



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 26, 2014 – 09:59 PM GMT

PDB ID : 1R1K
Title : Crystal structure of the ligand-binding domains of the heterodimer EcR/USP bound to ponasterone A
Authors : Billas, I.M.L.; Iwema, T.; Garnier, J.-M.; Mitschler, A.; Rochel, N.; Moras, D.; Structural Proteomics in Europe (SPINE)
Deposited on : 2003-09-24
Resolution : 2.90 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>

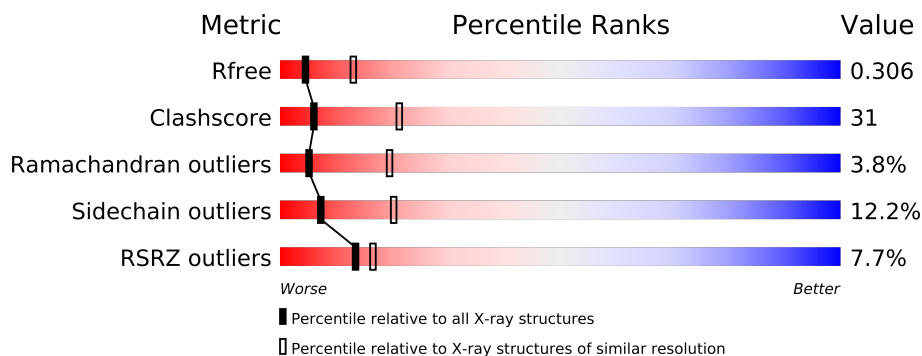
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.90 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	1053 (2.90-2.90)
Clashscore	79885	1326 (2.90-2.90)
Ramachandran outliers	78287	1290 (2.90-2.90)
Sidechain outliers	78261	1292 (2.90-2.90)
RSRZ outliers	66119	1054 (2.90-2.90)

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	D	266	
2	A	263	

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

Mol	Type	Chain	Res	Geometry	Electron density
3	P1A	D	3000	-	X
4	EPH	A	4000	-	X

2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 3887 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.

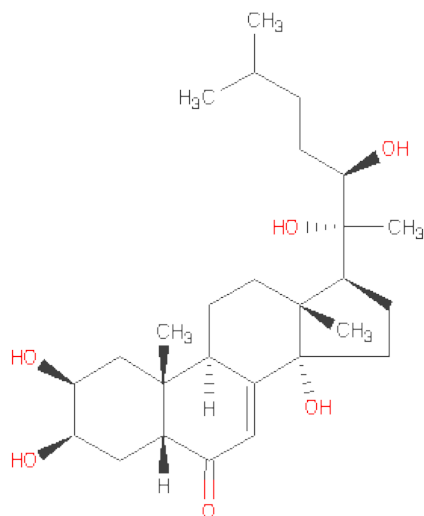
- Molecule 1 is a protein called Ecdysone receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	236	Total	C	N	O	S	0	0	0
			1872	1194	316	346	16			

- Molecule 2 is a protein called ULTRASPIRACLE PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	241	Total	C	N	O	S	0	0	0
			1925	1233	335	345	12			

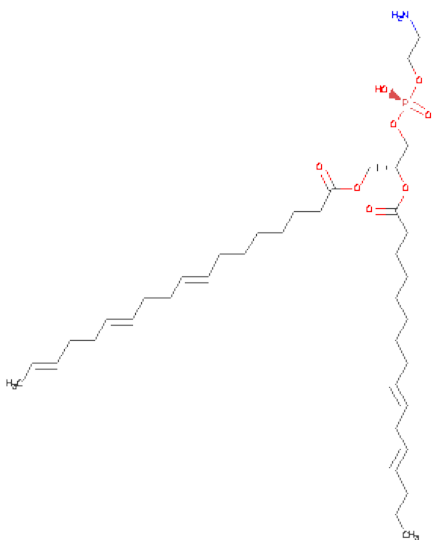
- Molecule 3 is 2,3,14,20,22-PENTAHYDROXYCHOLEST-7-EN-6-ONE (three-letter code: P1A) (formula: $C_{27}H_{44}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	D	1	Total	C	O	0	0
			33	27	6		

- Molecule 4 is L-ALPHA-PHOSPHATIDYL-BETA-OLEOYL-GAMMA-PALMITOYL-PH

OSPHATIDYLETHANOLAMINE (three-letter code: EPH) (formula: $C_{39}H_{68}NO_8P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			49	39	1	8	1		

- Molecule 5 is water.

