB2	1:A:335:ASN:HD21	1.56	0.70
1:A:335:ASN:HD22	1:A:336:SER:H	1.39	0.70
3:D:357:ARG:HH21	3:D:358:LYS:HD2	1.56	0.70
1:A:233:GLU:CG	7:A:103:HOH:O	1.65	0.67
3:D:474:LYS:HE3	7:D:568:HOH:O	1.95	0.67
3:D:450:VAL:O	3:D:454:GLN:HG2	1.96	0.65
3:D:448:GLU:O	3:D:452:LEU:HD13	1.97	0.64
1:A:335:ASN:HD22	1:A:335:ASN:N	1.94	0.64
1:A:298:VAL:HG13	2:B:690:LEU:HD23	1.81	0.63
1:A:306:ASN:OD1	5:A:463:L79:H10	1.99	0.61
3:D:207:GLU:HG3	3:D:209:ALA:H	1.67	0.60
3:D:442:LEU:O	3:D:446:VAL:HG23	2.02	0.59
3:D:335:ASN:ND2	3:D:337:ASP:H	1.99	0.59
1:A:233:GLU:O	1:A:237:GLU:HG3	2.02	0.59
2:B:688:LYS:O	2:B:692:ARG:HB2	2.03	0.58
1:A:228:GLU:O	1:A:229:ASP:HB2	2.03	0.57
3:D:363:PHE:CZ	3:D:452:LEU:HB3	2.39	0.57
3:D:335:ASN:HD21	3:D:338:GLY:N	2.03	0.57
1:A:335:ASN:ND2	1:A:336:SER:N	2.52	0.57
3:D:370:PHE:HB2	3:D:445:ILE:HD11	1.86	0.56
1:A:333:HIS:HB2	1:A:335:ASN:ND2	2.21	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:D:325:ILE:HD11	3:D:392:ILE:HG13	1.89	0.54
3:D:379:LEU:HD11	3:D:435:LEU:HD13	1.89	0.54
3:D:335:ASN:HD21	3:D:338:GLY:H	1.50	0.54
3:D:233:ALA:O	3:D:237:LEU:HD22	2.09	0.53
3:D:427:GLU:HG2	7:D:573:HOH:O	2.08	0.53
1:A:338:HIS:CE1	1:A:343:GLY:HA3	2.43	0.53
3:D:476:LEU:HG	3:D:477:TYR:HD1	1.74	0.53
1:A:436:LEU:HD11	5:A:463:L79:H9	1.91	0.52
1:A:310:ILE:HA	1:A:313:PHE:CE2	2.44	0.52
1:A:353:LEU:O	1:A:357:MET:HG3	2.08	0.52
3:D:286:GLN:OE1	3:D:465:LEU:HD12	2.09	0.51
1:A:230:MET:SD	1:A:287:PRO:HG2	2.50	0.51
1:A:378:PRO:HG3	1:A:390:GLU:OE1	2.09	0.51
3:D:425:HIS:HB3	3:D:428:SER:OG	2.11	0.51
3:D:335:ASN:C	3:D:335:ASN:ND2	2.62	0.51
3:D:244:LYS:HB2	3:D:244:LYS:NZ	2.25	0.50
3:D:306:PHE:O	3:D:309:LEU:HB2	2.12	0.50
3:D:306:PHE:HA	3:D:309:LEU:HD22	1.94	0.50
3:D:403:VAL:O	3:D:407:GLU:HG3	2.12	0.50
1:A:228:GLU:O	1:A:228:GLU:HG3	2.12	0.49
3:D:353:LEU:HD13	3:D:364:MET:HG3	1.93	0.49
3:D:242:THR:HG23	3:D:242:THR:O	2.12	0.49
3:D:465:LEU:O	3:D:470:GLN:NE2	2.46	0.48
1:A:394:GLU:HA	1:A:397:TYR:CD2	2.48	0.48
1:A:335:ASN:HD22	1:A:336:SER:N	2.08	0.48