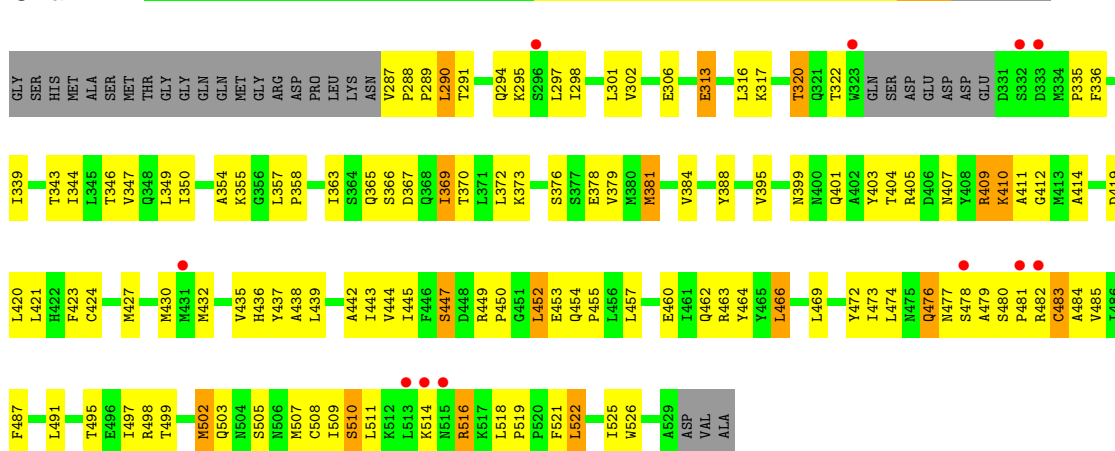
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	5	Total	O	0	0
			5	5		
5	D	3	Total	O	0	0
			3	3		

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.

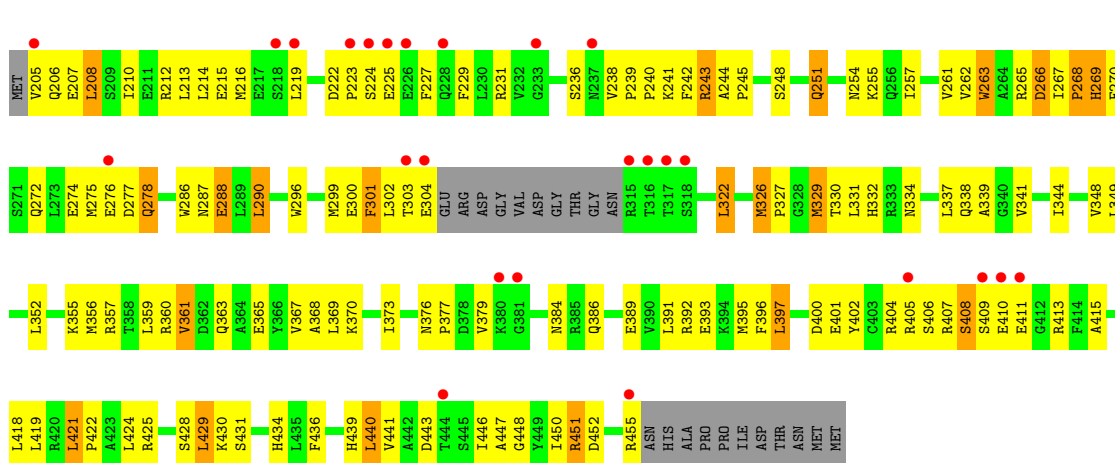
• Molecule 1: Ecdysone receptor

Chain D:



• Molecule 2: ULTRASPIRACLE PROTEIN

Chain A:



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	57.40Å 57.40Å 302.70Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.90 19.86 – 2.90	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-2.90) 99.6 (19.86-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.66 (at 2.88Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.243 , 0.304 0.247 , 0.306	Depositor DCC
R_{free} test set	1388 reflections (10.18%)	DCC
Wilson B-factor (Å ²)	78.2	Xtriage
Anisotropy	0.512	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 58.5	EDS
Estimated twinning fraction	0.039 for -h,-k,l	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 13666 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3887	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.90% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: P1A, EPH

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	D	0.55	0/1905	0.72	0/2579
2	A	0.51	0/1960	0.70	0/2646
All	All	0.53	0/3865	0.71	0/5225

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	D	1872	0	1880	130	0
2	A	1925	0	1970	119	1
3	D	33	0	43	3	0
4	A	49	0	67	5	0
5	A	5	0	0	1	0
5	D	3	0	0	0	0
All	All	3887	0	3960	242	1

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 31.

All (242) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:480:SER:HB2	1:D:481:PRO:HD2	1.51	0.93
1:D:403:TYR:HA	1:D:407:ASN:HD22	1.35	0.89
1:D:478:SER:O	1:D:480:SER:N	2.08	0.87
1:D:453:GLU:C	1:D:455:PRO:HD3	1.95	0.86
2:A:262:VAL:HG23	2:A:265:ARG:NH2	1.92	0.85
1:D:346:THR:O	1:D:350:ILE:HG13	1.78	0.82
2:A:376:ASN:HD22	2:A:379:VAL:HG23	1.43	0.82
2:A:451:ARG:HD2	2:A:452:ASP:N	1.96	0.81
1:D:365:GLN:O	1:D:369:ILE:HG22	1.81	0.80
2:A:431:SER:HB2	4:A:4000:EPH:H142	1.64	0.79
1:D:335:PRO:HB2	1:D:411:ALA:HB2	1.65	0.79
1:D:452:LEU:HB3	1:D:455:PRO:HG3	1.68	0.76
1:D:480:SER:CB	1:D:481:PRO:HD2	2.14	0.76
1:D:403:TYR:HA	1:D:407:ASN:ND2	2.01	0.75
1:D:502:MET:CE	2:A:429:LEU:HB3	2.17	0.74
2:A:406:SER:O	2:A:407:ARG:HG3	1.89	0.73
2:A:334:ASN:O	2:A:338:GLN:HG2	1.91	0.71
1:D:290:LEU:HB2	1:D:294:GLN:HE21	1.56	0.70
2:A:262:VAL:HG23	2:A:265:ARG:HH21	1.54	0.69
2:A:363:GLN:O	2:A:367:VAL:HG13	1.92	0.69
2:A:341:VAL:HG12	2:A:434:HIS:ND1	2.08	0.69
2:A:376:ASN:ND2	2:A:379:VAL:HG23	2.08	0.69
1:D:499:THR:HA	2:A:429:LEU:HD21	1.75	0.69
2:A:389:GLU:O	2:A:393:GLU:HG3	1.93	0.68
2:A:275:MET:HA	2:A:278:GLN:HB2	1.75	0.68
1:D:346:THR:HG22	1:D:350:ILE:HD11	1.76	0.68
1:D:447:SER:HA	1:D:462:GLN:HE22	1.59	0.68
1:D:474:LEU:HD23	1:D:483:CYS:SG	2.34	0.67
1:D:378:GLU:HB3	1:D:497:ILE:CG2	2.25	0.67
1:D:399:ASN:HD21	1:D:401:GLN:HG3	1.58	0.67
1:D:313:GLU:H	1:D:313:GLU:CD	1.99	0.66
1:D:502:MET:HE1	2:A:429:LEU:HB3	1.75	0.66
1:D:298:ILE:O	1:D:302:VAL:HG23	1.95	0.66
2:A:263:TRP:CZ2	2:A:370:LYS:HE2	2.31	0.65
1:D:444:VAL:O	1:D:447:SER:HB2	1.96	0.65
1:D:499:THR:O	1:D:503:GLN:HG3	1.95	0.65
2:A:206:GLN:NE2	2:A:212:ARG:HD3	2.11	0.65
2:A:213:LEU:HA	2:A:216:MET:SD	2.37	0.65
2:A:270:PHE:CE2	2:A:278:GLN:HG2	2.31	0.65
2:A:222:ASP:O	2:A:224:SER:N	2.30	0.65
1:D:290:LEU:CD1	1:D:295:LYS:HB2	2.26	0.64
1:D:357:LEU:HD11	1:D:445:ILE:HD13	1.77	0.64