1:A:293:PRO:HD2	1:A:296:ASP:OD2	2.13	0.48
1:A:335:ASN:N	1:A:335:ASN:ND2	2.61	0.48
3:D:451:GLN:O	3:D:454:GLN:HB2	2.14	0.47
3:D:476:LEU:HG	3:D:477:TYR:CD1	2.49	0.47
3:D:207:GLU:HG3	3:D:209:ALA:N	2.30	0.47
1:A:408:TYR:HB3	1:A:411:GLN:HG3	1.97	0.47
3:D:234:ARG:HG2	3:D:234:ARG:HH11	1.80	0.47
3:D:404:LYS:HB3	3:D:405:PRO:HD3	1.97	0.46
1:A:233:GLU:H	1:A:233:GLU:CD	2.19	0.46
1:A:426:ARG:NH1	1:A:426:ARG:HB2	2.27	0.46
3:D:325:ILE:HD13	3:D:388:ILE:HG23	1.98	0.45
3:D:279:ILE:O	3:D:282:PHE:HB3	2.16	0.45
1:A:228:GLU:O	1:A:229:ASP:CB	2.63	0.45
3:D:325:ILE:HD11	3:D:392:ILE:CG1	2.46	0.45
1:A:302:ARG:CZ	2:B:687:HIS:NE2	2.80	0.45
1:A:435:HIS:O	1:A:438:PHE:HB3	2.17	0.45
3:D:208:SER:O	3:D:212:ARG:HG2	2.17	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:362:MET:HA	1:A:366:GLU:OE2	2.17	0.44
1:A:263:ASP:O	1:A:266:THR:HB	2.17	0.44
1:A:452:MET:O	1:A:456:GLU:HG3	2.17	0.44
3:D:244:LYS:O	3:D:244:LYS:HG3	2.17	0.44
1:A:338:HIS:ND1	1:A:343:GLY:HA3	2.33	0.44
2:B:691:HIS:CE1	2:B:695:GLN:NE2	2.86	0.44
1:A:433:LEU:HD22	1:A:437:PHE:CE2	2.53	0.43
1:A:455:LEU:HD12	1:A:455:LEU:HA	1.84	0.43
3:D:322:VAL:HG21	3:D:472:ILE:HD13	2.00	0.43
1:A:302:ARG:HH11	1:A:302:ARG:HG2	1.83	0.43
3:D:276:GLU:OE1	3:D:276:GLU:HA	2.18	0.43
3:D:334:MET:HG2	3:D:339:VAL:HB	2.00	0.43
3:D:322:VAL:O	3:D:326:ILE:HG13	2.19	0.42
3:D:251:ASP:HA	3:D:352:PHE:CD1	2.55	0.42
1:A:313:PHE:C	1:A:313:PHE:CD1	2.93	0.42
1:A:356:LYS:NZ	3:D:403:VAL:HG11	2.34	0.42
1:A:321:LYS:HB3	1:A:322:ASP:H	1.70	0.41
1:A:240:LEU:HA	1:A:243:GLU:HG3	2.02	0.41
3:D:363:PHE:CE1	3:D:452:LEU:HB3	2.56	0.41
2:B:691:HIS:CE1	2:B:695:GLN:HE22	2.39	0.41
1:A:302:ARG:NH1	2:B:687:HIS:CD2	2.89	0.41
1:A:294:LEU:HD21	2:B:691:HIS:HD2	1.86	0.41
3:D:330:LEU:O	3:D:334:MET:HG3	2.20	0.41
1:A:334:ARG:HD2	1:A:347:ASP:OD1	2.20	0.41
1:A:243:GLU:CD	1:A:364:LYS:NZ	2.75	0.41
3:D:290:VAL:HG12	3:D:294:GLN:HE21	1.87	0.40
1:A:264:PRO:O	1:A:268:ILE:HG13	2.20	0.40
3:D:368:PHE:O	3:D:372:VAL:HG23	2.21	0.40