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:447:SER:HA	1:D:462:GLN:NE2	2.13	0.64
2:A:434:HIS:HB3	4:A:4000:EPH:H112	1.79	0.64
2:A:286:TRP:O	2:A:290:LEU:HB2	1.99	0.62
2:A:369:LEU:O	2:A:373:ILE:HG13	1.99	0.62
2:A:344:ILE:HD11	2:A:431:SER:HB3	1.82	0.62
2:A:287:ASN:ND2	2:A:428:SER:HB2	2.14	0.62
2:A:268:PRO:O	2:A:269:HIS:HB2	1.99	0.62
1:D:381:MET:HB3	1:D:497:ILE:HD12	1.82	0.61
1:D:399:ASN:HD21	1:D:401:GLN:CG	2.13	0.61
1:D:409:ARG:HA	1:D:414:ALA:HB2	1.82	0.60
1:D:369:ILE:HD11	1:D:373:LYS:NZ	2.17	0.60
1:D:297:LEU:CD1	1:D:358:PRO:HG2	2.32	0.60
2:A:214:LEU:HD12	2:A:367:VAL:HG21	1.81	0.60
2:A:213:LEU:O	2:A:216:MET:N	2.31	0.60
1:D:423:PHE:CZ	1:D:497:ILE:HD11	2.37	0.59
2:A:216:MET:HE1	2:A:267:ILE:HA	1.85	0.59
1:D:363:ILE:CG2	1:D:367:ASP:HB2	2.31	0.59
1:D:287:VAL:N	1:D:288:PRO:CD	2.65	0.59
2:A:360:ARG:HG2	2:A:360:ARG:O	2.03	0.58
1:D:499:THR:OG1	2:A:425:ARG:HD2	2.02	0.58
1:D:469:LEU:HG	1:D:473:ILE:HD11	1.85	0.58
1:D:290:LEU:O	1:D:290:LEU:HD12	2.04	0.58
2:A:270:PHE:O	2:A:278:GLN:NE2	2.37	0.58
1:D:404:THR:H	1:D:407:ASN:ND2	2.01	0.58
1:D:290:LEU:HB2	1:D:294:GLN:NE2	2.17	0.57
2:A:244:ALA:N	2:A:245:PRO:HD2	2.19	0.57
2:A:299:MET:HG2	2:A:357:ARG:HD3	1.87	0.57
1:D:453:GLU:O	1:D:455:PRO:HD3	2.04	0.57
1:D:409:ARG:HD2	1:D:414:ALA:HB2	1.85	0.57
2:A:238:VAL:HG13	4:A:4000:EPH:H62	1.86	0.57
1:D:290:LEU:HD13	1:D:295:LYS:HB2	1.87	0.56
1:D:480:SER:O	1:D:483:CYS:HB2	2.06	0.56
1:D:476:GLN:HE21	1:D:476:GLN:HA	1.70	0.56
1:D:502:MET:HE2	2:A:429:LEU:HB3	1.88	0.56
1:D:376:SER:HG	1:D:526:TRP:HE1	1.54	0.55
1:D:511:LEU:HB2	1:D:518:LEU:HD21	1.89	0.55
1:D:511:LEU:CB	1:D:518:LEU:HD21	2.37	0.55
1:D:290:LEU:HD11	1:D:295:LYS:HB2	1.88	0.55
1:D:399:ASN:ND2	1:D:401:GLN:CG	2.70	0.55
1:D:320:THR:O	1:D:320:THR:HG23	2.07	0.55
2:A:331:LEU:HD13	4:A:4000:EPH:H363	1.88	0.54
2:A:287:ASN:HD22	2:A:428:SER:HB2	1.73	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:A:288:GLU:HG2	2:A:424:LEU:HG	1.90	0.54
1:D:290:LEU:HD21	1:D:295:LYS:HD3	1.90	0.54
1:D:297:LEU:HD13	1:D:358:PRO:HG2	1.90	0.53
2:A:275:MET:O	2:A:276:GLU:C	2.47	0.53
2:A:448:GLY:O	2:A:452:ASP:N	2.38	0.53
2:A:404:ARG:O	2:A:408:SER:HB2	2.09	0.53
1:D:399:ASN:ND2	1:D:401:GLN:HG2	2.23	0.53
2:A:241:LYS:HE3	2:A:242:PHE:CE1	2.44	0.53
1:D:435:VAL:HG21	1:D:472:TYR:CE1	2.43	0.53
1:D:369:ILE:HD11	1:D:373:LYS:HZ2	1.72	0.53
1:D:509:ILE:C	1:D:511:LEU:H	2.12	0.53
1:D:518:LEU:O	1:D:519:PRO:C	2.45	0.53
2:A:300:GLU:O	2:A:302:LEU:N	2.42	0.52
2:A:392:ARG:HG2	2:A:396:PHE:CE2	2.43	0.52
2:A:332:HIS:CD2	2:A:334:ASN:H	2.27	0.52
2:A:269:HIS:HA	2:A:272:GLN:HG2	1.92	0.52
1:D:290:LEU:CB	1:D:294:GLN:NE2	2.73	0.52
1:D:313:GLU:N	1:D:313:GLU:CD	2.63	0.52
2:A:361:VAL:HA	2:A:365:GLU:OE2	2.09	0.52
1:D:367:ASP:O	1:D:370:THR:HB	2.09	0.52
2:A:300:GLU:C	2:A:302:LEU:H	2.13	0.52
1:D:499:THR:HA	2:A:429:LEU:CD2	2.40	0.51
2:A:269:HIS:HB2	2:A:391:LEU:HD21	1.93	0.51
1:D:480:SER:CB	1:D:481:PRO:CD	2.85	0.50
2:A:208:LEU:CD2	2:A:395:MET:SD	3.00	0.50
1:D:290:LEU:CB	1:D:294:GLN:HE21	2.24	0.50
2:A:265:ARG:HA	2:A:270:PHE:CD2	2.46	0.50
1:D:449:ARG:H	1:D:452:LEU:HD22	1.76	0.50
2:A:296:TRP:O	2:A:299:MET:HB2	2.12	0.50
2:A:225:GLU:HA	2:A:225:GLU:OE2	2.11	0.50
1:D:378:GLU:O	1:D:381:MET:HB2	2.11	0.49
1:D:381:MET:HB3	1:D:497:ILE:CD1	2.42	0.49
2:A:326:MET:O	2:A:327:PRO:C	2.48	0.49
2:A:407:ARG:O	2:A:410:GLU:HG2	2.12	0.49
2:A:208:LEU:HB2	2:A:268:PRO:HB2	1.94	0.49
1:D:344:ILE:HG12	1:D:522:LEU:HD22	1.93	0.49
2:A:447:ALA:O	2:A:450:ILE:HB	2.12	0.49
1:D:454:GLN:O	1:D:457:LEU:HB3	2.13	0.49
1:D:476:GLN:HE21	1:D:476:GLN:CA	2.25	0.49
2:A:418:LEU:HA	2:A:421:LEU:HD22	1.95	0.49
1:D:355:LYS:C	1:D:357:LEU:H	2.16	0.49
2:A:208:LEU:HD21	2:A:395:MET:SD	2.52	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:297:LEU:HD22	1:D:464:TYR:HE2	1.77	0.49
2:A:262:VAL:O	2:A:262:VAL:HG22	2.12	0.49
2:A:268:PRO:O	2:A:269:HIS:CB	2.60	0.49
1:D:476:GLN:NE2	1:D:476:GLN:HA	2.28	0.49
2:A:395:MET:HE3	2:A:395:MET:HA	1.94	0.49
1:D:363:ILE:HG22	1:D:367:ASP:HB2	1.94	0.49
2:A:210:ILE:O	2:A:214:LEU:HD13	2.13	0.48
1:D:379:VAL:C	1:D:381:MET:N	2.65	0.48
1:D:511:LEU:HD23	1:D:514:LYS:HE2	1.95	0.48
1:D:443:ILE:HD13	1:D:466:LEU:HD23	1.96	0.48
1:D:478:SER:C	1:D:480:SER:N	2.66	0.48
1:D:442:ALA:HA	1:D:445:ILE:HD12	1.96	0.48
1:D:287:VAL:N	1:D:288:PRO:HD3	2.29	0.47
2:A:339:ALA:HB2	4:A:4000:EPH:H202	1.95	0.47
1:D:435:VAL:HG21	1:D:472:TYR:CZ	2.49	0.47
1:D:497:ILE:HA	1:D:497:ILE:HD13	1.69	0.47
1:D:454:GLN:N	1:D:455:PRO:HD3	2.30	0.47
1:D:343:THR:O	1:D:347:VAL:HG23	2.15	0.47
2:A:341:VAL:HG12	2:A:434:HIS:CE1	2.50	0.47
2:A:400:ASP:O	2:A:404:ARG:HG3	2.15	0.47
2:A:415:ALA:O	2:A:419:LEU:HB2	2.14	0.47
2:A:406:SER:C	2:A:407:ARG:HG3	2.35	0.47
2:A:413:ARG:HG3	2:A:413:ARG:HH11	1.80	0.47
1:D:369:ILE:O	1:D:373:LYS:HB2	2.16	0.46
2:A:213:LEU:HD23	2:A:216:MET:SD	2.55	0.46
2:A:242:PHE:C	2:A:245:PRO:HD2	2.36	0.46
2:A:274:GLU:O	2:A:277:ASP:HB2	2.16	0.46
2:A:413:ARG:HG3	2:A:413:ARG:NH1	2.30	0.46
2:A:236:SER:HB2	2:A:238:VAL:HG23	1.97	0.46
2:A:341:VAL:HG12	2:A:434:HIS:HD1	1.80	0.46
1:D:508:CYS:O	1:D:518:LEU:HD11	2.15	0.46
2:A:227:PHE:HB3	2:A:229:PHE:CE1	2.50	0.46
2:A:301:PHE:CD2	2:A:322:LEU:HD22	2.50	0.46
1:D:384:VAL:HG23	1:D:395:VAL:HG11	1.98	0.46
1:D:432:MET:HB2	1:D:437:TYR:CE1	2.51	0.46
1:D:409:ARG:O	1:D:412:GLY:N	2.35	0.46
1:D:481:PRO:O	1:D:483:CYS:N	2.49	0.45
1:D:339:ILE:HG23	3:D:3000:P1A:C16	2.47	0.45
2:A:265:ARG:HA	2:A:270:PHE:HD2	1.80	0.45
1:D:466:LEU:HD21	1:D:491:LEU:HD21	1.99	0.45
2:A:216:MET:HA	2:A:219:LEU:HD12	1.98	0.45
2:A:439:HIS:HB3	5:A:1009:HOH:O	2.16	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:A:401:GLU:HG3	2:A:402:TYR:N	2.30	0.45
2:A:376:ASN:HA	2:A:377:PRO:HD3	1.69	0.45
1:D:379:VAL:C	1:D:381:MET:H	2.19	0.45
2:A:296:TRP:CZ3	2:A:361:VAL:HG22	2.51	0.45
2:A:348:VAL:O	2:A:352:LEU:HB2	2.16	0.45
1:D:366:SER:O	1:D:369:ILE:HG23	2.17	0.45
1:D:460:GLU:C	1:D:462:GLN:N	2.68	0.45
1:D:381:MET:SD	1:D:420:LEU:HD11	2.57	0.45