All (2) 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	Distance(Å)	Clash(Å)
1:A:240:LEU:CD1	2:B:692:ARG:NE[3_645]	1.66	0.54
1:A:240:LEU:CD1	2:B:692:ARG:CD[3_645]	1.76	0.44

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	207/242 (86%)	189 (91%)	13 (6%)	5 (2%)	9	9
2	B	7/25 (28%)	7 (100%)	0	0	100	100
3	D	249/284 (88%)	236 (95%)	9 (4%)	4 (2%)	14	18
4	E	5/23 (22%)	5 (100%)	0	0	100	100
All	All	468/574 (82%)	437 (93%)	22 (5%)	9 (2%)	12	14

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	321	LYS
1	A	446	PRO
3	D	358	LYS
1	A	444	ASP
3	D	244	LYS
3	D	464	SER
1	A	442	ILE
3	D	463	MET
1	A	445	THR

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	175/208 (84%)	167 (95%)	8 (5%)	37	55
2	B	8/24 (33%)	8 (100%)	0	100	100
3	D	219/254 (86%)	214 (98%)	5 (2%)	63	82
All	All	402/486 (83%)	389 (97%)	13 (3%)	51	72

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	229	ASP
1	A	236	LEU

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Mol	Chain	Res	Type
1	A	244	PRO
1	A	325	LEU
1	A	335	ASN
1	A	410	GLU
1	A	426	ARG
1	A	455	LEU
3	D	237	LEU
3	D	244	LYS
3	D	276	GLU
3	D	309	LEU
3	D	335	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	267	ASN
1	A	270	GLN
1	A	275	GLN
1	A	333	HIS
1	A	335	ASN
2	B	691	HIS
2	B	695	GLN
3	D	294	GLN
3	D	308	ASN
3	D	335	ASN
3	D	451	GLN

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 ⓘ

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
5	L79	A	463	-	31,31,31	3.74	17 (54%)	48,48,48	3.23	17 (35%)
6	570	D	478	-	45,45,45	2.68	11 (24%)	59,61,61	1.37	7 (11%)

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
5	L79	A	463	-	3/3/7/12	0/9/50/50	0/0/4/4
6	570	D	478	-	-	0/30/30/30	0/5/5/5

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	D	478	570	C3I-C3G	-14.37	1.34	1.47
5	A	463	L79	C3-C2	-10.69	1.39	1.54
5	A	463	L79	C9-C8	8.62	1.53	1.36
5	A	463	L79	C10-C11	7.09	1.53	1.34
5	A	463	L79	C-C3	-6.93	1.39	1.54
5	A	463	L79	C-C1	-5.56	1.39	1.54
5	A	463	L79	O-C2	-4.89	1.37	1.44
5	A	463	L79	C10-C9	4.79	1.53	1.40
6	D	478	570	C1F-C1A	-3.20	1.35	1.41
6	D	478	570	C1F-C1G	2.98	1.56	1.49
6	D	478	570	C1E-C1F	-2.89	1.34	1.39
5	A	463	L79	C13-C12	-2.89	1.35	1.39
5	A	463	L79	O-C1	-2.80	1.37	1.44
5	A	463	L79	C13-C14	-2.73	1.35	1.39
5	A	463	L79	C24-C19	-2.68	1.35	1.39
6	D	478	570	C1B-C1A	-2.60	1.35	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	463	L79	C19-C14	-2.57	1.35	1.40
6	D	478	570	C1M-C1H	-2.52	1.34	1.39
5	A	463	L79	C25-C12	-2.51	1.34	1.39
6	D	478	570	C1I-C1H	-2.43	1.35	1.39
5	A	463	L79	C1-C4	-2.41	1.37	1.50
6	D	478	570	C1A-N	2.40	1.43	1.37
5	A	463	L79	O1-C6	-2.32	1.23	1.30
5	A	463	L79	C25-C24	-2.14	1.34	1.38
6	D	478	570	C3J-C3I	-2.14	1.34	1.39
6	D	478	570	CE2-CD2	-2.06	1.34	1.38
5	A	463	L79	C12-C8	2.05	1.51	1.48
6	D	478	570	C3N-C3I	-2.02	1.34	1.39

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	463	L79	O-C2-C11	11.49	126.85	110.26
5	A	463	L79	O-C2-C3	7.72	108.77	101.07
5	A	463	L79	O-C1-C4	7.33	127.84	109.12
5	A	463	L79	C3-C2-C11	6.09	124.38	114.78
5	A	463	L79	C2-C11-C10	-5.27	113.69	121.32
5	A	463	L79	C3-C8-C9	-4.88	112.45	119.72
6	D	478	570	O3F-C3G-C3I	4.68	124.06	119.23
5	A	463	L79	C2-C3-C8	4.51	124.19	111.05
5	A	463	L79	C16-C15-C14	4.42	117.50	110.38
5	A	463	L79	C22-C20-C19	4.19	117.14	110.38
5	A	463	L79	C-C3-C2	4.08	107.87	101.45
6	D	478	570	C1F-C1G-C1H	3.60	125.89	119.65
5	A	463	L79	C12-C8-C9	-3.49	115.73	121.93
6	D	478	570	C1A-C1F-C1G	3.13	124.25	121.93
5	A	463	L79	C12-C8-C3	3.11	124.10	119.07
6	D	478	570	O1G-C1G-C1H	-2.94	115.31	120.02
6	D	478	570	O3F-C3G-N3H	-2.82	109.08	117.79
5	A	463	L79	C3-C-C1	2.70	106.41	102.50
5	A	463	L79	C-C3-C8	2.58	127.94	117.81
5	A	463	L79	O-C1-C	2.53	109.60	103.58
6	D	478	570	CG-CB-CA	2.35	119.93	113.31
5	A	463	L79	C10-C9-C8	-2.13	118.52	121.14
6	D	478	570	O3F-C3D-C3E	2.09	122.82	117.43
5	A	463	L79	C25-C12-C13	-2.04	116.67	119.24

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
5	A	463	L79	C3
5	A	463	L79	C2
5	A	463	L79	C1

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 ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.