1:D:507:MET:O	1:D:511:LEU:HG	2.17	0.45
1:D:509:ILE:O	1:D:511:LEU:N	2.50	0.45
1:D:344:ILE:HD11	1:D:519:PRO:CG	2.47	0.45
1:D:336:PHE:CZ	1:D:516:ARG:HG3	2.52	0.44
2:A:393:GLU:O	2:A:397:LEU:HD22	2.18	0.44
2:A:407:ARG:HB3	2:A:410:GLU:HG2	1.99	0.44
2:A:401:GLU:O	2:A:404:ARG:HB2	2.18	0.44
2:A:216:MET:CE	2:A:266:ASP:O	2.66	0.44
1:D:339:ILE:HG23	3:D:3000:P1A:H162	2.00	0.44
2:A:356:MET:O	2:A:357:ARG:C	2.55	0.44
1:D:453:GLU:HB3	1:D:454:GLN:NE2	2.33	0.44
1:D:354:ALA:CB	1:D:372:LEU:HD21	2.47	0.44
2:A:299:MET:O	2:A:302:LEU:HB2	2.18	0.44
2:A:446:ILE:O	2:A:450:ILE:HG13	2.17	0.44
2:A:240:PRO:HA	2:A:243:ARG:HG2	1.99	0.44
2:A:455:ARG:HA	2:A:455:ARG:HD3	1.84	0.44
2:A:205:VAL:HG22	2:A:206:GLN:N	2.33	0.43
1:D:378:GLU:OE1	1:D:498:ARG:HD2	2.17	0.43
2:A:216:MET:HE1	2:A:266:ASP:O	2.18	0.43
1:D:484:ALA:O	1:D:487:PHE:N	2.50	0.43
1:D:369:ILE:O	1:D:373:LYS:HG3	2.18	0.43
1:D:354:ALA:HA	1:D:357:LEU:HD12	1.99	0.43
2:A:272:GLN:NE2	2:A:272:GLN:HA	2.33	0.43
1:D:469:LEU:O	1:D:473:ILE:HG13	2.19	0.43
2:A:208:LEU:CB	2:A:268:PRO:HB2	2.48	0.43
1:D:381:MET:CB	1:D:497:ILE:HD12	2.47	0.43
1:D:313:GLU:O	1:D:317:LYS:HB2	2.19	0.43
1:D:510:SER:O	1:D:514:LYS:HG3	2.19	0.43
1:D:344:ILE:HG23	1:D:521:PHE:CD2	2.53	0.43
2:A:257:ILE:O	2:A:261:VAL:HG23	2.19	0.42
1:D:449:ARG:HA	1:D:450:PRO:HD3	1.86	0.42
2:A:368:ALA:O	2:A:369:LEU:C	2.57	0.42
2:A:401:GLU:O	2:A:405:ARG:N	2.49	0.42
2:A:418:LEU:HA	2:A:418:LEU:HD23	1.87	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:A:251:GLN:HB3	2:A:251:GLN:HE21	1.54	0.42
1:D:409:ARG:O	1:D:410:LYS:C	2.56	0.42
2:A:212:ARG:O	2:A:215:GLU:HB2	2.19	0.42
1:D:509:ILE:C	1:D:511:LEU:N	2.73	0.42
1:D:301:LEU:HB3	1:D:438:ALA:HB1	2.02	0.42
2:A:207:GLU:OE2	2:A:207:GLU:HA	2.20	0.42
1:D:423:PHE:O	1:D:427:MET:HG2	2.19	0.42
2:A:303:THR:O	2:A:304:GLU:O	2.37	0.42
1:D:388:TYR:CG	1:D:424:CYS:HB3	2.55	0.42
2:A:208:LEU:HD12	2:A:391:LEU:HD22	2.01	0.41
1:D:295:LYS:HD2	1:D:295:LYS:HA	1.82	0.41
1:D:373:LYS:HD2	1:D:525:ILE:O	2.20	0.41
2:A:254:ASN:O	2:A:255:LYS:C	2.58	0.41
2:A:451:ARG:NH1	2:A:452:ASP:OD1	2.53	0.41
1:D:502:MET:HE1	2:A:429:LEU:CB	2.46	0.41
2:A:440:LEU:HA	2:A:440:LEU:HD12	1.92	0.41
2:A:436:PHE:HE2	2:A:441:VAL:HB	1.85	0.41
1:D:346:THR:OG1	3:D:3000:P1A:H9	2.21	0.41
2:A:407:ARG:O	2:A:409:SER:N	2.54	0.41
1:D:357:LEU:CD1	1:D:445:ILE:HD13	2.50	0.41
2:A:244:ALA:N	2:A:245:PRO:CD	2.82	0.41
1:D:354:ALA:HB1	1:D:372:LEU:HD21	2.02	0.41
1:D:357:LEU:HA	1:D:358:PRO:HD3	1.89	0.41
1:D:521:PHE:CD2	1:D:522:LEU:HD13	2.57	0.40
1:D:516:ARG:HD2	1:D:516:ARG:HA	1.72	0.40
2:A:451:ARG:HD2	2:A:452:ASP:CA	2.51	0.40
2:A:238:VAL:O	2:A:239:PRO:C	2.60	0.40
1:D:367:ASP:O	1:D:370:THR:N	2.55	0.40
2:A:322:LEU:O	2:A:349:LEU:HD11	2.22	0.40
2:A:239:PRO:HA	2:A:240:PRO:HD3	1.87	0.40
1:D:474:LEU:HD23	1:D:474:LEU:HA	1.91	0.40
1:D:301:LEU:HD21	1:D:357:LEU:HD23	2.04	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	Distance(Å)	Clash(Å)
2:A:329:MET:SD	2:A:329:MET:SD[4_557]	2.08	0.12

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	D	232/266 (87%)	193 (83%)	29 (12%)	10 (4%)	4	15
2	A	237/263 (90%)	197 (83%)	32 (14%)	8 (3%)	6	23
All	All	469/529 (89%)	390 (83%)	61 (13%)	18 (4%)	5	19

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	322	THR
1	D	479	ALA
2	A	301	PHE
1	D	482	ARG
2	A	408	SER
2	A	443	ASP
1	D	289	PRO
1	D	291	THR
1	D	510	SER
1	D	409	ARG
2	A	268	PRO
2	A	451	ARG
1	D	410	LYS
1	D	436	HIS
2	A	269	HIS
2	A	223	PRO
2	A	422	PRO
1	D	485	VAL

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	D	205/235 (87%)	180 (88%)	25 (12%)	7	20
2	A	212/232 (91%)	186 (88%)	26 (12%)	7	20
All	All	417/467 (89%)	366 (88%)	51 (12%)	7	20

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	290	LEU
1	D	306	GLU
1	D	313	GLU
1	D	316	LEU
1	D	320	THR
1	D	349	LEU
1	D	369	ILE
1	D	381	MET
1	D	405	ARG
1	D	419	ASP
1	D	421	LEU
1	D	430	MET
1	D	439	LEU
1	D	447	SER
1	D	452	LEU
1	D	463	ARG
1	D	466	LEU
1	D	476	GLN
1	D	477	ASN
1	D	483	CYS
1	D	495	THR
1	D	502	MET
1	D	505	SER
1	D	516	ARG
1	D	522	LEU
2	A	208	LEU
2	A	231	ARG
2	A	243	ARG
2	A	248	SER
2	A	251	GLN
2	A	263	TRP
2	A	266	ASP
2	A	278	GLN
2	A	288	GLU
2	A	290	LEU

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Mol	Chain	Res	Type
2	A	322	LEU
2	A	326	MET
2	A	329	MET
2	A	330	THR
2	A	337	LEU
2	A	355	LYS
2	A	359	LEU
2	A	361	VAL
2	A	384	ASN
2	A	386	GLN
2	A	397	LEU
2	A	411	GLU
2	A	421	LEU
2	A	429	LEU
2	A	430	LYS
2	A	440	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	294	GLN
1	D	305	GLN
1	D	348	GLN
1	D	368	GLN
1	D	407	ASN
1	D	454	GLN
1	D	462	GLN
1	D	476	GLN
2	A	206	GLN
2	A	228	GLN
2	A	251	GLN
2	A	256	GLN
2	A	272	GLN
2	A	287	ASN
2	A	338	GLN
2	A	343	GLN
2	A	363	GLN
2	A	376	ASN
2	A	386	GLN
2	A	439	HIS

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	$\# Z > 2$
4	EPH	A	4000	-	48,48,48	1.66	8 (16%)	53,53,53	1.93	12 (22%)
3	P1A	D	3000	-	36,36,36	3.32	17 (47%)	60,60,60	2.29	17 (28%)

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
4	EPH	A	4000	-	-	0/52/52/52	0/0/0/0
3	P1A	D	3000	-	-	0/17/87/87	0/0/4/4

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	3000	P1A	C20-C22	8.03	1.68	1.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	3000	P1A	C1-C10	6.35	1.64	1.54
3	D	3000	P1A	C9-C8	6.20	1.65	1.52
3	D	3000	P1A	O3-C3	-5.66	1.30	1.43
3	D	3000	P1A	C3-C2	5.53	1.60	1.52
3	D	3000	P1A	C21-C20	5.45	1.63	1.52
3	D	3000	P1A	C5-C6	4.96	1.59	1.51
3	D	3000	P1A	C23-C22	4.65	1.60	1.52
3	D	3000	P1A	C10-C9	4.51	1.64	1.56
4	A	4000	EPH	C25-C24	4.44	1.56	1.31
4	A	4000	EPH	C16-C15	4.35	1.55	1.31
3	D	3000	P1A	C19-C10	4.32	1.62	1.54
4	A	4000	EPH	C29-C28	4.21	1.55	1.31
3	D	3000	P1A	C12-C13	3.98	1.62	1.54
4	A	4000	EPH	C13-C12	3.72	1.52	1.31
4	A	4000	EPH	C32-C33	3.51	1.54	1.29
4	A	4000	EPH	P1-O8	-3.23	1.44	1.59
3	D	3000	P1A	O20-C20	-3.23	1.38	1.44
3	D	3000	P1A	C11-C9	3.12	1.59	1.53
3	D	3000	P1A	O6-C6	-2.84	1.18	1.22
3	D	3000	P1A	C7-C6	2.68	1.51	1.45
4	A	4000	EPH	P1-O6	-2.47	1.44	1.55
4	A	4000	EPH	O8-C38	2.19	1.54	1.44
3	D	3000	P1A	C16-C17	2.15	1.58	1.54
3	D	3000	P1A	C18-C13	2.08	1.58	1.54

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	4000	EPH	O8-P1-O5	7.71	126.67	104.68
3	D	3000	P1A	C21-C20-C22	7.71	114.95	110.28
3	D	3000	P1A	C24-C23-C22	-5.79	106.50	112.57
3	D	3000	P1A	C18-C13-C14	5.55	114.91	109.14
3	D	3000	P1A	C20-C17-C13	5.31	126.58	120.59
4	A	4000	EPH	O8-P1-O7	-4.95	89.97	109.37
4	A	4000	EPH	O5-P1-O7	-4.91	90.12	109.37
3	D	3000	P1A	O14-C14-C13	3.96	115.65	107.94
3	D	3000	P1A	O14-C14-C8	-3.89	98.67	105.75
4	A	4000	EPH	O6-P1-O7	-3.87	90.57	112.21
3	D	3000	P1A	O14-C14-C15	-3.73	102.42	110.69
3	D	3000	P1A	C19-C10-C9	3.39	115.96	109.94
4	A	4000	EPH	O6-P1-O8	3.29	125.08	108.51
3	D	3000	P1A	C10-C1-C2	-2.99	109.17	114.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	3000	P1A	O6-C6-C7	-2.97	116.44	121.55
4	A	4000	EPH	C1-O2-C4	2.92	125.67	117.13
3	D	3000	P1A	C4-C5-C10	2.89	116.93	112.95
3	D	3000	P1A	C19-C10-C5	-2.75	105.05	109.94
3	D	3000	P1A	C4-C3-C2	2.34	113.36	110.37
4	A	4000	EPH	C11-C12-C13	-2.34	112.19	125.43
3	D	3000	P1A	C12-C13-C14	-2.25	104.42	107.66
4	A	4000	EPH	C31-C30-C29	2.22	119.01	112.88
4	A	4000	EPH	P1-O8-C38	-2.19	109.32	121.66
3	D	3000	P1A	O22-C22-C23	-2.19	105.12	109.11
4	A	4000	EPH	C14-C13-C12	-2.17	111.50	124.86
4	A	4000	EPH	C17-C16-C15	-2.15	113.23	125.43
4	A	4000	EPH	C23-C24-C25	-2.10	113.52	125.43
3	D	3000	P1A	O20-C20-C17	-2.08	105.26	109.02
3	D	3000	P1A	C10-C9-C8	-2.08	109.16	112.46

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	D	236/266 (88%)	0.11	11 (4%) 30 37	40, 71, 98, 113	0
2	A	241/263 (91%)	0.26	25 (10%) 7 9	42, 70, 107, 122	0
All	All	477/529 (90%)	0.18	36 (7%) 13 17	40, 70, 104, 122	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	A	317	THR	5.5
2	A	316	THR	5.5
2	A	226	GLU	4.7
1	D	333	ASP	4.4
2	A	205	VAL	4.2
1	D	332	SER	3.8
1	D	323	TRP	3.8
2	A	225	GLU	3.7
2	A	233	GLY	3.7
2	A	315	ARG	3.6
1	D	431	MET	3.4
1	D	478	SER	3.4
2	A	219	LEU	3.4
2	A	223	PRO	3.3
2	A	405	ARG	3.2
2	A	410	GLU	3.1
1	D	482	ARG	3.0
2	A	455	ARG	2.9
2	A	411	GLU	2.9
2	A	318	SER	2.8
2	A	303	THR	2.8
2	A	224	SER	2.8
1	D	481	PRO	2.6
2	A	237	ASN	2.5

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Mol	Chain	Res	Type	RSRZ
1	D	514	LYS	2.4
1	D	513	LEU	2.4
2	A	304	GLU	2.4
2	A	218	SER	2.4
2	A	228	GLN	2.4
2	A	380	LYS	2.4
2	A	444	THR	2.3
2	A	276	GLU	2.2
2	A	409	SER	2.2
1	D	515	ASN	2.1
1	D	296	SER	2.0
2	A	381	GLY	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(Å ²)	Q<0.9
4	EPH	A	4000	49/49	0.35	3.69	86,92,111,111	0
3	P1A	D	3000	33/33	0.29	2.65	43,46,55,56	0

6.5 Other polymers ⓘ

There are no such residues in this entry.